

# Introduction to supercomputing: numerical experiments on shared and distributed memory cluster

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# Plan

- Login to cluster
- User setup
- Run // tests
- Plots with timing and speedup

## Instructions:

[http://cluster2.inm.ras.ru/~hpc19/0\\_INDEX\\_0.html](http://cluster2.inm.ras.ru/~hpc19/0_INDEX_0.html)

<http://dodo.inm.ras.ru/~konshin/tr/2020-Sechenov-China-exp.pdf>

## Login to cluster

```
ssh hpc19@dodo.inm.ras.ru
```

```
Password: *****
```

```
To user setup:                xsetup nickname
To request the computing resources: xqs
To list the parallel tests available: xlist
To run parallel test (on nodes):   xrun test
To construct the plots (on h2):    xplot
To view plots:      http://cluster2.inm.ras.ru/~hpc19/nickname/*.png
To view sources:   http://cluster2.inm.ras.ru/~hpc19/sources/*.c
To see this help info:                xinfo
hpc19@h2:~/public_html>
```

*j Now you are at the host computer h2 in account hpc19 !*

## User setup: **xsetup**

```
hpc19@h2:~/public_html> xsetup nickname  
Setup for user: nickname  
hpc19@h2:~/public_html/nickname>
```

Here *nickname* is your personal identification name or surname

*j* Now you are at **your own personal local directory** !

## Computing resources request: **xqs**

```
hpc19@h2:~/public_html/nickname> xqs
```

```
qsub: waiting for job 52193.h2.hpc to start
```

```
qsub: job 52193.h2.hpc ready
```

```
hpc19@cl1n036:~/public_html/nickname>
```

*j Now you are at your local directory at your personal cluster node !*

Linear algebra: daxpy, norm, mvm

## List of parallel tests: **xlist**

```
hpc19@cl1n036:~/public_html/nickname> xlist
```

```
omp_daxpy      omp_norm      omp_mvm
mpi_daxpy      mpi_norm      mpi_mvm-bcast  mpi_mvm-isend
t1a t2a t2b t2c t2d t3b t3c t3d t4b t4c t4d t5h t5i t5j t6k t6l
```

Types of tests:

- `omp_*` — tests on OpenMP (shared memory)
- `mpi_*` — tests on MPI (distributed memory)
- `t1a ... t6l` — INMOST tests (distributed memory)
- `t3d t4d t5h t5i t5j t6k t6l` — INMOST tests with `*.jpg` output

List of examples:

- `*_daxpy`:  $\alpha \cdot \vec{x} + \vec{y}$
- `*_norm`:  $\|\vec{x}\|$
- `*_mvm*`:  $y = A \cdot x$

*`j *_bcast` and `*_isend` are different implementations of the same example !*



## Parallel run: `xrun`

```
hpc19@cl1n036:~/public_html/nickname> xrun omp_daxpy
```

```
#OMP daxpy: N=1048576 M=3 np=12 ::: p time(sec) perf(MFLOPS) S=t(1)/t(p)
```

1	0.021050	149.437	1.000
2	0.014708	213.873	1.431
3	0.007322	429.634	2.875
4	0.005641	557.627	3.732
5	0.008197	383.745	2.568
6	0.006517	482.684	3.230
8	0.010810	291.000	1.947
10	0.003337	942.588	6.308
12	0.004306	730.547	4.889

*i // run of daxpy on shared memory with OpenMP !*

## Construct the plots: **xplot**

Then, go back to h2 (exit) and run xplot to construct the plots for timing and speedup

```
hpc19@cl1n036:~/public_html/nickname> exit
qsub: job 52193.h2.hpc completed
```

```
hpc19@h2:~/public_html/nickname> xplot
```

```
-rw-r-r- 1 hpc19 users 4796 Jan 14 01:12 s.png
-rw-r-r- 1 hpc19 users 4695 Jan 14 01:12 t.png
```

See resulting "\*.png" plots in:

[http://cluster2.inm.ras.ru/~hpc19/\*nickname\*](http://cluster2.inm.ras.ru/~hpc19/<i>nickname</i>)

