

Design and Implementation of a Parallel Threshold Markowitz Algorithm

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Introduction

Goal: solve highly unsymmetric sparse systems.

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 - ▶ chemical engineering
 - ▶ linear programming
 - ▶ economic modelling
 - ▶ power systems

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 - ▶ linear programming
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 - ▶ power systems

Both code and matrices can be very **evil**

Highly unsymmetric systems

Highly unsymmetric systems

$$si(A) = \frac{\text{number}_{i \neq j} \{a_{ij} * a_{ji} \neq 0\}}{nz\{A\}}$$

Highly unsymmetric systems

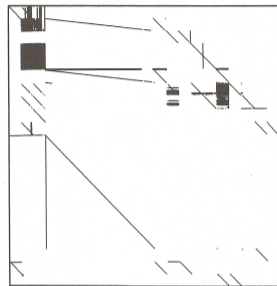
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A matrix A is highly unsymmetric if
 $si(A) < 0.9$

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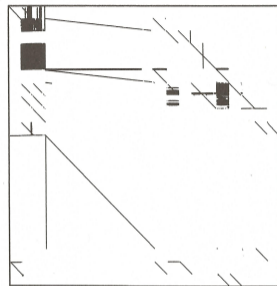


Matrix from econometric model of SE Asia

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Matrix from econometric model of SE Asia

State of the art solvers: MA48, UMFPACK, SuperLU, MUMPS.

Threshold Markowitz pivoting

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- ▶ Threshold value
Consider only entries a_{ij} that satisfy

$$|a_{ij}| \geq u * \max_k |a_{kj}|, k = 1, n$$

where u is a threshold parameter $0 < u \leq 1.0$.

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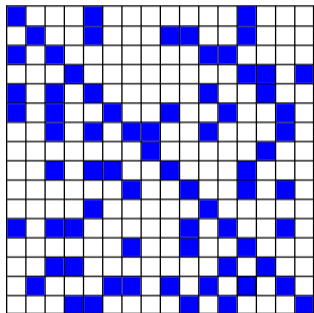
- ▶ Markowitz threshold

If there are r_i entries in row i and c_j entries in column j , the Markowitz count for the entry a_{ij} is given by $M(a_{ij}) = (r_i - 1)(c_j - 1)$. In each column, consider only the entries that satisfy the threshold test $M(a_{ij}) \leq \alpha * \min_{\text{markowitz}}$ where $\min_{\text{markowitz}}$ is the minimum Markowitz cost for an entry in the matrix.

High-level algorithm description

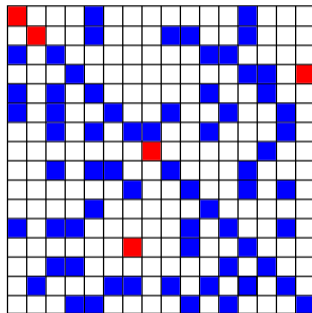
while $size(A) > 1$ **do**

end while



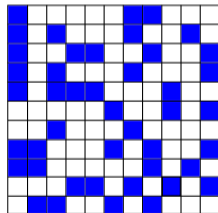
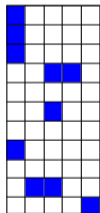
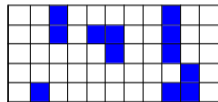
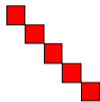
High-level algorithm description

```
while  $size(A) > 1$  do  
    Find a set of independent pivots  
end while
```

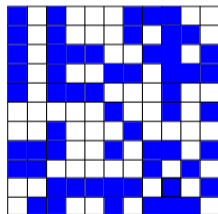
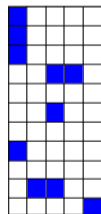
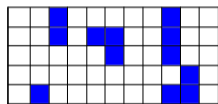


