The Portable Extensible Toolkit for Scientific Computing

Matthew Knepley

Mathematics and Computer Science Division
Argonne National Laboratory

PETSc Tutorial
Exascale Computing Project Annual Meeting
Houston, TX January 14, 2019
Never believe *anything*,

unless you can run it.
Never believe *anything*,

unless you can run it.
The PETSc Team

Matt Knepley  Barry Smith  Satish Balay

Hong Zhang  Jed Brown  Lisandro Dalcin

Stefano Zampini  Mark Adams  Toby Issac
Timeline (Young People)

Lisandro
Jed
Shri
Peter
Jason
Mark
Patrick
Michael
Toby
Karl
Stefano
Dave

PETSc-3

PETSc
What I Need From You

- Tell me if you do not understand
- Tell me if an example does not work
- Suggest better wording or figures
- Followup problems at petsc-maint@mcs.anl.gov
Ask Questions!!!

- Helps me understand what you are missing
- Helps you clarify misunderstandings
- Helps others with the same question
How We Can Help at the Tutorial

- Point out relevant documentation
- Quickly answer questions
- Help install
- Guide design of large scale codes
- Answer email at petsc-maint@mcs.anl.gov
How We Can Help at the Tutorial

- Point out relevant documentation
- Quickly answer questions
- Help install
- Guide design of large scale codes
- Answer email at petsc-maint@mcs.anl.gov
How We Can Help at the Tutorial

- Point out relevant documentation
- Quickly answer questions
- Help install
- Guide design of large scale codes
- Answer email at petsc-maint@mcs.anl.gov
How We Can Help at the Tutorial

- Point out relevant documentation
- Quickly answer questions
- Help install
- Guide design of large scale codes
- Answer email at petsc-maint@mcs.anl.gov
Getting Started with PETSc

- Who uses PETSc?
- Stuff for Windows
- How can I get PETSc?
- How do I Configure PETSc?
- How do I Build PETSc?
- How do I run an example?
- How do I get more help?

PETSc Integration

Advanced Solvers

More Stuff
How did PETSc Originate?

PETSc was developed as a Platform for Experimentation

We want to experiment with different

- Models
- Discretizations
- Solvers
- Algorithms
  - which blur these boundaries
Developing parallel, nontrivial PDE solvers that deliver high performance is still difficult and requires months (or even years) of concentrated effort.

PETSc is a toolkit that can ease these difficulties and reduce the development time, but it is not a black-box PDE solver, nor a silver bullet.

— Barry Smith
You want to think about how you decompose your data structures, how you think about them globally. [...] If you were building a house, you’d start with a set of blueprints that give you a picture of what the whole house looks like. You wouldn’t start with a bunch of tiles and say, “Well I’ll put this tile down on the ground, and then I’ll find a tile to go next to it.” But all too many people try to build their parallel programs by creating the smallest possible tiles and then trying to have the structure of their code emerge from the chaos of all these little pieces. You have to have an organizing principle if you’re going to survive making your code parallel.

(http://www.rce-cast.com/Podcast/rce-28-mpich2.html)
What is PETSc?

A freely available and supported research code for the parallel solution of nonlinear algebraic equations

Free
- Download from http://www.mcs.anl.gov/petsc
- Free for everyone, including industrial users

Supported
- Hyperlinked manual, examples, and manual pages for all routines
- Hundreds of tutorial-style examples
- Support via email: petsc-maint@mcs.anl.gov

Usable from C, C++, Fortran 77/90, Matlab, Julia, and Python
What is PETSc?

- Portable to any parallel system supporting MPI, including:
  - Tightly coupled systems
    - Cray XT6, BG/Q, NVIDIA Fermi, K Computer
  - Loosely coupled systems, such as networks of workstations
    - IBM, Mac, iPad/iPhone, PCs running Linux or Windows

PETSc History

- Begun September 1991
- Over 60,000 downloads since 1995 (version 2)
- Currently 400 per month

PETSc Funding and Support

- Department of Energy
  - ECP, AMR Program, SciDAC, MICS Program, INL Reactor Program
- National Science Foundation
  - SI2, CIG, CISE, Multidisciplinary Challenge Program
- Intel Parallel Computing Center
1 Getting Started with PETSc
   • Who uses PETSc?
     • Stuff for Windows
     • How can I get PETSc?
     • How do I Configure PETSc?
     • How do I Build PETSc?
     • How do I run an example?
     • How do I get more help?
Who Uses PETSc?

Computational Scientists

- Earth Science
  - PyLith (CIG)
  - Underworld (Monash)
  - Salvus (ETHZ)
  - TerraFERMA (LDEO, Columbia, Oxford)

- Multiphysics
  - MOOSE
  - GRINS

- Subsurface Flow and Porous Media
  - PFLOTRAN (DOE)
  - STOMP (DOE)
Who Uses PETSc?

Computational Scientists

- CFD
  - IBAMR
  - Fluidity
  - OpenFVM

- Fusion
  - XGC
  - BOUT++
  - NIMROD
  - $M3D - C^1$
Who Uses PETSc?

Algorithm Developers

- **Iterative methods**
  - Deflated GMRES
  - LGMRES
  - QCG
  - SpecEst

- **Preconditioning researchers**
  - FETI-DP *(Klawonn and Rheinbach)*
  - STRUMPACK *(Ghysels and Li)*
  - HPDDM *(Jolivet and Nataf)*
  - ParPre *(Eijkhout)*
Algorithm Developers

- Discretization
  - Firedrake
  - FEniCS
  - libMesh
  - Deal II
  - PETSc-FEM
  - OOFEM
  - PetRBF

- Outer Loop Solvers
  - Eigensolvers (SLEPc)
  - Optimization (PERMON)
What Can We Handle?

- PETSc has run implicit problems with over 500 billion unknowns
  - UNIC on BG/P and XT5
  - PFLOTRAN for flow in porous media

- PETSc has run on over 1,500,000 cores efficiently
  - Gordon Bell Prize Mantle Convection on IBM BG/Q Sequoia

- PETSc applications have run at 23% of peak (600 Teraflops)
  - Jed Brown on NERSC Edison
  - HPGMG code
PETSc has run implicit problems with over 500 billion unknowns
  - UNIC on BG/P and XT5
  - PFLOTRAN for flow in porous media

PETSc has run on over 1,500,000 cores efficiently
  - Gordon Bell Prize Mantle Convection on IBM BG/Q Sequoia

PETSc applications have run at 23% of peak (600 Teraflops)
  - Jed Brown on NERSC Edison
  - HPGMG code
What Can We Handle?

- PETSc has run implicit problems with over 500 billion unknowns
  - UNIC on BG/P and XT5
  - PFLOTRTRAN for flow in porous media

- PETSc has run on over 1,500,000 cores efficiently
  - Gordon Bell Prize Mantle Convection on IBM BG/Q Sequoia

- PETSc applications have run at 23% of peak (600 Teraflops)
  - Jed Brown on NERSC Edison
  - HPGMG code
PyLith

- Multiple problems
  - Dynamic rupture
  - Quasi-static relaxation
- Multiple models
  - Nonlinear visco-plastic
  - Finite deformation
  - Fault constitutive models
- Multiple meshes
  - 1D, 2D, 3D
  - Hex and tet meshes
- Parallel
  - PETSc solvers
  - DMPLex mesh management

\(^a\) Aagaard, Knepley, Williams
Multiple Mesh Types

- Triangular
- Tetrahedral
- Rectangular
- Hexahedral
Magma Dynamics

- Couples scales
  - Subduction
  - Magma Migration

- Physics
  - Incompressible fluid
  - Porous solid
  - Variable porosity

- Deforming matrix
  - Compaction pressure

- Code generation
  - FEniCS

- Multiphysics Preconditioning
  - PETSc FieldSplit

\[ ^{a} \text{Katz} \]
Magma Dynamics

- Couples scales
  - Subduction
  - Magma Migration
- Physics
  - Incompressible fluid
  - Porous solid
  - Variable porosity
- Deforming matrix
  - Compaction pressure
- Code generation
  - FEniCS
- Multiphysics Preconditioning
  - PETSc FieldSplit

\(^a\)Katz, Speigelman
Fracture Mechanics

- Full variational formulation
  - Phase field
  - Linear or Quadratic penalty
- Uses TAO optimization
  - Necessary for linear penalty
  - Backtacking
- No prescribed cracks (movie)
  - Arbitrary crack geometry
  - Arbitrary intersections
- Multiple materials
  - Composite toughness

\[^{\text{aBourdin}}\]
Fracture Mechanics

1 Bourdin
Vortex Method

$t = 000$

- Incompressible Flow
  - Gaussian vortex blobs
  - High Re

- PetFMM
  - 2D/3D domains
  - Automatic load balancing
  - Variety of kernels
  - Optimized with templates

- PetRBF
  - Variety of RBFs
  - Uses PETSc solvers
  - Scalable preconditioner

- Parallelism
  - MPI
  - GPU

---

*Cruz, Yokota, Barba, Knepley*
Vortex Method

$t = 100$

- Incompressible Flow
  - Gaussian vortex blobs
  - High Re

- PetFMM
  - 2D/3D domains
  - Automatic load balancing
  - Variety of kernels
  - Optimized with templates

- PetRBF
  - Variety of RBFs
  - Uses PETSc solvers
  - Scalable preconditioner

- Parallelism
  - MPI
  - GPU

---

*aCruz, Yokota, Barba, Knepley*
Vortex Method

\( t = 200 \)

- Incompressible Flow
  - Gaussian vortex blobs
  - High Re
- PetFMM
  - 2D/3D domains
  - Automatic load balancing
  - Variety of kernels
  - Optimized with templates
- PetRBF
  - Variety of RBFs
  - Uses PETSc solvers
  - Scalable preconditioner
- Parallelism
  - MPI
  - GPU

\(^a\)Cruz, Yokota, Barba, Knepley
Vortex Method
\( t = 300 \)

- Incompressible Flow
  - Gaussian vortex blobs
  - High Re

- PetFMM
  - 2D/3D domains
  - Automatic load balancing
  - Variety of kernels
  - Optimized with templates

- PetRBF
  - Variety of RBFs
  - Uses PETSc solvers
  - Scalable preconditioner

- Parallelism
  - MPI
  - GPU

---

\(^a\)Cruz, Yokota, Barba, Knepley

---
Vortex Method

\( t = 400 \)

- Incompressible Flow
  - Gaussian vortex blobs
  - High Re

- PetFMM
  - 2D/3D domains
  - Automatic load balancing
  - Variety of kernels
  - Optimized with templates

- PetRBF
  - Variety of RBFs
  - Uses PETSc solvers
  - Scalable preconditioner

- Parallelism
  - MPI
  - GPU

- \(^a\)Cruz, Yokota, Barba, Knepley
Vortex Method

\[ t = 500 \]

- Incompressible Flow
  - Gaussian vortex blobs
  - High Re
- PetFMM
  - 2D/3D domains
  - Automatic load balancing
  - Variety of kernels
  - Optimized with templates
- PetRBF
  - Variety of RBFs
  - Uses PETSc solvers
  - Scalable preconditioner
- Parallelism
  - MPI
  - GPU

\(^a\)Cruz, Yokota, Barba, Knepley
Vortex Method

\( t = 600 \)

- Incompressible Flow
  - Gaussian vortex blobs
  - High Re

- PetFMM
  - 2D/3D domains
  - Automatic load balancing
  - Variety of kernels
  - Optimized with templates

- PetRBF
  - Variety of RBFs
  - Uses PETSc solvers
  - Scalable preconditioner

- Parallelism
  - MPI
  - GPU

---

\(^a\) Cruz, Yokota, Barba, Knepley
Vortex Method

\( t = 700 \)

- Incompressible Flow
  - Gaussian vortex blobs
  - High Re

- PetFMM
  - 2D/3D domains
  - Automatic load balancing
  - Variety of kernels
  - Optimized with templates