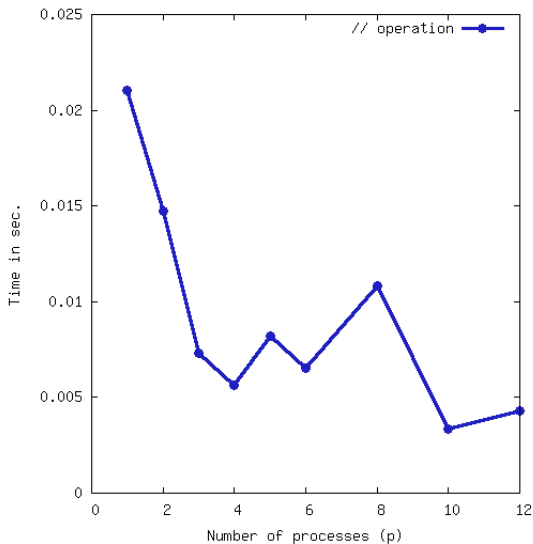
where "s.png" is for Speedup and "t.png" is for Time plots

*; Two \*.png files with plots were created on h2 host !*

*; These files can be seen in regular browser via internet !*

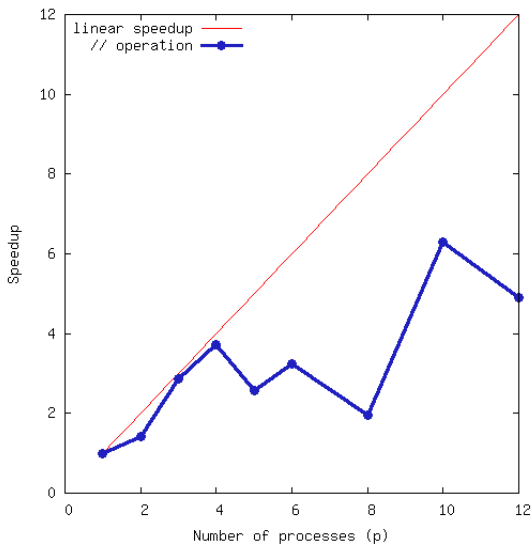
*; Do not forget to press Ctrl+<R> (Reload) in time to update !*

OMP daxpy:  $\alpha \cdot \vec{x} + \vec{y}$



*j* Time in sec.: <http://cluster2.inm.ras.ru/~hpc19/nickname/t.png> !

OMP daxpy:  $\alpha \cdot \vec{x} + \vec{y}$



*j Speedup: <http://cluster2.inm.ras.ru/~hpc19/nickname/s.png> !*

## The same test daxpy on MPI

```
hpc19@h2:~/public_html/nickname> xqs
```

```
hpc19@cl1n036:~/public_html/nickname> xrun mpi_daxpy
```

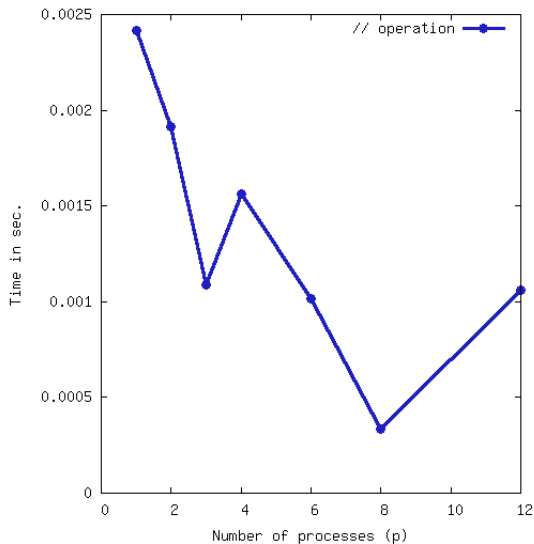
```
hpc19@cl1n036:~/public_html/nickname> exit
```

```
hpc19@h2:~/public_html/nickname> xplot
```

See results in:

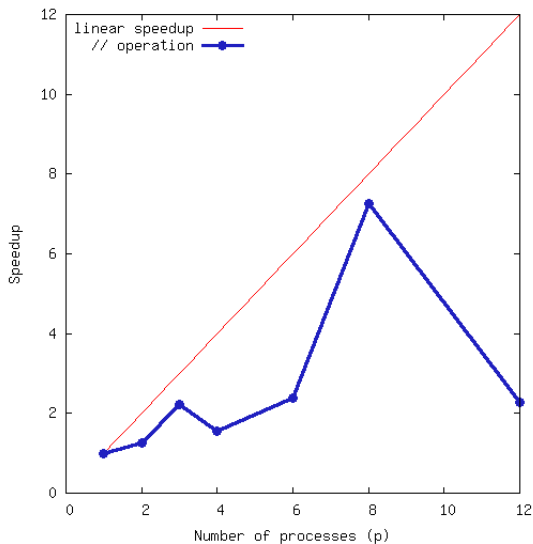
```
http://cluster2.inm.ras.ru/~hpc19/nickname/\*.png
```

MPI daxpy:  $\alpha \cdot \vec{x} + \vec{y}$



*j* Time in sec.: <http://cluster2.inm.ras.ru/~hpc19/nickname/t.png> !

MPI daxpy:  $\alpha \cdot \vec{x} + \vec{y}$



*j Speedup: <http://cluster2.inm.ras.ru/~hpc19/nickname/s.png> !*

## The other tests...

```
hpc19@cl1n036:~/public_html/nickname> xlist
```

```
omp_daxpy      omp_norm      omp_mvm
```

```
mpi_daxpy      mpi_norm      mpi_mvm-bcast  mpi_mvm-isend
```

*j* To save the plot rename \*.png file by command: `mv s.png s-new.png !`



## There are some advanced commans...

Some additional commands:

`xget test` - get test source (on h2 or nodes)

`xcc test` - compile the local test source (on h2)

`xrun` - run the local parallel test (on nodes)

*j Add “-v1” to the name of the test — it will be much more interesting :) !*

## Mathematical physics: INMOST

## INMOST tests

```
hpc19@cl1n036:~/public_html/nickname> xlist
```

```
.....
```

```
t3d t4d t5h t5i t5j t6k t6l
```

*j Pictures: [http://cluster2.inm.ras.ru/~hpc19/nickname/JPG/\\*.jpg](http://cluster2.inm.ras.ru/~hpc19/nickname/JPG/*.jpg) !*

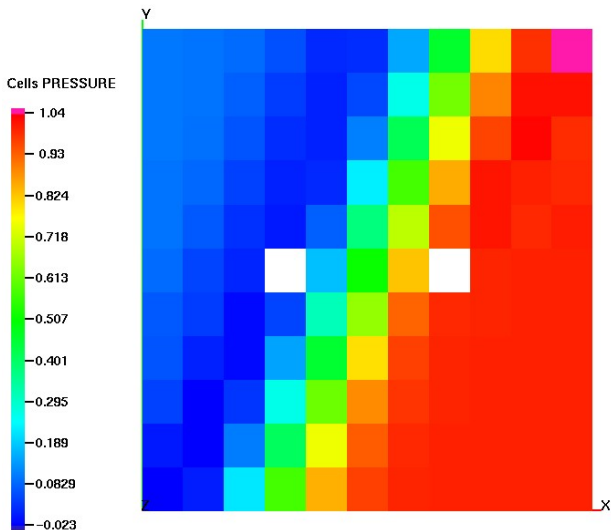
*j **exit** and **xplot** are NOT required,  
while **Ctrl+<R>** is required !*