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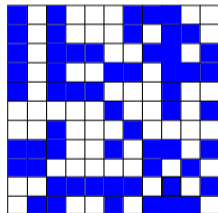
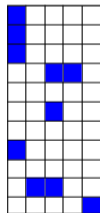
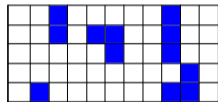
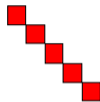
High-level algorithm description



while $size(A) > 1$ **do**
 Find a set of independent pivots
 Update the trailing matrix
end while

High-level algorithm description

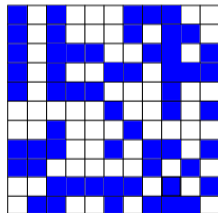
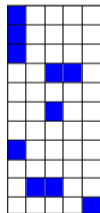
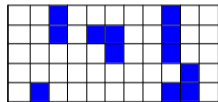
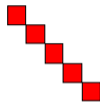
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while  $\text{density}(A) < \text{eps}$  do  
  Find a set of independent pivots  
  Update the trailing matrix  
end while
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High-level algorithm description

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while  $\text{density}(A) < \text{eps}$  do  
  Find a set of independent pivots  
  Update the trailing matrix  
end while
```

Switch to dense factorization.



Luby's Algorithm for a **Maximal Independent Set (MIS)**

Input $G = (V, E)$ an undirected graph

Output $I \subseteq V$, an MIS

$I \leftarrow \emptyset$

$G' = (V', E') \leftarrow G = (V, E)$

while $V' \neq \emptyset$ **do**

assign random score to each node in V'

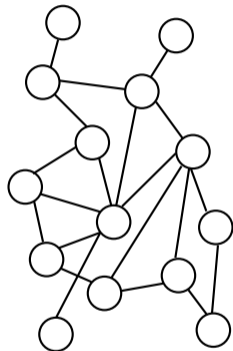
$I' \leftarrow$ nodes having highest score among their neighbours

$I \leftarrow I \cup I'$

$Y \leftarrow I' \cup N(I')$

Set $G' = (V', E')$ to the induced subgraph on $V' - Y$

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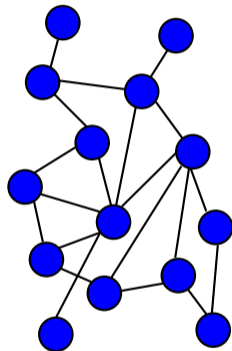
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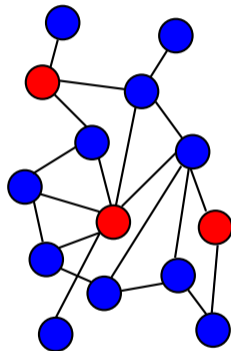