- PetRBF
  - Variety of RBFs
  - Uses PETSc solvers
  - Scalable preconditioner

- Parallelism
  - MPI
  - GPU

\(^a\text{Cruz, Yokota, Barba, Knepley}\)
Vortex Method

Incompressible Flow
- Gaussian vortex blobs
- High Re

PetFMM
- 2D/3D domains
- Automatic load balancing
- Variety of kernels
- Optimized with templates

PetRBF
- Variety of RBFs
- Uses PETSc solvers
- Scalable preconditioner

Parallelism
- MPI
- GPU

\(^a\text{Cruz, Yokota, Barba, Knepley}\)
Real-time Surgery

- Brain Surgery
  - Elastic deformation
  - Overlaid on MRI
  - Guides surgeon

- Laser Thermal Therapy
  - PDE constrained optimization
  - Per-patient calibration
  - Thermal inverse problem

\(^a\)Warfield, Ferrant, et.al.
Real-time Surgery

- Brain Surgery
  - Elastic deformation
  - Overlaid on MRI
  - Guides surgeon

- Laser Thermal Therapy
  - PDE constrained optimization
  - Per-patient calibration
  - Thermal inverse problem

\[a\text{Fuentes, Oden, et.al.}\]
1 Getting Started with PETSc
- Who uses PETSc?
- **Stuff for Windows**
  - How can I get PETSc?
  - How do I Configure PETSc?
  - How do I Build PETSc?
  - How do I run an example?
  - How do I get more help?
Questions for Windows Users

- Have you installed cygwin?
  - Need python, make, and build-utils packages

- Will you use the GNU compilers?
  - If not, remove `link.exe`
  - If MS, check compilers from `cmd window` and use `win32fe`

- Which MPI will you use?
  - You can use `--with-mpi=0`
  - If MS, need to install MPICH2
  - If GNU, can use `--download-mpich`

- Minimal build works on Linux subsystem
Getting Started with PETSc

- Who uses PETSc?
- Stuff for Windows
- **How can I get PETSc?**
- How do I Configure PETSc?
- How do I Build PETSc?
- How do I run an example?
- How do I get more help?
Getting Started with PETSc

How can I get PETSc?

**Downloading PETSc**

- The latest tarball is on the PETSc site: 
  http://www.mcs.anl.gov/petsc/download

- There is a **Debian package** (`aptitude install petsc-dev`)

- There is a **Git development repository**
Cloning PETSc

- The full development repository is open to the public
  - [https://bitbucket.org/petsc/petsc/](https://bitbucket.org/petsc/petsc/)

- Why is this better?
  - You can clone to any release (or any specific ChangeSet)
  - You can easily rollback changes (or releases)
  - You can get fixes from us the same day
  - You can easily submit changes using a pull request

- All releases are just tags:
  - Source at tag v3.10.3
Unpacking PETSc

- Just clone development repository
  - `git clone http://bitbucket.org/petsc/petsc.git`
  - `git checkout -rv3.10.3`

  or

- Unpack the tarball
  - `tar xzf petsc.tar.gz`
Exercise 1

Download and Unpack PETSc!
Outline

1 Getting Started with PETSc
   - Who uses PETSc?
   - Stuff for Windows
   - How can I get PETSc?
   - How do I Configure PETSc?
   - How do I Build PETSc?
   - How do I run an example?
   - How do I get more help?
Configuring PETSc

- Set `$PETSC_DIR` to the installation root directory
- Run the configuration utility
  - `$PETSC_DIR/configure`
  - `$PETSC_DIR/configure --help`
  - `$PETSC_DIR/configure --download-mpich`
  - `$PETSC_DIR/configure --prefix=/usr`
- There are many examples in `$PETSC_DIR/config/examples`
- Config files in `$PETSC_DIR/$PETSC_ARCH/lib/petsc/conf`
  - Config header in `$PETSC_DIR/$PETSC_ARCH/include`
  - `$PETSC_ARCH` has a default if not specified
You can easily reconfigure with the same options

- 

Can maintain several different configurations

- ./configure -PETSC_ARCH=arch-linux-opt --with-debugging=0

All configuration information is in the logfile

- 

ALWAYS send this file with bug reports
Configuring PETSc for FEM

```bash
$PETSC_DIR/configure
  --download-triangle --download-ctetgen --download-p4est
  --download-eigen --download-pragmatic
  --download-chaco --download-metis --download-parmetis
  --download-hdf5 --download-netcdf --download-pnetcdf
  --download-exodusii --download-med
```
Getting Started with PETSc

How do I Configure PETSc?

Configuring PETSc for FEM

$PETSC_DIR/configure

–download-triangle –download-ctetgen –download-p4est
–download-eigen –download-pragmatic
–download-chaco –download-metis –download-parmetis
–download-hdf5 –download-netcdf –download-pnetcdf
–download-exodusii –download-med
$PETSC_DIR/configure
   --download-triangle --download-ctetgen --download-p4est
   --download-eigen --download-pragmatic
   --download-chaco --download-metis --download-parmetis
   --download-hdf5 --download-netcdf --download-pnetcdf
   --download-exodusii --download-med
$PETSC_DIR/configure

    --download-triangle --download-ctetgen --download-p4est
    --download-eigen --download-pragmatic
    --download-chaco --download-metis --download-parmetis
    --download-hdf5 --download-netcdf --download-pnetcdf
    --download-exodusii --download-med
$PETSC_DIR/configure

–with-cuda
–with-cudac='nvcc -m64' –with-cuda-arch=sm_10
–with-opencl
–with-opencl-include=/System/Library/Frameworks/OpenCL.framework/Headers/
–with-opencl-lib=/System/Library/Frameworks/OpenCL.framework/OpenCL
–with-precision=single
$PETSC_DIR/configure

–with-cuda
–with-cudac=’nvcc -m64’ –with-cuda-arch=sm_10
–with-opencl
–with-opencl-include=/System/Library/Frameworks/OpenCL.framework/Headers/
–with-opencl-lib=/System/Library/Frameworks/OpenCL.framework/OpenCL
–with-precision=single
$PETSC_DIR/configure

   --with-cuda
   --with-cudac='nvcc -m64' --with-cuda-arch=sm_10
   --with-opencl
      --with-opencl-include=/System/Library/Frameworks/OpenCL.framework/Headers/
      --with-opencl-lib=/System/Library/Frameworks/OpenCL.framework/OpenCL
   --with-precision=single
Configuring PETSc for Accelerators

$PETSC_DIR/configure

- `--with-cuda`
- `--with-cudac='nvcc -m64'`  `--with-cuda-arch=sm_10`
- `--with-opencl`
  - `--with-opencl-include=/System/Library/Frameworks/OpenCL.framework/Headers/`
  - `--with-opencl-lib=/System/Library/Frameworks/OpenCL.framework/OpenCL`
- `--with-precision=single`
$PETSC_DIR/configure

–with-cuda
–with-cudac=’nvcc -m64’ –with-cuda-arch=sm_10
–with-opencl
–with-opencl-include=/System/Library/Frameworks/OpenCL.framework/Headers/
–with-opencl-lib=/System/Library/Frameworks/OpenCL.framework/OpenCL
–with-precision=single
Starting in 2.2.1, some packages are automatically
  Downloaded
  Configured and Built (in $PETSC_DIR/externalpackages)
  Installed with PETSc
Currently works for
  petsc4py, mpi4py
  PETSc documentation utilities (Sowing, c2html)
  BLAS, LAPACK, Elemental, ScaLAPACK
  MPICH, OpenMPI
  ParMetis, Chaco, Jostle, Party, Scotch, Zoltan
  SuiteSparse, MUMPS, SuperLU, SuperLU_Dist, PaStiX, Pardiso
  HYPRE, ML
  BLOPEX, FFTW, STRUMPACK, SPAI, CUSP, Sundials
  Triangle, TetGen, p4est, Pragmatic
  HDF5, NetCDF, ExodusII
  AfterImage, gifLib, libjpeg, opengl
  GMP, MPFR
  ConcurrencyKit, hwloc
Exercise 2

Configure your downloaded PETSc.
Outline

1 Getting Started with PETSc
   - Who uses PETSc?
   - Stuff for Windows
   - How can I get PETSc?
   - How do I Configure PETSc?
   - How do I Build PETSc?
   - How do I run an example?
   - How do I get more help?
There is now One True Way to build PETSc:

- make
- make install if you configured with --prefix
- Check build when done with make check

Can build multiple configurations

- PETSC_ARCH=arch-linux-opt make
- Libraries are in $PETSC_DIR/$PETSC_ARCH/lib/

Complete log for each build is in logfile

- ./$PETSC_ARCH/lib/petsc/conf/make.log
- ALWAYS send this with bug reports
Build your configured PETSc.
Reconfigure PETSc to use ParMetis.

1. `linux-debug/lib/petsc/conf/reconfigure-linux-debug.py`
   - `--PETSC_ARCH=arch-linux-parmetis`
   - `--download-metis --download-parmetis`

2. `PETSC_ARCH=linux-parmetis make`

3. `PETSC_ARCH=linux-parmetis make check`
Outline

1. Getting Started with PETSc
   - Who uses PETSc?
   - Stuff for Windows
   - How can I get PETSc?
   - How do I Configure PETSc?
   - How do I Build PETSc?
   - How do I run an example?
   - How do I get more help?
Running PETSc

- Try running PETSc examples first
  - `cd $PETSC_DIR/src/snes/examples/tutorials`

- Build examples using make targets
  - `make ex5`

- Run examples using the make target
  - `make runex5`

- Can also run using MPI directly
  - `mpirun ./ex5 -snes_max_it 5`
  - `mpiexec ./ex5 -snes_monitor`
Running PETSc

- PETSc has a new test infrastructure
  - Described in Manual Section 1.3 and the Developer’s Guide
- Run all tests
  - `make PETSC_ARCH=arch-myarch test`
- Run a specific example
  - `make -f gmakefile test search='vec_vec_tutorials-ex6'`
- Run a set of similar examples
  - `make -f gmakefile test globsearch='ts*'
  - `make -f gmakefile test globsearch='ts_tutorials-ex11_*'
  - `make -f gmakefile test argsearch='cuda'`
Using MPI

- The **Message Passing Interface** is:
  - a library for parallel communication
  - a system for launching parallel jobs (mpirun/mpiexec)
  - a community standard

- Launching jobs is easy
  - `mpiexec -n 4 ./ex5`

- You should never have to make MPI calls when using PETSc
  - Almost never
Communicator

- A context (or scope) for parallel communication ("Who can I talk to")
- There are two defaults:
  - yourself (PETSC_COMM_SELF),
  - and everyone launched (PETSC_COMM_WORLD)
- Can create new communicators by splitting existing ones
- Every PETSc object has a communicator
- Set PETSC_COMM_WORLD to put all of PETSc in a subcomm

Point-to-point communication

- Happens between two processes (like in MatMult())

Reduction or scan operations

- Happens among all processes (like in VecDot())
Getting Started with PETSc

How do I run an example?