## INMOST tests: description

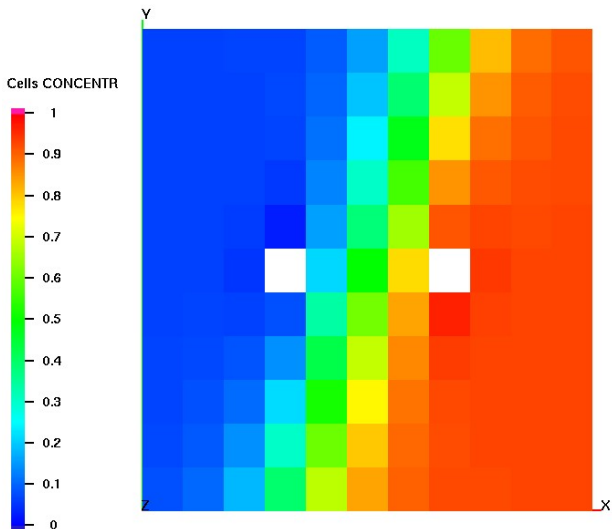
- **t3d** — anisotropic diffusion with two wells with a mimetic scheme (reference operator method)
- **t4d** — two-well anisotropic diffusion with a non-linear finite-volume scheme
- **t5h** — oil in the reservoir with anisotropic permeability that starts the flow in a zigzag
- **t5i** — oil on several layers of SPE10 test
- **t5j** — oil in the real Norn field
- **t6k** — incompressible fluid in the cavity
- **t6l** — incompressible fluid in a stepped region (L-shaped domain)

*j INMOST tests were prepared by K.M. Terekhov !*

### t3d: two-well anisotropic diffusion with a mimetic scheme (reference operator method)

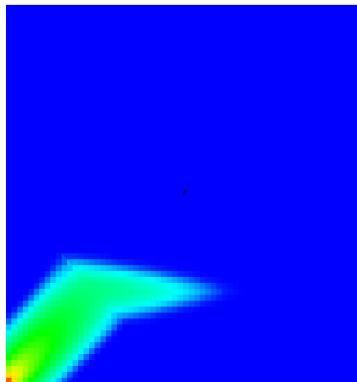
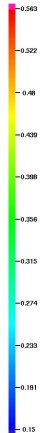