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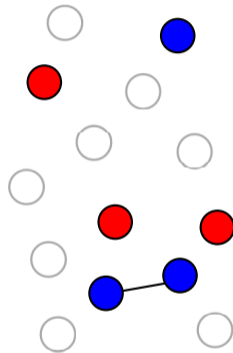
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Output $I \subseteq V$, an MIS

$I \leftarrow \emptyset$

$G' = (V', E') \leftarrow G = (V, E)$

while $V' \neq \emptyset$ **do**

assign random score to each node in V'

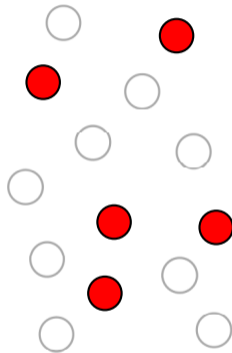
$I' \leftarrow$ *nodes having highest score among their neighbours*

$I \leftarrow I \cup I'$

$Y \leftarrow I' \cup N(I')$

 Set $G' = (V', E')$ to the induced subgraph on $V' - Y$

end while



We then continue the process to obtain an MIS with 5 nodes.

Luby's Algorithm for a Maximal Independent Set (MIS)

Input $G = (V, E)$ an undirected graph

Output $I \subseteq V$, an MIS

$I \leftarrow \emptyset$

$G' = (V', E') \leftarrow G = (V, E)$

while $V' \neq \emptyset$ **do**

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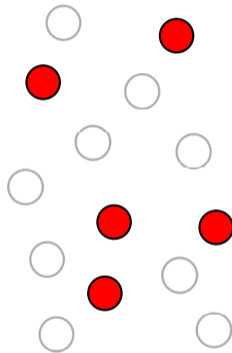
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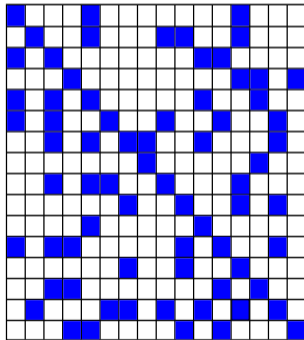
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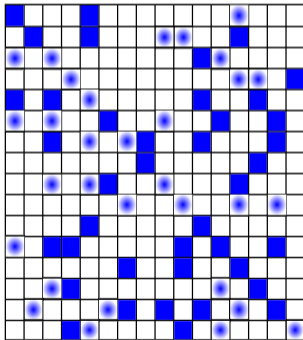
We have adopted the Luby's algorithm for directed graphs.

Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library



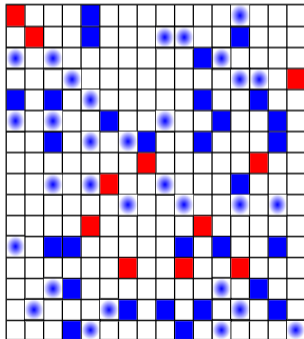
Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library

Discard all numerically ineligible entries and calculate the minimum Markowitz cost.



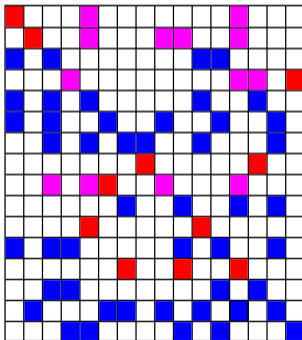
Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library

Choose potential pivots that satisfy the Markowitz threshold test, one for each column and assign a score for each one. The pivot score is associated with the column.