Common Viewing Options

- Gives a text representation
  - `vec_view`
- Generally views subobjects too
  - `snes_view`
- Can visualize some objects
  - `mat_view` draw::
- Alternative formats
  - `vec_view` binary:sol.bin:, `vec_view` ::matlab, `vec_view` socket
- Sometimes provides extra information
  - `mat_view` ::ascii_info, `mat_view` ::ascii_info_detailed
- Use `help` to see all options
Common Monitoring Options

- **Display the residual**
  - `-ksp_monitor`, **graphically** `-ksp_monitor_draw`

- **Can disable dynamically**
  - `-ksp_monitors_cancel`

- **Does not display subsolvers**
  - `-snes_monitor`

- **Can use the true residual**
  - `-ksp_monitor_true_residual`

- **Can display different subobjects**
  - `-snes_monitor_residual`, `-snes_monitor_solution`, `-snes_monitor_solution_update`
  - `-snes_monitor_range`
  - `-ksp_gmres_krylov_monitor`

- **Can display the spectrum**
  - `-ksp_monitor_singular_value`
Getting Started with PETSc

- Who uses PETSc?
- Stuff for Windows
- How can I get PETSc?
- How do I Configure PETSc?
- How do I Build PETSc?
- How do I run an example?
- How do I get more help?
Getting More Help

- http://www.mcs.anl.gov/petsc
- Hyperlinked documentation
  - Manual
  - Manual pages for every method
  - HTML of all example code (linked to manual pages)
- FAQ
- Full support at petsc-maint@mcs.anl.gov
- Knowledgeable users
  - David Keyes
  - Lawrence Mitchell
  - Brad Aagaard
  - Gerard Gorman
  - Paul Bauman
  - Marc Spiegelman
Outline

1. Getting Started with PETSc

2. PETSc Integration
   - Initial Operations
   - Vector Algebra
   - Matrix Algebra
   - Algebraic Solvers
   - Debugging PETSc
   - Profiling PETSc
   - Data Layout and Traversal

3. Advanced Solvers

4. More Stuff
Outline

PETSc Integration

- Initial Operations
  - Vector Algebra
  - Matrix Algebra
  - Algebraic Solvers
  - Debugging PETSc
  - Profiling PETSc
  - Data Layout and Traversal
Application Integration

- Be willing to experiment with algorithms
  - No optimality without interplay between physics and algorithmics
- Adopt flexible, extensible programming
  - Algorithms and data structures not hardwired
- Be willing to play with the real code
  - Toy models are rarely helpful
- If possible, profile before integration
  - Automatic in PETSc
PETSc is a set of library interfaces

- We do not seize `main()`
- We do not control output
- We propagate errors from underlying packages
- We present the same interfaces in:
  - C
  - C++
  - F77
  - F90
  - Python

See Gropp in *SIAM, OO Methods for Interop SciEng, ’99*
Integration Stages

- Version Control
  - It is impossible to overemphasize
  - We use Git
- Initialization
  - Linking to PETSc
- Profiling
  - Profile before changing
  - Also incorporate command line processing
- Linear Algebra
  - First PETSc data structures
- Solvers
  - Very easy after linear algebra is integrated
**Initialization**

- **Call** PetscInitialize()
  - Setup static data and services
  - Setup MPI if it is not already
- **Call** PetscFinalize()
  - Calculates logging summary
  - Shutdown and release resources
- Checks compile and link
Use `-log_view` for a performance profile
- Event timing
- Event flops
- Memory usage
- MPI messages

This used to be `-log_summary`

Call `PetscLogStagePush()` and `PetscLogStagePop()`
- User can add new stages

Call `PetscLogEventBegin()` and `PetscLogEventEnd()`
- User can add new events
Command Line Processing

- Check for an option
  - PetscOptionsHasName()

- Retrieve a value
  - PetscOptionsGetInt(), PetscOptionsGetIntArray()

- Set a value
  - PetscOptionsSetValue()

- Check for unused options
  - -options_left

- Clear, alias, reject, etc.

- Modern form uses
  - PetscOptionsBegin(), PetscOptionsEnd()
  - PetscOptionsInt(), PetscOptionsReal()
  - Integrates with -help
Outline

2 PETSc Integration

- Initial Operations
- Vector Algebra
- Matrix Algebra
- Algebraic Solvers
- Debugging PETSc
- Profiling PETSc
- Data Layout and Traversal
What are PETSc vectors?

- Fundamental objects representing solutions, right-hand sides, and coefficients.
- Each process locally owns a subvector of contiguous global data.
How do I create vectors?

- `VecCreate(MPI_Comm comm, Vec*v)`
- `VecSetSizes(Vec v, PetscInt n, PetscInt N)`
- `VecSetType(Vec v, VecType typeName)`
- `VecSetFromOptions(Vec v)`
  - Can set the type at runtime
A PETSc Vec

- Supports all vector space operations
  - VecDot(), VecNorm(), VecScale()
- Has a direct interface to the values
  - VecGetArray(), VecGetArrayF90()
- Has unusual operations
  - VecSqrtAbs(), VecStrideGather()
- Communicates automatically during assembly
- Has customizable communication (PetscSF, VecScatter)
Processes may set an arbitrary entry
  - Must use proper interface
Entries need not be generated locally
  - Local meaning the process on which they are stored
PETSc automatically moves data if necessary
  - Happens during the assembly phase
Vector Assembly

- A three step process
  - Each process sets or adds values
  - Begin communication to send values to the correct process
  - Complete the communication

\[
\text{VecSetValues}(\text{Vec } v, \text{PetscInt } n, \text{PetscInt } \text{rows}[],
\text{PetscScalar } \text{values}[], \text{InsertMode } \text{mode})
\]

- Mode is either \text{INSERT\_VALUES} or \text{ADD\_VALUES}
- Two phases allow overlap of communication and computation
  - VecAssemblyBegin(v)
  - VecAssemblyEnd(v)
One Way to Set the Elements of a Vector

```c
ierr = VecGetSize(x, &N); CHKERRQ(ierr);
ierr = MPI_Comm_rank(PETSC_COMM_WORLD, &rank); CHKERRQ(ierr);
if (rank == 0) {
    val = 0.0;
    for (i = 0; i < N; ++i) {
        ierr = VecSetValues(x, 1, &i, &val, INSERT_VALUES); CHKERRQ(ierr);
        val += 10.0;
    }
}
/* These routines ensure that the data is
   distributed to the other processes */
ierr = VecAssemblyBegin(x); CHKERRQ(ierr);
ierr = VecAssemblyEnd(x); CHKERRQ(ierr);
```
One Way to Set the Elements of a Vector

```c
VecGetSize(x, &N);
MPI_Comm_rank(PETSC_COMM_WORLD, &rank);
if (rank == 0) {
    val = 0.0;
    for (i = 0; i < N; ++i) {
        VecSetValues(x, 1, &i, &val, INSERT_VALUES);
        val += 10.0;
    }
}
/* These routines ensure that the data is distributed to the other processes */
VecAssemblyBegin(x);
VecAssemblyEnd(x);
```
A Better Way to Set the Elements of a Vector

```c
VecGetOwnershipRange(x, &low, &high);
val = low * 10.0;
for (i = low; i < high; ++i) {
    VecSetValues(x, 1, &i, &val, INSERT_VALUES);
    val += 10.0;
}
/* No data will be communicated here */
VecAssemblyBegin(x);
VecAssemblyEnd(x);
```
## Selected Vector Operations

<table>
<thead>
<tr>
<th>Function Name</th>
<th>Operation</th>
</tr>
</thead>
<tbody>
<tr>
<td>VecAXPY(Vec y, PetscScalar a, Vec x)</td>
<td>$y = y + a \cdot x$</td>
</tr>
<tr>
<td>VecAYPX(Vec y, PetscScalar a, Vec x)</td>
<td>$y = x + a \cdot y$</td>
</tr>
<tr>
<td>VecWAYPX(Vec w, PetscScalar a, Vec x, Vec y)</td>
<td>$w = y + a \cdot x$</td>
</tr>
<tr>
<td>VecScale(Vec x, PetscScalar a)</td>
<td>$x = a \cdot x$</td>
</tr>
<tr>
<td>VecCopy(Vec y, Vec x)</td>
<td>$y = x$</td>
</tr>
<tr>
<td>VecPointwiseMult(Vec w, Vec x, Vec y)</td>
<td>$w_i = x_i \cdot y_i$</td>
</tr>
<tr>
<td>VecMax(Vec x, PetscInt *idx, PetscScalar *r)</td>
<td>$r = \max r_i$</td>
</tr>
<tr>
<td>VecShift(Vec x, PetscScalar r)</td>
<td>$x_i = x_i + r$</td>
</tr>
<tr>
<td>VecAbs(Vec x)</td>
<td>$x_i =</td>
</tr>
<tr>
<td>VecNorm(Vec x, NormType type, PetscReal *r)</td>
<td>$r = |x|$</td>
</tr>
</tbody>
</table>
Working With Local Vectors

It is sometimes more efficient to directly access local storage of a `Vec`

- PETSc allows you to access the local storage with
  - `VecGetArray(Vec, double *[])`

- You must return the array to PETSc when you finish
  - `VecRestoreArray(Vec, double *[])`

- Allows PETSc to handle data structure conversions
  - Commonly, these routines are fast and do not involve a copy
VecGetArray in C

```c
Vec v;
PetscScalar *array;
PetscInt n, i;

VecGetArray(v, &array);
VecGetLocalSize(v, &n);
PetscSynchronizedPrintf(PETSC_COMM_WORLD,
    "First element of local array is %f\n", array[0]);
PetscSynchronizedFlush(PETSC_COMM_WORLD);
for (i = 0; i < n; ++i) {
    array[i] += (PetscScalar) rank;
}
VecRestoreArray(v, &array);
```
VecGetArray in F77

```
#include "finclude/petsc.h"

Vec v;
PetscScalar array(1)
PetscOffset offset
PetscInt n, i
PetscErrorCode ierr

call VecGetArray(v, array, offset, ierr)
call VecGetLocalSize(v, n, ierr)
do i=1,n
    array(i+offset) = array(i+offset) + rank
end do
call VecRestoreArray(v, array, offset, ierr)
```
VecGetArray in F90

```c
#include "finclude/petsc.h90"

Vec v;
PetscScalar pointer :: array(:)
PetscInt n, i
PetscErrorCode ierr

call VecGetArrayF90(v, array, ierr)
call VecGetLocalSize(v, n, ierr)
do i=1,n
    array(i) = array(i) + rank
end do
call VecRestoreArrayF90(v, array, ierr)
```
VecGetArray in Python

```
with v as a:
    for i in range(len(a)):
        a[i] = 5.0*i
```
```c
DM da;
Vec v;
DMDALocalInfo *info;
PetscScalar **array;

DMDAVecGetArray(da, v, &array);
for (j = info->ys; j < info->ys+info->ym; ++j) {
    for (i = info->xs; i < info->xs+info->xm; ++i) {
        u = x[j][i];
        uxx = (2.0*u - x[j][i-1] - x[j][i+1])*hydhx;
        uyy = (2.0*u - x[j-1][i] - x[j+1][i])*hxdhy;
        f[j][i] = uxx + uyy;
    }
}
DMDAVecRestoreArray(da, v, &array);
```
PETSc Integration  Vector Algebra

DMDAVecGetArray in F90

```fortran
DM da
Vec v
PetscScalar,pointer :: array(:,:,:)

call DMDAGetCorners(da,xs,ys,PETSC_NULL_INTEGER,
xm,ym,PETSC_NULL_INTEGER,ierr)
call DMDAVecGetArrayF90(da,v,array,ierr);
do i = xs,xs+xm
   do j = ys,ys+ym
      u = x(i,j)
      uxx = (2.0*u - x(i-1,j) - x(i+1,j))*hydhx;
      uyy = (2.0*u - x(i,j-1) - x(i,j+1)*hxdhy;
      f(i,j) = uxx + uyy;
   enddo
endo
call DMDAVecRestoreArrayF90(da,v,array,ierr);
```

Matt (ANL)
Outline

2 PETSc Integration
- Initial Operations
- Vector Algebra
- **Matrix Algebra**
  - Algebraic Solvers
  - Debugging PETSc
  - Profiling PETSc
  - Data Layout and Traversal
What are PETSc matrices?

- Fundamental objects for storing stiffness matrices and Jacobians
- Each process locally owns a contiguous set of rows
- Supports many data types
  - AIJ, Block AIJ, Symmetric AIJ, Block Matrix, etc.
- Supports structures for many packages
  - Elemental, MUMPS, SuperLU, UMFPack, PasTiX
How do I create matrices?

