

Parallel processing and matrix solver interface: mt xp

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1 Parallel processing library: MPICH

1.1 Download

1. Download a source file from MPICH web:

```
https://www.mpich.org/downloads/
mpich-3.3.1.tar.gz (26MB)
```

2. Make a directory and expand the source file

```
mkdir mpich
cd mpich
mv ~/Download./mpich-3.3.1.tar.gz .
tar xzf mpich-3.3.1.tar.gz
cd mpich-3.3.1
```

1.2 Configure

1. Create a configure script “run” by your favorite editor

```
CC=gcc CFLAGS="-m64" CXX=c++ CXXFLAGS="-m64" FC=gfortran
FFLAGS="-m64" ./configure --prefix=/usr/local/mpich331-gfortran-gcc
--enable-cxx --enable-fast --enable-romio --disable-shared
Install directory: /usr/local/mpich331-gfortran-gcc
```

2. Make “run” executable

```
chmod 755 run
```

3. Configure

```
./run
```

1.3 Compile and install

1. Compile

```
make
```

2. Install

```
sudo make install
```

2 Parallel matrix solver library: PETSc

2.1 Download

1. Download a source file from PETSc web site

```
https://www.mcs.anl.gov/petsc/download/  
petsc-3.11.3.tar.gz (33.2MB)
```

2. Make a directory and expand the source file

```
sudo mkdir /opt  
cd /opt  
sudo mkdir petsc  
sudo chown admin:admin petsc \qquad ("admin" should be replaced by you  
user name)  
cd petsc  
mv ~/Download/petsc-3.11.3.tar.gz .  
tar xzf petsc-3.11.3.tar.gz
```

2.2 Configure

1. Create a configure script “gfortran.py” by your favorite editor

```
#!/usr/bin/env python  
  
# Build PETSc, with gfortran  
  
configure_options = [  
    '--with-mpi=1',  
    '--with-mpi-dir=/usr/local/mpich331-gfortran-gcc',  
    '--with-shared-libraries=0',  
    '--with-cxx-dialect=C++11',  
    '--download-mpich=0',  
    '--download-hypre=0',  
    '--download-fblaslapack=1',  
    '--download-spooles=1',  
    '--download-superlu=1',  
    '--download-metis=1',  
    '--download-parmetis=1',
```

```

    '--download-superlu_dist=1',
    '--download-blacs=1',
    '--download-scalapack=1',
    '--download-mumps=1'
]

if __name__ == '__main__':
    import sys,os
    sys.path.insert(0,os.path.abspath('config'))
    import configure
    configure.petsc_configure(configure_options)

```

2. Make “gfortran.py” executable

```
chmod 755 gfortran.py
```

3. Setup environment variables for PETSc

```
export PETSC_DIR=/opt/petsc/petsc-3.11.3
export PETSC_ARCH=gfortran
```

4. Configure

```
./gfortran.py
```

2.3 Compile

1. Compile

```
make all
```

3 Compile of task/mtxp module

3.1 Update setup file

1. Goto mt xp directory

```
cd task/mtxp
```

2. Create a setup file

```
cp make.mtxp.org make.mtxp
```

3. Edit the setup file “make.mtxp”

- If MPI is not available, remove comment mark “#” on lines 4–9
- If MPI is available but PETs not, remove comment mark “#” on lines 12–17
- If MPI and PETSc are available, remove comment mark “#” on lines 21, 38,39, 41–45

3.2 Compile

1. Compile

```
make
```

4 Compile of task/fp and related modules

4.1 Update Makefile

1. Change directory

```
cd ../fp
```

2. Edit Makefile

- To use serial band matrix solver, remove comment mark “#” on lines 4-5.
- To use serial iterative solver, remove comment mark “#” on lines 6-7.
- To use parallel direct solver MUMPS, remove comment mark “#” on lines 8-9.
- To use parallel direct solver library PETSc, remove comment mark “#” on lines 10-11. .

4.2 Compile

1. Compile related modules and task/fp files

```
make
```