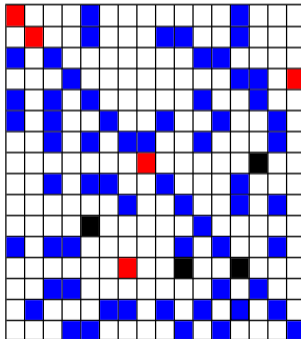
Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library

For each potential pivot scan its row, comparing its score with the score of the columns with entries in the row.



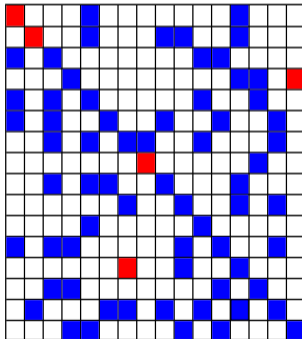
Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library

Discard the column with a potential pivot with lower score.



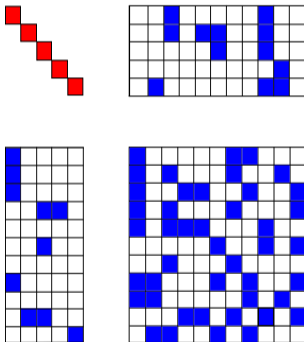
Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library

The set of independent pivots.



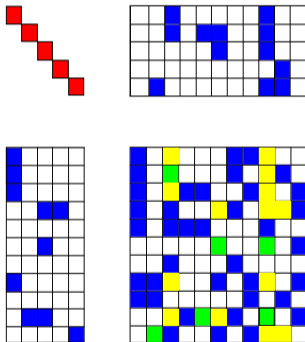
Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library

This results in a reordered matrix of the form:



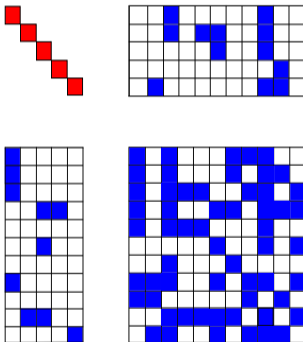
Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library

The trailing matrix is then updated by a sparse matrix-matrix multiply.



Parallel Solver for Highly Unsymmetric Matrices (ParSHUM) library

The pivot search is then repeated on the reduced matrix.



Experimental set-up

HPC2N platform at
Umeå University

Each node:

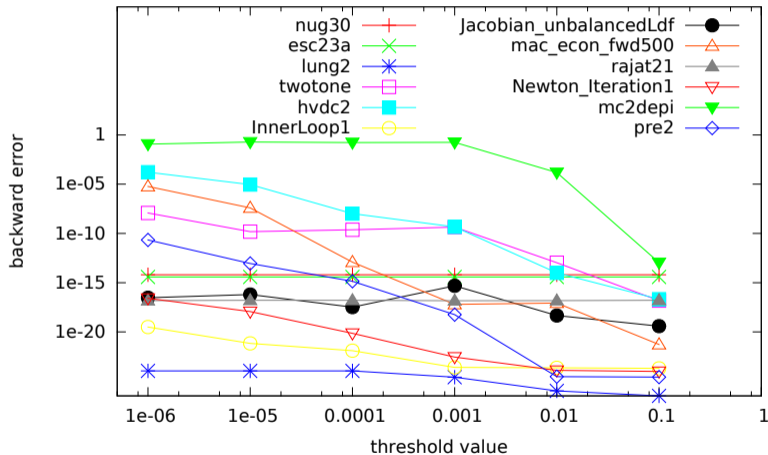
28 Intel Xeon E5-2690v4

35 MB of shared L3 cache

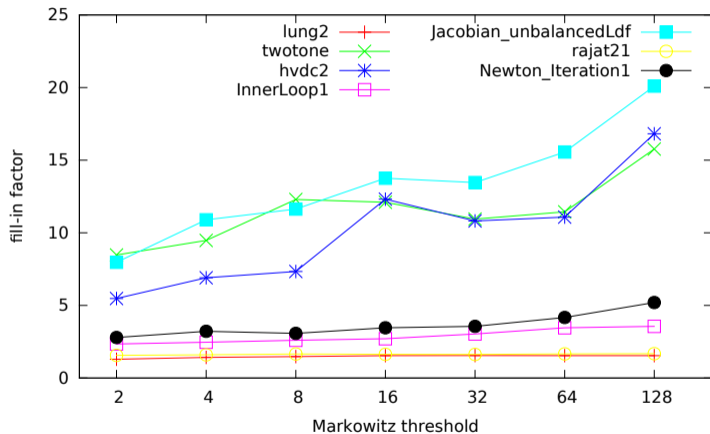
128 GB RAM memory

Matrix	Order $\times 10^3$	Entries $\times 10^6$	<i>si</i>
nug30	52.4	0.24	0.00
esc32a	63.6	0.31	0.00
lung2	109	0.49	0.57
twotone	120	1.22	0.26
hvdc2	190	1.35	0.99
InnerLoop1	197	0.75	0.44
Jacobian_unbalancedLdf	203	2.41	0.80
mac_econ_fwd500	206	1.27	0.07
rajat21	411	1.89	0.76
Newton_iteration1	427	2.38	0.14
esc64a	504	2.40	0.00
mc2depi	525	2.10	0.00
pre2	659	5.96	0.36