- **MatCreate(MPI_Comm comm, Mat *A)**
- **MatSetSizes(Mat A, PetscInt m, PetscInt n, PetscInt M, PetscInt N)**
- **MatSetType(Mat A, MatType typeName)**
- **MatSetFromOptions(Mat A)**
  - Can set the type at runtime
- **MatSeqAIJPreallocation(Mat A, PetscInt nz, const PetscInt nnz[])**
- **MatXAIJPreallocation(Mat A, bs, dnz[], onz[], dnzu[], onzu[])**
- **MatSetValues(Mat A, m, rows[], n, cols [], values [], InsertMode)**
  - **MUST** be used, but does automatic communication
Matrix Polymorphism

The PETSc \texttt{Mat} has a single user interface,
\begin{itemize}
  \item Matrix assembly
    \begin{itemize}
      \item \texttt{MatSetValues()}
      \item \texttt{MatGetLocalSubMatrix()}
    \end{itemize}
  \item Matrix-vector multiplication
    \begin{itemize}
      \item \texttt{MatMult()}
    \end{itemize}
  \item Matrix viewing
    \begin{itemize}
      \item \texttt{MatView()}
    \end{itemize}
\end{itemize}
but multiple underlying implementations.
\begin{itemize}
  \item AIJ, Block AIJ, Symmetric Block AIJ,
  \item Dense
  \item Matrix-Free
  \item etc.
\end{itemize}
A matrix is defined by its \texttt{interface}, not by its \texttt{data structure}.
Matrix Assembly

- A three step process
  - Each process sets or adds values
  - Begin communication to send values to the correct process
  - Complete the communication

  `MatSetValues(A, m, rows[], n, cols [], values [], mode)`
  - `mode` is either `INSERT_VALUES` or `ADD_VALUES`
  - Logically dense block of values

- Two phase assembly allows overlap of communication and computation
  - `MatAssemblyBegin(A, type)`
  - `MatAssemblyEnd(A, type)`
  - `type` is either `MAT_FLUSH_ASSEMBLY` or `MAT_FINAL_ASSEMBLY`
One Way to Set the Elements of a Matrix

Simple 3-point stencil for 1D Laplacian

```c
v[0] = -1.0; v[1] = 2.0; v[2] = -1.0;
if (rank == 0) {
    for (row = 0; row < N; row++) {
        cols[0] = row - 1; cols[1] = row; cols[2] = row + 1;
        if (row == 0) {
            MatSetValues(A, 1, &row, 2, &cols[1], &v[1], INSERT_VALUES);
        } else if (row == N - 1) {
            MatSetValues(A, 1, &row, 2, cols, v, INSERT_VALUES);
        } else {
            MatSetValues(A, 1, &row, 3, cols, v, INSERT_VALUES);
        }
    }
}
MatAssemblyBegin(A, MAT_FINAL_ASSEMBLY);
MatAssemblyEnd(A, MAT_FINAL_ASSEMBLY);
```
Parallel Sparse Matrix Layout

- **proc 0**: diagonal blocks
- **proc 1**: offdiagonal blocks
- **proc 2**: diagonal blocks
- **proc 3**: offdiagonal blocks
- **proc 4**: diagonal blocks
- **proc 5**: offdiagonal blocks

**Legend**
- Blue: diagonal blocks
- Green: offdiagonal blocks
A Better Way to Set the Elements of a Matrix
Simple 3-point stencil for 1D Laplacian

\[ v[0] = -1.0; \quad v[1] = 2.0; \quad v[2] = -1.0; \]

\text{MatGetOwnershipRange}(A, \&\text{start}, \&\text{end});
\text{for}(row = \text{start}; \quad row < \text{end}; \quad row++) {
  cols[0] = row - 1; \quad cols[1] = row; \quad cols[2] = row + 1;
  \text{if}(row == 0) {
    \text{MatSetValues}(A, 1, \&row, 2, \&cols[1], \&v[1], \text{INSERT_VALUES});
  } \quad \text{else if}(row == N-1) {
    \text{MatSetValues}(A, 1, \&row, 2, cols, v, \text{INSERT_VALUES});
  } \quad \text{else} {
    \text{MatSetValues}(A, 1, \&row, 3, cols, v, \text{INSERT_VALUES});
  }
}
\text{MatAssemblyBegin}(A, \text{MAT_FINAL_ASSEMBLY});
\text{MatAssemblyEnd}(A, \text{MAT_FINAL_ASSEMBLY});
Why Are PETSc Matrices That Way?

- No one data structure is appropriate for all problems
  - Blocked and diagonal formats provide performance benefits
  - PETSc has many formats
  - Makes it easy to add new data structures

- Assembly is difficult enough without worrying about partitioning
  - PETSc provides parallel assembly routines
  - High performance still requires making most operations local
  - However, programs can be incrementally developed.
  - `MatPartitioning` and `MatOrdering` can help
  - It’s better to partition and reorder the underlying grid

- Matrix decomposition in contiguous chunks is simple
  - Makes interoperation with other codes easier
  - For other ordering, PETSc provides “Application Orderings” (AO)
Outline

2. PETSc Integration
   - Initial Operations
   - Vector Algebra
   - Matrix Algebra
   - Algebraic Solvers
     - Debugging PETSc
     - Profiling PETSc
     - Data Layout and Traversal
Experimentation is Essential!

Proof is not currently enough to examine solvers


Using PETSc linear algebra, just add:

- KSPSetOperators(ksp, A, M, flag)
- KSPSolve(ksp, b, x)

Can access subobjects

- KSPGetPC(ksp, &pc)

Preconditioners must obey PETSc interface

- Basically just the KSP interface

Can change solver dynamically from the command line

- -ksp_type bicgstab
Nonlinear Solvers

Using PETSc linear algebra, just add:
- `SNESSetFunction(snes, r, residualFunc, ctx)`
- `SNESSetJacobian(snes, A, M, jacFunc, ctx)`
- `SNESolve(snes, b, x)`

Can access subobjects
- `SNESGetKSP(snes, &ksp)`

Can customize subobjects from the cmd line
- Set the subdomain preconditioner to ILU with `−sub_pc_type ilu`
Basic Solver Usage

Use `SNESSetFromOptions()` so that everything is set dynamically

- **Set the type**
  - Use `−snes_type` (or take the default)

- **Set the preconditioner**
  - Use `−npc_snes_type` (or take the default)

- **Override the tolerances**
  - Use `−snes_rtol` and `−snes_atol`

- **View the solver to make sure you have the one you expect**
  - Use `−snes_view`

- **For debugging, monitor the residual decrease**
  - Use `−snes_monitor`
  - Use `−ksp_monitor` to see the underlying linear solver
Complete table of solvers

Sequential LU
- ESSL (IBM)
- SuperLU (Sherry Li, LBNL)
- Suitesparse (Tim Davis, U. of Florida)
- LUSOL (MINOS, Michael Saunders, Stanford)
- PILUT (Hypre, David Hysom, LLNL)

Parallel LU
- Elemental/Clique (Jack Poulson, Google)
- MUMPS (Patrick Amestoy, IRIT)
- SuperLU_Dist (Jim Demmel and Sherry Li, LBNL)
- Pardiso (MKL, Intel)
- STRUMPACK (Pieter Ghysels, LBNL)

Parallel Cholesky
- Elemental (Jack Poulson, Google)
- DSCPACK (Padma Raghavan, Penn. State)
- MUMPS (Patrick Amestoy, Toulouse)
Complete table of solvers

- Parallel Algebraic Multigrid
  - GAMG (Mark Adams, LBNL)
  - BoomerAMG (Hypre, LLNL)
  - ML (Trilinos, Ray Tuminaro and Jonathan Hu, SNL)
- Parallel BDDC (Stefano Zampini, KAUST)
- Parallel ILU, PaStiX (Faverge Mathieu, INRIA)
- Parallel Redistribution (Dave May, Oxford and Patrick Sanan, USI)
- Parallel Sparse Approximate Inverse
  - Parasails (Hypre, Edmund Chow, LLNL)
  - SPAI 3.0 (Marcus Grote and Barnard, NYU)
MPI_Comm  comm;
SNES  snes;
DM  dm;
Vec  u;

SNESCreate(comm, &snes);
SNESSetDM(snes, dm);
SNESSetFromOptions(snes);
DMCreateGlobalVector(dm, &u);
SNESSolve(snes, NULL, u);
Solver use in SNES ex62

Solver code does not change for different algorithms:

```c
SNES snes;
DM dm;
Vec u;
PetscErrorCode ierr;

ierr = SNESCreate(PETSC_COMM_WORLD, &snes);CHKERRQ(ierr);
ierr = SNESSetDM(snes, dm);CHKERRQ(ierr);
/* Specify residual computation */
ierr = SNESSetFromOptions(snes);CHKERRQ(ierr); /* Configure solver */
ierr = DMCreateGlobalVector(dm, &u);CHKERRQ(ierr);
ierr = SNES Solve(snes, PETSC_NULL, u);CHKERRQ(ierr);
```

- Never recompile! all configuration is dynamic
- DM controls data layout and communication
- Type of nested solvers can be changed at runtime
Solver use in SNES ex62

I will omit error checking and declarations:

```c
SNESCreate(PETSC_COMM_WORLD, &snes);
SNESSetDM(snes, dm);
/* Specify residual computation */
SNESSetFromOptions(snes); /* Configure solver */
DMCreateGlobalVector(dm, &u);
SNESSolve(snes, PETSC_NULL, u);
```
**PETSc Integration**

**Algebraic Solvers**

**Solver use in SNES ex62**

The configuration API can also be used:

```c
SNESCreate(PETSC_COMM_WORLD, &snes);
SNESSetDM(snes, dm);
/* Specify residual computation */
SNESNGMRESRestartType(snes, SNES_NGMRES_RESTART_PERIODIC);
SNESSetFromOptions(snes);
DMCreateGlobalVector(dm, &u);
SNESolve(snes, PETSC_NULL, u);
```

- Ignored when not applicable (no ugly check)
- Type safety of arguments is retained
- No downcasting
Adding a prefix namespaces command line options:

```c
SNESCreate(PETSC_COMM_WORLD, &snes);
SNESSetDM(snes, dm);
/* Specify residual computation */
SNESSetOptionsPrefix(snes, "stokes_");
SNESSetFromOptions(snes);
DMCreateGlobalVector(dm, &u);
SNES Solve(snes, PETSC_NULL, u);
```

- `-stokes_snes_type qn` changes the solver type,

whereas `-snes_type qn` does not
User provides a function to compute the residual:

\[
r = F(u)
\]

- User handles parallel communication
- User handles domain geometry and discretization
PETSc Integration  Algebraic Solvers

Solver use in SNES ex62

DM allows the user to compute only on a local patch:

```c
SNESCreate(PETSC_COMM_WORLD, &snes);
SNESSetDM(snes, dm);
SNESSetFromOptions(snes);
DMCreateGlobalVector(dm, &u);
SNESolve(snes, PETSC_NULL, u);
DMSNESSetLocalFunction(dm, FormFunctionLocal);
```

- Code looks serial to the user
- PETSc handles global residual assembly
- Also works for unstructured meshes
Solver use in SNES ex62

Optionally, the user can also provide a Jacobian:

```c
SNESCreate(PETSC_COMM_WORLD, &snes);
SNESSetDM(snes, dm);
SNESSetFromOptions(snes);
DMCreateGlobalVector(dm, &u);
SNESolve(snes, PETSC_NULL, u);

DMSNESSetLocalFunction(dm, FormFunctionLocal);
DMSNESSetLocalJacobian(dm, FormJacobianLocal);
```