# t4d: two-well anisotropic diffusion with a non-linear finite-volume scheme

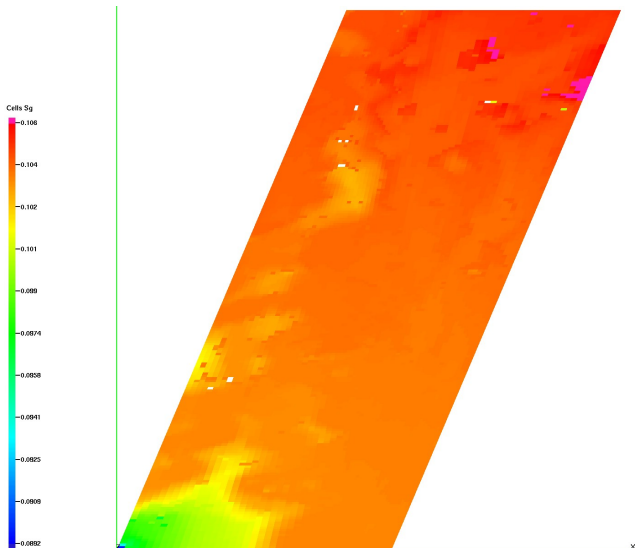


**t5h:** oil in the reservoir with anisotropic permeability  
that starts the zigzag flow

Cells Sw

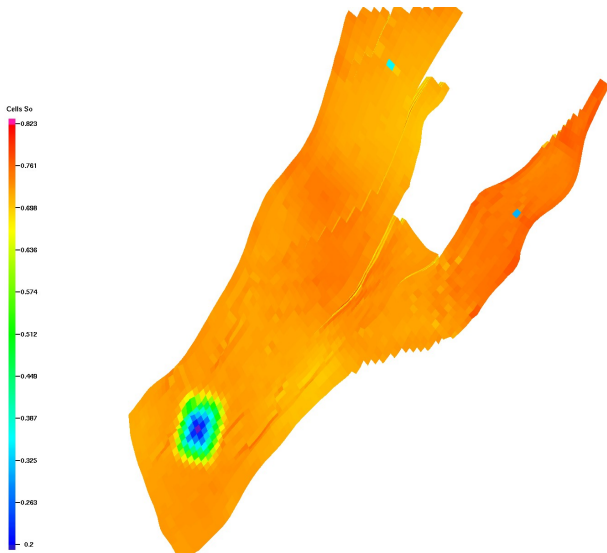


## t5i: oil in several layers of SPE10 test

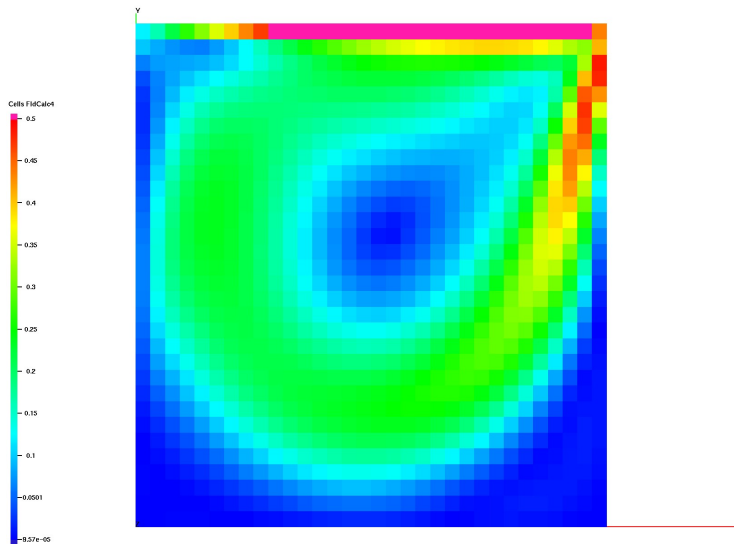




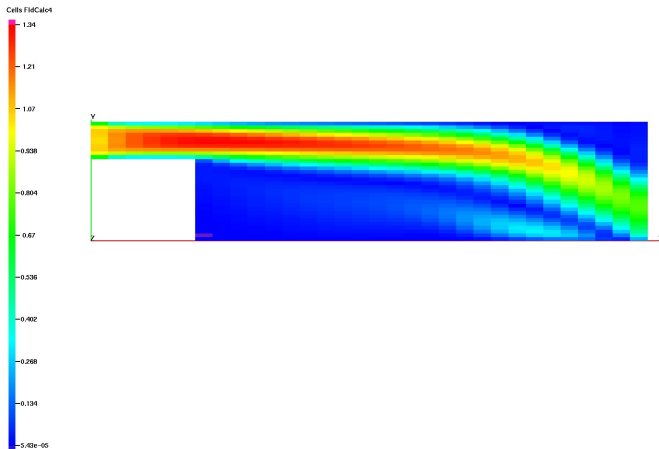
## t5j: oil in the real Norn field



# t6k: incompressible fluid in a cavity



# t6l: incompressible fluid in a stepped region (L-shaped domain)



## INMOST

<http://inmost.org> — homepage

Documentation:

<https://github.com/INMOST-DEV/INMOST/wiki> — wiki

<http://www.inmost.org/Doxygen/html/annotated.html> — doxygen

# INMOST: download and compile

Wiki instructions:

<https://github.com/INMOST-DEV/INMOST/wiki/0400-Compilation-Linux>

```
$ git clone https://github.com/INMOST-DEV/INMOST
$ mkdir INMOST/build
$ cd INMOST/build
$ cmake -DUSE_MPI=ON -DCOMPILER_EXAMPLES=ON ..
$ make
```

## INMOST: examples

Wiki instructions:

<https://github.com/INMOST-DEV/INMOST/wiki/1800-Examples>

```
$ cd Examples
```

```
### Parallel Grid Generation
```

```
$ mpiexec -np 4 GridGen/GridGen 4 32 32 32 grid.pvtk
```

```
### Parallel Finite Volume Discretization
```

```
$ mpiexec -np 4 FVDiscr/FVDiscr grid.pvtk A.mtx b.rhs
```

```
### Solve the Matrix Stored in MTX Format
```

```
$ mpiexec -np 4 MatSolve/MatSolve -m A.mtx -b b.rhs
```

```
### If you have ParaView installed:
```

```
$ paraview --data=grid.pvtk
```

```
  P_OWNER_PROCESSOR - partitioning to processors
```

```
$ paraview --data=result.pvtk
```

```
  Solution - the solution to the problem
```