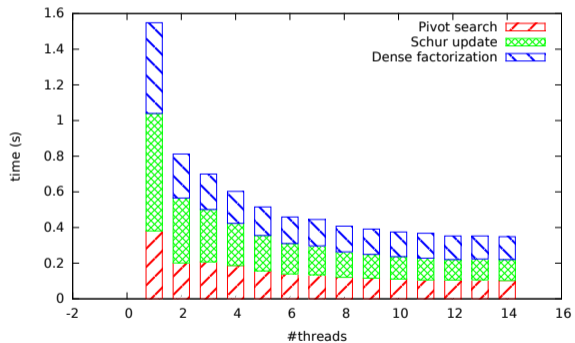
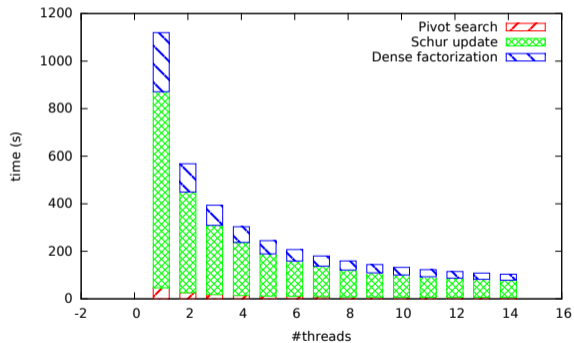
Effect of the threshold value on the backward error



Effect of the Markowitz threshold on the fill-in factor ($nz(LU)/nz(A)$)



Scalability on the mc2depi and twotone matrices



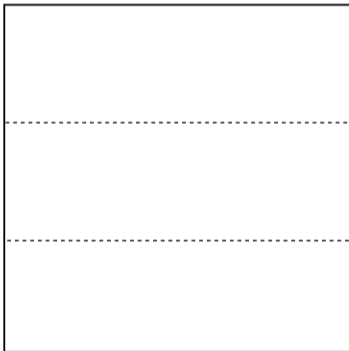
Comparison of ParSHUM, MUMPS and UMFPACK on single NUMA node (14 cores)

Matrix	ParSHUM		MUMPS		UMFPACK	
	time	fill-in	time	fill-in	time	fill-in
nug30	0.71	142.	6.19	730	48.15	463.
esc32a	0.46	70.9	10.59	758	71.0	341.
lung2	0.04	1.48	—	—	0.10	1.44
twotone	0.35	7.37	2.22	25.4	0.49	5.45
hvdc2	0.38	6.91	1.13	2.14	0.38	2.06
InnerLoop1	0.13	2.71	1.64	3.90	0.30	2.48
Jacobian_unbalancedLdf	0.99	10.9	1.52	3.45	0.74	3.88
mac_econ_fwd500	33.5	450.	10.5	56.4	4.62	57.5
rajat21	0.11	1.64	—	—	44.5	1.89
Newton_Iteration1	0.46	3.46	4.26	6.02	1.00	2.55
esc64a	12.8	155	—	2151	—	—
mc2depi	104.	302.	4.60	25	4.36	38.6
pre2	13.0	57.0	9.73	18.4	25.8	32.3

Table: The execution time and the fill-in factor for ParSHUM, MUMPS and UMFPACK.

Towards distributed memory: Singly Bordered Block decomposition using Zoltan

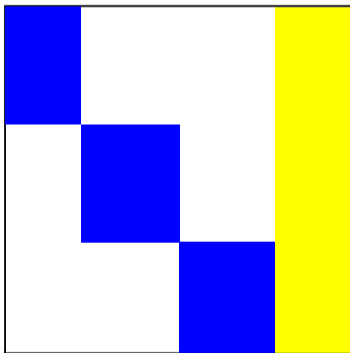
Global view



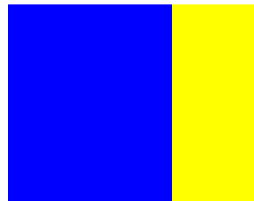
Local view

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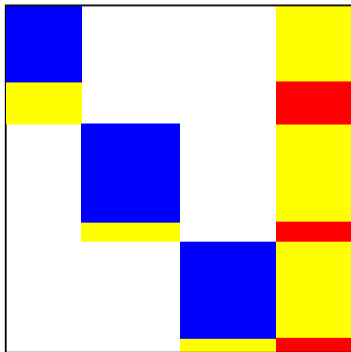


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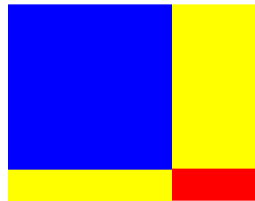


Towards distributed memory: Singly Bordered Block decomposition using Zoltan

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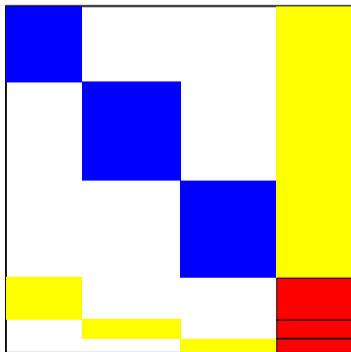


Local view



Towards distributed memory: Singly Bordered Block decomposition using Zoltan