SNES ex62 allows both
- finite difference (JFNK), and
- FEM action
versions of the Jacobian.
Solver use in SNES ex62

Convenience form uses Plex defaults:

```c
SNESCreate(PETSC_COMM_WORLD, &snes);
SNESSetDM(snes, dm);
SNESSetFromOptions(snes);
DMCreateGlobalVector(dm, &u);
SNESolve(snes, PETSC_NULL, u);

DM<PlexSetSNESLocalFEM(dm,&user,&user,&user);
```

This also handles Dirichlet boundary conditions.
The DM also handles storage:

```c
CreateMesh(PETSC_COMM_WORLD, &user, &dm);
DMCreateLocalVector(dm, &lu);
DMCreateGlobalVector(dm, &u);
DMCreateMatrix(dm, &J);
```

- DM can create local and global vectors
- Matrices are correctly preallocated
- Easy supported for discretization
Outline

PETSc Integration
- Initial Operations
- Vector Algebra
- Matrix Algebra
- Algebraic Solvers

Debugging PETSc
- Profiling PETSc
- Data Layout and Traversal
Correctness Debugging

- Automatic generation of tracebacks
- Detecting memory corruption and leaks
- Optional user-defined error handlers
Interacting with the Debugger

- Launch the debugger
  - `start_in_debugger [gdb, dbx, noxterm]`
  - `on_error_attach_debugger [gdb, dbx, noxterm]`

- Attach the debugger only to some parallel processes
  - `debugger_nodes 0,1`

- Set the display (often necessary on a cluster)
  - `display khan.mcs.anl.gov:0.0`
Debugging Tips

- Put a breakpoint in `PetscError()` to catch errors as they occur
- PETSc tracks memory overwrites at both ends of arrays
  - The `CHKMEMQ` macro causes a check of all allocated memory
  - Track memory overwrites by bracketing them with `CHKMEMQ`
- PETSc checks for leaked memory
  - Use `PetscMalloc()` and `PetscFree()` for all allocation
  - Print unfreed memory on `PetscFinalize()` with `-malloc_dump`
- Simply the best tool today is `valgrind`
  - It checks memory access, cache performance, memory usage, etc.
  - [http://www.valgrind.org](http://www.valgrind.org)
  - Need `--trace-children=yes` when running under MPI
Use the debugger to find a SEGV
Locate a memory overwrite using CHKMEMQ.

- Get the example
  - hg clone -r1 http://petsc.cs.iit.edu/petsc/SimpleTutorial
- Build the example `make`
- Run it and watch the fireworks
  - mpiexec -n 2 ./bin/ex5 -use_coords
- Run it under the debugger and correct the error
  - mpiexec -n 2 ./bin/ex5 -use_coords -start_in_debugger
  - hg update -r2
- Build it and run again smoothly
Outline

2. PETSc Integration
   - Initial Operations
   - Vector Algebra
   - Matrix Algebra
   - Algebraic Solvers
   - Debugging PETSc
   - Profiling PETSc
   - Data Layout and Traversal
PETSc has integrated profiling
- Option \(-\text{log\_view}\) prints a report on \(\text{PetscFinalize()}\)

PETSc allows user-defined events
- Events report time, calls, flops, communication, etc.
- Memory usage is tracked by object

Profiling is separated into stages
- Event statistics are aggregated by stage
Using Stages and Events

- **Use** PetscLogStageRegister() to create a new stage
  - Stages are identifier by an integer handle

- **Use** PetscLogStagePush/Pop() to manage stages
  - Stages may be nested, but will not aggregate in a nested fashion

- **Use** PetscLogEventRegister() to create a new stage
  - Events also have an associated class

- **Use** PetscLogEventBegin/End() to manage events
  - Events may also be nested and will aggregate in a nested fashion
  - Can use PetscLogFlops() to log user flops
int stageNum;

PetscLogStageRegister(&stageNum, "name");
PetscLogStagePush(stageNum);

/* Code to Monitor */
PetscLogStagePop();
PETSc Integration

Profiling PETSc

Adding A Logging Stage

Python

```
with PETSc.LogStage('Fluid Stage') as fluidStage:
    # All operations will be aggregated in fluidStage
    fluid.solve()
```
static int USER_EVENT;

PetscLogEventRegister(&USER_EVENT, "name", CLS_ID);
PetscLogEventBegin(USER_EVENT, 0, 0, 0, 0);

/* Code to Monitor */

PetscLogFlops(user_event_flops);
PetscLogEventEnd(USER_EVENT, 0, 0, 0, 0);
with PETSc.logEvent('Reconstruction') as recEvent:
    # All operations are timed in recEvent
    reconstruct(sol)
    # Flops are logged to recEvent
    PETSc.Log.logFlops(user_event_flops)
Adding A Logging Class

```c
static int CLASS_ID;
PetscLogClassRegister(&CLASS_ID, "name");
```

- Class ID identifies a class uniquely
- Must initialize before creating any objects of this type
PETSc Integration
Profiling PETSc

Matrix Memory Preallocation

- PETSc sparse matrices are dynamic data structures
  - can add additional nonzeros freely
- Dynamically adding many nonzeros
  - requires additional memory allocations
  - requires copies
  - can kill performance
- Memory preallocation provides
  - the freedom of dynamic data structures
  - good performance
- Easiest solution is to replicate the assembly code
  - Remove computation, but preserve the indexing code
  - Store set of columns for each row
- Call preallocation routines for all datatypes
  - \texttt{MatSeqAIJSetPreallocation()}
  - \texttt{MatMPIAIJSetPreallocation()}
  - Only the relevant data will be used
Matrix Memory Preallocation
Sequential Sparse Matrices

**MatSeqAIJPreallocation**(

\textbf{MatA}, \textbf{int} \textbf{nz}, \textbf{int} \textbf{nnz}[])

- **nz**: expected number of nonzeros in any row
- **nnz(i)**: expected number of nonzeros in row \(i\)
Each process locally owns a submatrix of contiguous global rows.

Each submatrix consists of diagonal and off-diagonal parts.

MatGetOwnershipRange(MatA, int *start, int *end)

**start**: first locally owned row of global matrix

**end-1**: last locally owned row of global matrix
Matrix Memory Preallocation
Parallel Sparse Matrices

MatMPIAIJPreallocation(MatA, int dnz, int dnnz[], int onz, int onnz[])

 dnz: expected number of nonzeros in any row in the diagonal block
 dnnz(i): expected number of nonzeros in row i in the diagonal block
 onz: expected number of nonzeros in any row in the offdiagonal portion
 onnz(i): expected number of nonzeros in row i in the offdiagonal portion
Matrix Memory Preallocation
Verifying Preallocation

- Use runtime option `-info`
- Output:
  ```plaintext
  [proc #] Matrix size: %d X %d; storage space: %d unneeded, %d used
  [proc #] Number of mallocs during MatSetValues() is %d
  ```

```
[merlin] mpirun ex2 -log_info
[0]MatAssemblyEnd_SeqAIJ: Matrix size: 56 X 56; storage space:
[0] 310 unneeded, 250 used
[0]MatAssemblyEnd_SeqAIJ: Number of mallocs during MatSetValues() is 0
[0]MatAssemblyEnd_SeqAIJ: Most nonzeros in any row is 5
[0]MatAIJ_CheckInode: Found 56 nodes out of 56 rows. Not using Inode routine
[0]MatAIJ_CheckInode: Found 56 nodes out of 56 rows. Not using Inode routine
Norm of error 0.000156044 iterations 6
[0]PetscFinalize: PETSc successfully ended!
```
Exercise 8

Return to Exercise 7 and add more profiling.

- Update to the next revision
  - hg update -r3
- Build, run, and look at the profiling report
  - make ex5
  - ./bin/ex5 -use_coords -log_summary
- Add a new stage for setup
- Add a new event for FormInitialGuess() and log the flops
- Build it again and look at the profiling report
Outline

2 PETSc Integration

- Initial Operations
- Vector Algebra
- Matrix Algebra
- Algebraic Solvers
- Debugging PETSc
- Profiling PETSc

Data Layout and Traversal
PETSc includes several tools for parallel data layout and traversal:

- **PetscSection**
  Data layout

- **PetscSF**
  Sharing and communication

- **DM**
  Topology and traversal
**Data Layout**

**PetscSection** defines a data layout

- maps $p \rightarrow (\text{off}, \text{off} + 1, \ldots, \text{off} + \text{dof})$
- where $p \in [\text{pStart}, \text{pEnd})$, called the **chart**
- ranges can be divided into parts, called **fields**
- prefix sums calculated automatically on setup
**Data Layout**

**PetscSection** defines a data layout

- `PetscSectionGetOffset()`, `PetscSectionGetDof()`
- where $p \in [p_{\text{Start}}, p_{\text{End}})$, called the *chart*
- ranges can be divided into parts, called *fields*
- prefix sums calculated automatically on setup
**Data Layout**

**PetscSection** defines a data layout

- maps \( p \rightarrow (off, off + 1, \ldots, off + dof) \)
- where \( p \in [pStart, pEnd) \), called the *chart*
- ranges can be divided into parts, called *fields*
- prefix sums calculated automatically on setup
**Data Layout**

**PetscSection** defines a data layout

- maps \( p \rightarrow (off, off + 1, \ldots, off + dof) \)

- \texttt{PetscSectionGetChart()}

- ranges can be divided into parts, called *fields*

- prefix sums calculated automatically on setup
**PetScSection** defines a data layout

- maps $p \rightarrow (off, off + 1, \ldots, off + dof)$
- where $p \in [pStart, pEnd)$, called the *chart*
- ranges can be divided into parts, called *fields*
- prefix sums calculated automatically on setup
**Data Layout**

**PetscSection** defines a data layout

- maps \( p \rightarrow (off, off + 1, \ldots, off + dof) \)
- where \( p \in [pStart, pEnd) \), called the chart

- `PetscSectionGetFieldOffset()`, `PetscSectionGetFieldDof()`

- prefix sums calculated automatically on setup
**Data Layout**

**PetscSection** defines a data layout

- maps $p \rightarrow (off, off + 1, \ldots, off + dof)$
- where $p \in [pStart, pEnd)$, called the *chart*
- ranges can be divided into parts, called *fields*
- prefix sums calculated automatically on setup
**Data Layout**

**PetscSection** defines a data layout

- maps $p \rightarrow (off, off + 1, \ldots, off + dof)$
- where $p \in [pStart, pEnd)$, called the chart
- ranges can be divided into parts, called fields

```c
PetscSectionSetUp()
```
**PetscSection** maps \( point \rightarrow (size, offset) \)

- If points are *processes*, it is **PetscLayout**
  - Could also be used for multicore layout

- Boundary conditions are just another **PetscSection**
  - Map points to number of constrained dofs
  - Offsets into integer array of constrained local dofs

- Fields are just another **PetscSection**
  - Map points to number of field dofs
  - Offsets into array with all fields
Decouples Mesh From Discretization
- Mesh does not need to know how dofs are generated, just how many are attached to each point.
- It does not matter whether you use FD, FVM, FEM, etc.

Decouples Mesh from Solver
- Solver gets the data layout and partitioning from Vec and Mat, nothing else from the mesh.
- Solver gets restriction/interpolation matrices from DM.

Decouples Discretization from Solver
- Solver gets the field division and blocking from Section
**PetscSection** can be used to segment data

- Use **Vec** and **IS** to store data
- Use point \( p \) instead of index \( i \)
- Maps to a set of values instead of just one

We provide a convenience method for extraction

```c
VecGetValuesSection(Vec v, PetscSection s, PetscInt p, PetscScalar **a);
```

which works in an analogous way to

```c
MatSetValuesStencil(Mat A, PetscInt nr, const MatStencil rs[],
                     PetscInt nc, const MatStencil cs[],
                     const PetscScalar v[], InsertMode m);
```
**PetscSection** can be used to segment data

- Use **Vec** and **IS** to store data
- Use point \( p \) instead of index \( i \)
- Maps to a set of values instead of just one

We can get the layout of coordinates over the mesh

```c
DMPlexGetCoordinateSection(DM dm, PetscSection *s);
```

where the data is stored in a **Vec**

```c
DMGetCoordinates(DM dm, Vec *coords);
```
**PetscSection** can be used to segment data

- Use **Vec** and **IS** to store data
- Use point \( p \) instead of index \( i \)
- Maps to a set of values instead of just one

We can retrieve FEM data from vector without complicated indexing,

\[
\text{DMPlexVecGetClosure}( \text{DM} \ dm, \text{PetscSection} \ s, \text{Vec} \ v, \\
\text{PetscInt} \ cell, \text{PetscInt} \ *, \text{PetscScalar} \ *a[]);
\]

and the same thing works for matrices

\[
\text{DMPlexMatSetClosure}( \text{DM} \ dm, \text{PetscSection} \ rs, \text{PetscSection} \ cs, \text{Mat} \ A, \\
\text{PetscInt} \ p, \text{const} \text{PetscScalar} \ v[], \text{InsertMode} \ m);
\]
High Level Interface

```c
DMPlexCreateSection(
    DM dm, PetscInt dim, PetscInt numFields,
    PetscInt numComp[], PetscInt numDof[],
    PetscInt numBC, PetscInt bcField[], IS bcPoints[],
    PetscSection *section);
```

<table>
<thead>
<tr>
<th>Discretization</th>
<th>Dof/Dimension</th>
</tr>
</thead>
<tbody>
<tr>
<td>$P_1 - P_0$</td>
<td>$[3 \ 0 \ 0 \ 0</td>
</tr>
<tr>
<td>$Q_2 - Q_1$</td>
<td>$[3 \ 3 \ 3 \ 3</td>
</tr>
<tr>
<td>$Q_2 - P_1^{\text{disc}}$</td>
<td>$[3 \ 3 \ 3 \ 3</td>
</tr>
</tbody>
</table>
Low Level Interface

PetscSectionCreate(PETSC_COMM_WORLD, &s);
PetscSectionSetNumFields(s, 2);
PetscSectionSetFieldComponents(s, 0, 3);
PetscSectionSetFieldComponents(s, 1, 1);
PetscSectionSetChart(s, cStart, vEnd);
for(PetscInt v = vStart; v < vEnd; ++v) {
    PetscSectionSetDof(s, v, 3);
    PetscSectionSetFieldDof(s, v, 0, 3);
}
for(PetscInt c = cStart; c < cEnd; ++c) {
    PetscSectionSetDof(s, c, 1);
    PetscSectionSetFieldDof(s, c, 1, 1);
}
PetscSectionSetUp(s);
**PetscSF** encodes a *star forest*:

- one-way communication pattern
- arbitrary datatype or struct
- message and one-sided implementations
- automatically builds two-sided info
We use PetscSF to describe shared points. Composing a point PetscSF and PetscSection, we can build

- a global section
- a PetscSF for shared dofs

This *composability* means we can build hierarchies of sections and pieces of sections.
We use **PetscSF** to describe shared points

Composing a point **PetscSF** and **PetscSection**, we can build

- `PetscSectionCreateGlobalSection()`

- a **PetscSF** for shared dofs

This *composability* means we can build hierarchies of sections and pieces of sections.
We use **PetscSF** to describe shared points

Composing a point **PetscSF** and **PetscSection**, we can build

- a global section
- a **PetscSF** for shared dofs

This *composability* means we can build hierarchies of sections and pieces of sections.
We use **PetscSF** to describe shared points

Composing a point **PetscSF** and **PetscSection**, we can build

- a global section
  
  **PetscSFCreateSectionSF()**

This *composability* means we can build hierarchies of sections and pieces of sections.
Communication Automation

<table>
<thead>
<tr>
<th>Point Space</th>
<th>Dof Space</th>
<th>Section</th>
<th>SF</th>
</tr>
</thead>
<tbody>
<tr>
<td>Solution Dofs</td>
<td>Adjacent Dofs</td>
<td>Jacobian Layout $\rightarrow$ Shared Adjacency</td>
<td></td>
</tr>
<tr>
<td>Mesh Points</td>
<td>Solution Dofs</td>
<td>Data Layout $\rightarrow$ Shared Dofs</td>
<td></td>
</tr>
<tr>
<td>Mesh Points</td>
<td>Mesh Points</td>
<td>Topology $\rightarrow$ Shared Topology</td>
<td></td>
</tr>
<tr>
<td>Processes</td>
<td>Mesh Points</td>
<td>Point Partition $\rightarrow$ Shared Points</td>
<td></td>
</tr>
<tr>
<td>Processes</td>
<td></td>
<td>Neighbors</td>
<td></td>
</tr>
</tbody>
</table>
The **DM** encodes point topology for traversal

- **DMDA**
  Cartesian grids, collocated layout
- **DMStag**
  Staggered grids, Section layout
- **DMplex**
  Arbitrary topology, Section layout
- **DMForest**
  Adaptive octree, Section layout
- **DMNetwork**
  Graph topology, Section layout
- **DMSwarm**
  Particles, struct/particle layout
1. Getting Started with PETSc
2. PETSc Integration
3. Advanced Solvers
   - Fieldsplit
   - Multigrid
   - Nonlinear Preconditioning
4. More Stuff
Outline

3 Advanced Solvers
- Fieldsplit
- Multigrid
- Nonlinear Preconditioning
FieldSplit Preconditioner

- **Analysis**
  - Use **IS**es to define **fields**
  - Decouples **PC** from problem definition

- **Synthesis**
  - Additive, Multiplicative, Schur
  - Commutes with Multigrid
FieldSplit Customization

Analysis
- `pc_fieldsplit_<split num>_fields 2,1,5`
- `pc_fieldsplit_detect_saddle_point`

Synthesis
- `pc_fieldsplit_type <additive, multiplicative, schur>`
- `pc_fieldsplit_diag_use_amat`
- `pc_fieldsplit_off_diag_use_amat`
  Use diagonal blocks of operator to build PC

Schur complements
- `pc_fieldsplit_schur_precondition <user,all,full,self,selfp>`
  How to build preconditioner for $S$
- `pc_fieldsplit_schur_factorization_type <diag,lower,upper,full>`
  Which off-diagonal parts of the block factorization to use
ex62: $P_2/P_1$ Stokes Problem on Unstructured Mesh

\[
\begin{pmatrix}
A & B \\
B^T & 0
\end{pmatrix}
\]
**Solver Configuration: No New Code**

**ex62**: $P_2/P_1$ Stokes Problem on Unstructured Mesh


- `ksp_type gmres` - `pc_type fieldsplit` - `pc_fieldsplit_type additive`
- `fieldsplit_velocity_ksp_type preonly` - `fieldsplit_velocity_pc_type lu`
- `fieldsplit_pressure_ksp_type preonly` - `fieldsplit_pressure_pc_type jacobi`

\[
\begin{pmatrix}
A & 0 \\
0 & I
\end{pmatrix}
\]
ex62: $P_2/P_1$ Stokes Problem on Unstructured Mesh


```plaintext
-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type additive
-fieldsplit_velocity_ksp_type preonly -fieldsplit_velocity_pc_type gamg
-fieldsplit_pressure_ksp_type preonly -fieldsplit_pressure_pc_type jacobi
```

\[
\begin{pmatrix}
\hat{A} & 0 \\
0 & I
\end{pmatrix}
\]
ex62: $P_2/P_1$ Stokes Problem on Unstructured Mesh


-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type multiplicative
-fieldsplit_velocity_ksp_type preonly -fieldsplit_velocity_pc_type gamg
-fieldsplit_pressure_ksp_type preonly -fieldsplit_pressure_pc_type jacobi

$\begin{pmatrix} \hat{A} & B \\ 0 & I \end{pmatrix}$
ex62: $P_2/P_1$ Stokes Problem on Unstructured Mesh


-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type multiplicative
- pc_fieldsplit_0_fields 1 - pc_fieldsplit_1_fields 0
- fieldsplit_velocity_ksp_type preonly - fieldsplit_velocity_pc_type gamg
- fieldsplit_pressure_ksp_type preonly - fieldsplit_pressure_pc_type jacobi

$$
\begin{pmatrix}
I & B^T \\
0 & \hat{A}
\end{pmatrix}
$$
Solver Configuration: No New Code

**ex62**: $P_2/P_1$ Stokes Problem on Unstructured Mesh


```
-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type schur
   -pc_fieldsplit_schur_factorization_type diag
   -fieldsplit_velocity_ksp_type preonly -fieldsplit_velocity_pc_type gamg
   -fieldsplit_pressure_ksp_type minres -fieldsplit_pressure_pc_type none
```

\[
\begin{pmatrix}
\hat{A} & 0 \\
0 & \hat{S}
\end{pmatrix}
\]
ex62: $P_2/P_1$ Stokes Problem on Unstructured Mesh

Lower Schur Complement, May and Moresi, PEPI, 2008.

-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type schur
-pc_fieldsplit_schur_factorization_type lower
-fieldsplit_velocity_ksp_type preonly -fieldsplit_velocity_pc_type gamg
-fieldsplit_pressure_ksp_type minres -fieldsplit_pressure_pc_type none

\[
\begin{pmatrix}
\hat{A} & 0 \\
B^T & \hat{S}
\end{pmatrix}
\]
Solver Configuration: No New Code

**ex62**: $P_2/P_1$ Stokes Problem on Unstructured Mesh


-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type schur -pc_fieldsplit_schur_factorization_type upper
-fieldsplit_velocity_ksp_type preonly -fieldsplit_velocity_pc_type gamg -fieldsplit_pressure_ksp_type minres -fieldsplit_pressure_pc_type none

$$
\begin{pmatrix}
\hat{A} & B \\
\hat{S} &
\end{pmatrix}
$$
**Solver Configuration: No New Code**

**ex62:** $P_2/P_1$ Stokes Problem on Unstructured Mesh

Uzawa Iteration, Uzawa, 1958

-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type schur
-pc_fieldsplit_schur_factorization_type upper
-fieldsplit_velocity_ksp_type preonly -fieldsplit_velocity_pc_type lu
-fieldsplit_pressure_ksp_type richardson -fieldsplit_pressure_pc_type jac
-fieldsplit_pressure_ksp_max_it 1

\[
\begin{pmatrix}
A & B \\
\hat{S}
\end{pmatrix}
\]
Solver Configuration: No New Code

ex62: $P_2/P_1$ Stokes Problem on Unstructured Mesh

Full Schur Complement, Schur, 1905.