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Each node:

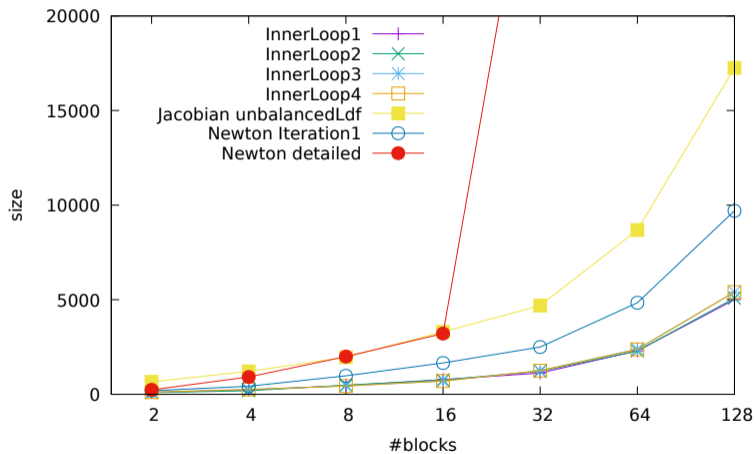
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Matrix	Order $\times 10^3$	Entries $\times 10^6$	<i>si</i>
InnerLoop1	197	0.75	0.44
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Jacobian_unbalancedLdf	203	2.41	0.80
Newton_iteration1	427	2.38	0.14
Newton_detailed	7355	24	0.29

Bordered Block size



Scalability

Matrix	#MPI processes			
	1	2	4	8
InnerLoop1	0.15	0.08	0.06	0.05
InnerLoop2	0.15	0.08	0.06	0.05
InnerLoop3	0.15	0.09	0.06	0.05
InnerLoop4	0.15	0.09	0.06	0.05
Jacobian_unbalancedLdf	1.11	0.52	0.35	0.18
Newton_iteration1	0.49	0.30	0.14	0.09
Newton_detailed	7.39	3.01	1.67	1.63

The execution time in seconds for ParSHUM on the test matrices partitioned in SBBD form. One numa node (14 cores) is used per process.

Comparison with MUMPS and SuperLU

Matrix	ParSHUM		MUMPS		SuperLU	
	time	fill-in	time	fill-in	time	fill-in
InnerLoop1	0.05	3.03	0.23	3.90	1.11	6.02
InnerLoop2	0.05	2.82	0.25	3.72	1.11	5.57
InnerLoop3	0.05	2.78	0.17	3.72	1.10	5.60
InnerLoop4	0.05	2.41	0.15	3.73	1.15	5.56
Jacobian_unbalancedLdf	0.18	6.75	0.28	3.45	1.05	3.59
Newton_iteration1	0.09	3.57	0.42	6.02	2.59	5.60
Newton_detailed	1.63	5.32	7.88	6.09	52.4	5.39

Conclusions and future work

Conclusions:

- ▶ We have developed a generalization of Luby's algorithm for directed graphs.
- ▶ We have used this to develop a multi-threaded threshold Markowitz code (ParSHUM library).
- ▶ In general it outperforms established codes.

Future work:

- ▶ Ongoing work on SBBD to exploit distributed systems.
- ▶ Develop a GPU version of the solver.

THANK YOU FOR YOUR ATTENTION