-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type schur -pc_fieldsplit_schur_factorization_type full
-fieldsplit_velocity_ksp_type preonly -fieldsplit_velocity_pc_type lu -fieldsplit_pressure_ksp_rtol 1e-10 -fieldsplit_pressure_pc_type jacobi

$$\begin{pmatrix}
I & 0 \\
B^T A^{-1} & I
\end{pmatrix}
\begin{pmatrix}
A & 0 \\
0 & S
\end{pmatrix}
\begin{pmatrix}
I & A^{-1}B \\
0 & I
\end{pmatrix}$$
**Solver Configuration: No New Code**

**ex62:** $P_2/P_1$ Stokes Problem on Unstructured Mesh


```
-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type schur
-pc_fieldsplit_schur_factorization_type full
-fieldsplit_velocity_ksp_type preonly -fieldsplit_velocity_pc_type lu
-fieldsplit_pressure_ksp_rtol 1e-10 -fieldsplit_pressure_pc_type jacobi
-fieldsplit_pressure_inner_ksp_type preonly
-fieldsplit_pressure_inner_pc_type jacobi
-fieldsplit_pressure_upper_ksp_type preonly
-fieldsplit_pressure_upper_pc_type jacobi
```

\[
\begin{pmatrix}
I & 0 \\
B^T A^{-1} & I
\end{pmatrix}
\begin{pmatrix}
A & 0 \\
0 & B^T D_A^{-1} B
\end{pmatrix}
\begin{pmatrix}
I & D_A^{-1} B \\
0 & I
\end{pmatrix}
\]
ex62: $P_2/P_1$ Stokes Problem on Unstructured Mesh


-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type schur
-pc_fieldsplit_schur_factorization_type full
-pc_fieldsplit_schur_precondition self
-fieldsplit_velocity_ksp_type gmres -fieldsplit_velocity_pc_type lu
-fieldsplit_pressure_ksp_rtol 1e-5 -fieldsplit_pressure_pc_type lsc

\[
\begin{pmatrix}
I & 0 \\
B^T A^{-1} & I
\end{pmatrix}
\begin{pmatrix}
A & 0 \\
0 & \hat{S}_{LSC}
\end{pmatrix}
\begin{pmatrix}
I & A^{-1}B \\
0 & I
\end{pmatrix}
\]
Solver Configuration: No New Code

ex31: $P_2/P_1$ Stokes Problem with Temperature on Unstructured Mesh


-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type additive
-pc_fieldsplit_0_fields 0,1 -pc_fieldsplit_1_fields 2
-fieldsplit_0_ksp_type fgmres -fieldsplit_0_pc_type fieldsplit
-fieldsplit_0_pc_fieldsplit_type schur
-fieldsplit_0_pc_fieldsplit_schur_factorization_type full
-fieldsplit_0_fieldsplit_velocity_ksp_type preonly
-fieldsplit_0_fieldsplit_velocity_pc_type lu
-fieldsplit_0_fieldsplit_pressure_ksp_rtol 1e-10
-fieldsplit_0_fieldsplit_pressure_pc_type jacobi
-fieldsplit_temperature_ksp_type preonly
-fieldsplit_temperature_pc_type lu

\[
\begin{pmatrix}
  I & 0 \\
  B^T A^{-1} & I \\
\end{pmatrix}
\begin{pmatrix}
  \hat{A} & 0 \\
  0 & \hat{S} \\
\end{pmatrix}
\begin{pmatrix}
  I & A^{-1} B \\
  0 & I \\
\end{pmatrix}
\begin{pmatrix}
  0 \\
  L_T \\
\end{pmatrix}
\]
**Solver Configuration: No New Code**

**ex31**: $P_2/P_1$ Stokes Problem with Temperature on Unstructured Mesh

Upper Schur Comp. + Full Schur Comp. + Least-Squares Comm.

```bash
-ksp_type fgmres -pc_type fieldsplit -pc_fieldsplit_type schur
-pc_fieldsplit_0_fields 0,1 -pc_fieldsplit_1_fields 2
-pc_fieldsplit_schur_factorization_type upper
-fieldsplit_0_ksp_type fgmres -fieldsplit_0_pc_type fieldsplit
-fieldsplit_0_pc_fieldsplit_type schur
-fieldsplit_0_pc_fieldsplit_schur_factorization_type full
-fieldsplit_0_fieldsplit_velocity_ksp_type preonly
-fieldsplit_0_fieldsplit_velocity_pc_type lu
-fieldsplit_0_fieldsplit_pressure_ksp_rtol 1e-10
-fieldsplit_0_fieldsplit_pressure_pc_type jacobi
-fieldsplit_temperature_ksp_type gmres
-fieldsplit_temperature_pc_type lsc

\[
\begin{pmatrix}
I & 0 \\
B^T A^{-1} & I
\end{pmatrix}
\begin{pmatrix}
\hat{A} & 0 \\
0 & \hat{S}
\end{pmatrix}
\begin{pmatrix}
I & A^{-1}B \\
0 & I
\end{pmatrix}
G
\begin{pmatrix}
\hat{S}_{LSC}
\end{pmatrix}
\]
```

Matt (ANL)
Jacobi

```
ex62
  -run_type full -bc_type dirichlet -show_solution 0
  -refinement_limit 0.00625 -interpolate 1
  -vel_petsc-space_order 2 -pres_petsc-space_order 1
  -snes_monitor_short -snes_converged_reason
    -snes_view
  -ksp_gmres_restart 100 -ksp_rtol 1.0e-9
    -ksp_monitor_short
  -pc_type jacobi
```
Block diagonal

ex62

-run_type full -bc_type dirichlet -show_solution 0
-refinement_limit 0.00625 -interpolate 1
-vel_petscspace_order 2 -pres_petscspace_order 1
-snes_monitor_short -snes_converged_reason
   -snes_view
-ksp_type fgmres -ksp_gmres_restart 100
   -ksp_rtol 1.0e-9 -ksp_monitor_short
-pc_type fieldsplit -pc_fieldsplit_type additive
-fieldsplit_velocity_pc_type lu
-fieldsplit_pressure_pc_type jacobi
Block triangular

ex62

- run_type full -bc_type dirichlet -show_solution 0
- refinement_limit 0.00625 -interpolate 1
- vel_petscspace_order 2 - pres_petscspace_order 1
- snes_monitor_short - snes_converged_reason
  - snes_view
- ksp_type fgmres - ksp_gmres_restart 100
  - ksp_rtol 1.0e-9 - ksp_monitor_short
- pc_type fieldsplit - pc_fieldsplit_type multiplicative
- fieldsplit_velocity_pc_type lu
- fieldsplit_pressure_pc_type jacobi
Diagonal Schur complement

```
ex62
  -run_type full -bc_type dirichlet -show_solution 0
  -refinement_limit 0.00625 -interpolate 1
  -vel_petscspace_order 2 -pres_petscspace_order 1
  -snes_monitor_short -snes_converged_reason
    -snes_view
  -ksp_type fgmres -ksp_gmres_restart 100
    -ksp_rtol 1.0e-9 -ksp_monitor_short
  -pc_type fieldsplit -pc_fieldsplit_type schur
    -pc_fieldsplit_schur_factorization_type diag
  -fieldsplit_velocity_ksp_type gmres
    -fieldsplit_velocity_pc_type lu
  -fieldsplit_pressure_ksp_rtol 1e-10
    -fieldsplit_pressure_pc_type jacobi
```
Upper triangular Schur complement

ex62

- run_type full -bc_type dirichlet -show_solution 0
- refinement_limit 0.00625 -interpolate 1
- vel_petscspace_order 2 -pres_petscspace_order 1
- snes_monitor_short -snes_converged_reason
  - snes_view
- ksp_type fgmres -ksp_gmres_restart 100
  - ksp_rtol 1.0e-9 -ksp_monitor_short
- pc_type fieldsplit -pc_fieldsplit_type schur
  - pc_fieldsplit_schur_factorization_type upper
- fieldsplit_velocity_ksp_type gmres
  - fieldsplit_velocity_pc_type lu
- fieldsplit_pressure_ksp_rtol 1e-10
  - fieldsplit_pressure_pc_type jacobi
Lower triangular Schur complement

```bash
ex62
  -run_type full -bc_type dirichlet -show_solution 0
  -refinement_limit 0.00625 -interpolate 1
  -vel_petscspace_order 2 -pres_petscspace_order 1
  -snes_monitor_short -snes_converged_reason
    -snes_view
  -ksp_type fgmres -ksp_gmres_restart 100
    -ksp_rtol 1.0e-9 -ksp_monitor_short
  -pc_type fieldsplit -pc_fieldsplit_type schur
    -pc_fieldsplit_schur_factorization_type lower
  -fieldsplit_velocity_ksp_type gmres
    -fieldsplit_velocity_pc_type lu
  -fieldsplit_pressure_ksp_rtol 1e-10
    -fieldsplit_pressure_pc_type jacobi
```
Full Schur complement

`ex62`

```bash
-run_type full -bc_type dirichlet -show_solution 0
-refinement_limit 0.00625 -interpolate 1
-vel_petscspace_order 2 -pres_petscspace_order 1
-snes_monitor_short -snes_converged_reason
    -snes_view
-ksp_type fgmres -ksp_gmres_restart 100
    -ksp_rtol 1.0e-9 -ksp_monitor_short
-pc_type fieldsplit -pc_fieldsplit_type schur
    -pc_fieldsplit_schur_factorization_type full
-fieldsplit_velocity_ksp_type gmres
    -fieldsplit_velocity_pc_type lu
-fieldsplit_pressure_ksp_rtol 1e-10
    -fieldsplit_pressure_pc_type jacobi
```
ex55: Allen-Cahn problem in 2D

- constant mobility
- triangular elements

Geometric multigrid method for saddle point variational inequalities:

```
./ex55 -ksp_type fgmres -pc_type mg -mg_levels_ksp_type fgmres
    -mg_levels_pc_type fieldsplit -mg_levels_pc_fieldsplit_detect_saddle_point
    -mg_levels_pc_fieldsplit_type schur -da_grid_x 65 -da_grid_y 65
    -mg_levels_pc_fieldsplit_factorization_type full
    -mg_levels_pc_fieldsplit_schur_precondition user
    -mg_levels_fieldsplit_1_ksp_type gmres -mg_coarse_ksp_type preonly
    -mg_levels_fieldsplit_1_pc_type none -mg_coarse_pc_type svd
    -mg_levels_fieldsplit_0_ksp_type preonly
    -mg_levels_fieldsplit_0_pc_type sor -pc_mg_levels 5
    -mg_levels_fieldsplit_0_pc_sor_forward -pc_mg_galerkin
    -snes_vi_monitor -ksp_monitor_true_residual -snes_atol 1.e-11
    -mg_levels_ksp_monitor -mg_levels_fieldsplit_ksp_monitor
    -mg_levels_ksp_max_it 2 -mg_levels_fieldsplit_ksp_max_it 5
```
ex55: Allen-Cahn problem in 2D

Run flexible GMRES with 5 levels of multigrid as the preconditioner

```
./ex55 -ksp_type fgmres -pc_type mg -pc_mg_levels 5
 -da_grid_x 65 -da_grid_y 65
```

Use the Galerkin process to compute the coarse grid operators

```
-pc_mg_galerkin
```

Use SVD as the coarse grid saddle point solver

```
-mg_coarse_ksp_type preonly -mg_coarse_pc_type svd
```
ex55: Allen-Cahn problem in 2D

Run flexible GMRES with 5 levels of multigrid as the preconditioner

```
./ex55 -ksp_type fgmres -pc_type mg -pc_mg_levels 5
   -da_grid_x 65 -da_grid_y 65
```

Use the Galerkin process to compute the coarse grid operators

```
-pc_mg_galerkin
```

Use SVD as the coarse grid saddle point solver

```
-mg_coarse_ksp_type preonly -mg_coarse_pc_type svd
```
**ex55**: Allen-Cahn problem in 2D

Run flexible GMRES with 5 levels of multigrid as the preconditioner

```
./ex55 -ksp_type fgmres -pc_type mg -pc_mg_levels 5 -da_grid_x 65 -da_grid_y 65
```

Use the Galerkin process to compute the coarse grid operators

```
-pc_mg_galerkin
```

Use SVD as the coarse grid saddle point solver

```
-mg_coarse_ksp_type preonly -mg_coarse_pc_type svd
```
**ex55: Allen-Cahn problem in 2D**

Run flexible GMRES with 5 levels of multigrid as the preconditioner

```
./ex55 -ksp_type fgmres -pc_type mg -pc_mg_levels 5 -da_grid_x 65 -da_grid_y 65
```

Use the Galerkin process to compute the coarse grid operators

```
-pc_mg_galerkin
```

Use SVD as the coarse grid saddle point solver

```
-mg_coarse_ksp_type preonly -mg_coarse_pc_type svd
```
Programming with Options

**ex55: Allen-Cahn problem in 2D**

Smoother: Flexible GMRES (2 iterates) with a Schur complement PC

- `mg_levels_ksp_type fgmres`  
- `mg_levels_pc_fieldsplit_detect_saddle_point`  
- `mg_levels_ksp_max_it 2`  
- `mg_levels_pc_type fieldsplit`  
- `mg_levels_pc_fieldsplit_type schur`  
- `mg_levels_pc_fieldsplit_factorization_type full`  
- `mg_levels_pc_fieldsplit_schur_precondition diag`

Schur complement solver: GMRES (5 iterates) with no preconditioner

- `mg_levels_fieldsplit_1_ksp_type gmres`  
- `mg_levels_fieldsplit_1_pc_type none`  
- `mg_levels_fieldsplit_ksp_max_it 5`

Schur complement action: Use only the lower diagonal part of A00

- `mg_levels_fieldsplit_0_ksp_type preonly`  
- `mg_levels_fieldsplit_0_pc_type sor`  
- `mg_levels_fieldsplit_0_pc_sor_forward`
**Programming with Options**

**ex55**: Allen-Cahn problem in 2D

**Smoother**: Flexible GMRES (2 iterates) with a Schur complement PC

```
-mg_levels_ksp_type fgmres -mg_levels_pc_fieldsplit_detect_saddle_point
-mg_levels_ksp_max_it 2 -mg_levels_pc_type fieldsplit
-mg_levels_pc_fieldsplit_type schur
-mg_levels_pc_fieldsplit_factorization_type full
-mg_levels_pc_fieldsplit_schur_precondition diag
```

**Schur complement solver**: GMRES (5 iterates) with no preconditioner

```
-mg_levels_fieldsplit_1_ksp_type gmres
-mg_levels_fieldsplit_1_pc_type none -mg_levels_fieldsplit_ksp_max_it 5
```

**Schur complement action**: Use only the lower diagonal part of A00

```
-mg_levels_fieldsplit_0_ksp_type preonly
-mg_levels_fieldsplit_0_pc_type sor
-mg_levels_fieldsplit_0_pc_sor_forward
```
ex55: Allen-Cahn problem in 2D

Smoother: Flexible GMRES (2 iterates) with a Schur complement PC

- `mg_levels_ksp_type fgmres`  
- `mg_levels_pc_fieldsplit_detect_saddle_point`  
- `mg_levels_ksp_max_it 2`  
- `mg_levels_pc_type fieldsplit`  
- `mg_levels_pc_fieldsplit_type schur`  
- `mg_levels_pc_fieldsplit_factorization_type full`  
- `mg_levels_pc_fieldsplit_schur_precondition diag`

Schur complement solver: GMRES (5 iterates) with no preconditioner

- `mg_levels_fieldsplit_1_ksp_type gmres`  
- `mg_levels_fieldsplit_1_pc_type none`  
- `mg_levels_fieldsplit_ksp_max_it 5`

Schur complement action: Use only the lower diagonal part of $A_{00}$

- `mg_levels_fieldsplit_0_ksp_type preonly`  
- `mg_levels_fieldsplit_0_pc_type sor`  
- `mg_levels_fieldsplit_0_pc_sor_forward`
Programming with Options

**ex55**: Allen-Cahn problem in 2D

Smoother: Flexible GMRES (2 iterates) with a Schur complement PC

```
-mg_levels_ksp_type fgmres -mg_levels_pc_fieldsplit_detect_saddle_point
-mg_levels_ksp_max_it 2 -mg_levels_pc_type fieldsplit
-mg_levels_pc_fieldsplit_type schur
-mg_levels_pc_fieldsplit_factorization_type full
-mg_levels_pc_fieldsplit_schur_precondition diag
```

Schur complement solver: GMRES (5 iterates) with no preconditioner

```
-mg_levels_fieldsplit_1_ksp_type gmres
-mg_levels_fieldsplit_1_pc_type none -mg_levels_fieldsplit_ksp_max_it 5
```

Schur complement action: Use only the lower diagonal part of A00

```
-mg_levels_fieldsplit_0_ksp_type preonly
-mg_levels_fieldsplit_0_pc_type sor
-mg_levels_fieldsplit_0_pc_sor_forward
```
Null spaces

For a single matrix, use

\texttt{MatSetNullSpace(J, nullSpace);}

to alter the \texttt{KSP}, and

\texttt{MatSetNearNullSpace(J, nearNullSpace);}

to set the coarse basis for AMG.

But this will not work for dynamically created operators.
Null spaces

For a single matrix, use

\texttt{MatSetNullSpace(J, nullSpace);}

to alter the \texttt{KSP}, and

\texttt{MatSetNearNullSpace(J, nearNullSpace);} 

to set the coarse basis for AMG.

But this will not work for dynamically created operators.
Can attach a nullspace to the **IS** that creates a split,

\[
\text{PetscObjectCompose}(\text{pressure IS}, "nullspace",
\quad (\text{PetscObject}) \text{nullSpacePres});
\]

If the **DM** makes the **IS**, use

\[
\text{PetscObject} \quad \text{pressure};
\]

\[
\text{DMGetField}(\text{dm}, 1, &\text{pressure});
\quad \text{PetscObjectCompose}(\text{pressure}, "nullspace",
\quad (\text{PetscObject}) \text{nullSpacePres});
\]
Advanced Solvers
- Fieldsplit
- Multigrid
- Nonlinear Preconditioning
Why not use AMG?

- Of course we will try AMG
  - GAMG, `pc_type gamg`
  - ML, `download-ml, pc_type ml`
  - BoomerAMG, `download-hypre, pc_type hypre`
    - `pc_hypre_type boomeramg`

- Problems with
  - vector character
  - anisotropy
  - scalability of setup time
Why not use AMG?

- Of course we will try AMG
  - GAMG, \texttt{-pc\_type gamg}
  - ML, \texttt{-download-ml, -pc\_type ml}
  - BoomerAMG, \texttt{-download-hypre, -pc\_type hypre}
    - \texttt{-pc\_hypre\_type boomeramg}

- Problems with
  - vector character
  - anisotropy
  - scalability of setup time
Why not use AMG?

- Of course we will try AMG
  - GAMG, -pc_type gamg
  - ML, -download-ml, -pc_type ml
  - BoomerAMG, -download-hypre, -pc_type hypre, -pc_hypre_type boomeramg

- Problems with
  - vector character
  - anisotropy
  - scalability of setup time
Multigrid with DM

Allows multigrid with some simple command line options

- \texttt{-pc\_type mg}, \texttt{-pc\_mg\_levels}
- \texttt{-pc\_mg\_type}, \texttt{-pc\_mg\_cycle\_type}, \texttt{-pc\_mg\_galerkin}
- \texttt{-mg\_levels\_1\_ksp\_type}, \texttt{-mg\_levels\_1\_pc\_type}
- \texttt{-mg\_coarse\_ksp\_type}, \texttt{-mg\_coarse\_pc\_type}
- \texttt{-da\_refine}, \texttt{-ksp\_view}

Interface also works with GAMG and 3rd party packages like ML
## A 2D Problem

Problem has:
- 1,640,961 unknowns (on the fine level)
- 8,199,681 nonzeros

### Options

<table>
<thead>
<tr>
<th>Option</th>
<th>Explanation</th>
</tr>
</thead>
<tbody>
<tr>
<td>./ex5 -da_grid_x 21 -da_grid_y 21</td>
<td>Original grid is 21x21</td>
</tr>
<tr>
<td>-ksp_rtol 1.0e-9</td>
<td>Solver tolerance</td>
</tr>
<tr>
<td>-da_refine 6</td>
<td>6 levels of refinement</td>
</tr>
<tr>
<td>-pc_type mg</td>
<td>4 levels of multigrid</td>
</tr>
<tr>
<td>-pc_mg_levels 4</td>
<td>Describe solver</td>
</tr>
<tr>
<td>-snes_monitor -snes_view</td>
<td></td>
</tr>
</tbody>
</table>
A 3D Problem

Problem has:
- 1,689,600 unknowns (on the fine level)
- 89,395,200 nonzeros

Options

<table>
<thead>
<tr>
<th>Options</th>
<th>Explanation</th>
</tr>
</thead>
<tbody>
<tr>
<td>./ex48 -M 5 -N 5</td>
<td>Coarse problem size</td>
</tr>
<tr>
<td>-da_refine 5</td>
<td>5 levels of refinement</td>
</tr>
<tr>
<td>-ksp_rtol 1.0e-9</td>
<td>Solver tolerance</td>
</tr>
<tr>
<td>-thi_mat_type baij</td>
<td>Needs SOR</td>
</tr>
<tr>
<td>-pc_type mg</td>
<td>4 levels of multigrid</td>
</tr>
<tr>
<td>-pc_mg_levels 4</td>
<td></td>
</tr>
<tr>
<td>-snes_monitor -snes_view</td>
<td>Describe solver</td>
</tr>
</tbody>
</table>
The Full Multigrid algorithm (FMG)

- V-cycle at each level,
- then interpolate to the next finer grid
- Can solve to discretization error with a single iteration
$C_{FMG} = \left(1 + \frac{1}{2^d} + \frac{1}{2^{2d}} + \ldots\right) C_V$

$= \sum_{n=0}^{\infty} \frac{1}{2^{nd}} C_V$

$= \frac{2^d}{2^d - 1} C_V$

$= \left(\frac{2^d}{2^d - 1}\right)^2 C_{\text{twolevel}}$.

1D FMG is $2 \times C_V$, 3D FMG is $\frac{8}{7} \times C_V$. 

Matt (ANL)
Full Multigrid Work

\[ C_{FMG} = \left( 1 + \frac{1}{2^d} + \frac{1}{2^{2d}} + \ldots \right) C_V \]
\[ = \sum_{n=0}^{\infty} \frac{1}{2^{nd}} C_V \]
\[ = \frac{2^d}{2^d - 1} C_V \]
\[ = \left( \frac{2^d}{2^d - 1} \right)^2 C_{\text{twolevel}}. \]

1D FMG is \( 2 \times C_V \), 3D FMG is \( \frac{8}{7} \times C_V \)
Suppose we have an order $\alpha$ method,

$$\|x - x_h\| < Ch^\alpha$$

FD and $P_1$ both have $\alpha = 2$
Full Multigrid Accuracy

$E_d$  Discretization Error
$E_a$  Algebraic Error

Choose iterative tolerance so that

$$E_a = rE_d \quad r < 1$$

and

$$E \leq E_d + E_a = (1 + r)Ch^\alpha$$
Suppose

- Finish V-cycle for $2h$ grid,
- Use as coarse correction for $h$ grid
- Perform final V-cycle for $h$ grid
- Need V-cycle error reduction factor $\eta$ to get $r$ reduction in $E_a$
Full Multigrid Accuracy

\[ \eta E_a < r Ch^\alpha \]
\[ \eta (E - E_d) < r Ch^\alpha \]
\[ \eta ((1 + r)C(2h)^\alpha - Ch^\alpha) < r Ch^\alpha \]
\[ \eta ((1 + r)2^\alpha - 1) < r \]
\[ \eta < \frac{1}{2^\alpha + \frac{2^\alpha - 1}{r}}. \]

If \( \alpha = 2 \) and \( r = \frac{1}{2} \), then \( \eta < \frac{1}{10} \).
Advanced Solvers

Multigrid

Full Multigrid Experiment
V-cycle

./ex5 -mms 1 -par 0.0 -da_refine 3 -snes_type newtonls -snes_max_it 1 -ksp_rtol 1e-10 -pc_type mg -snes_monitor_short -ksp_monitor_short gives

0 SNES Function norm 0.0287773
0 KSP Residual norm 0.793727
1 KSP Residual norm 0.00047526
2 KSP Residual norm 4.18007e-06
3 KSP Residual norm 1.1668e-07
4 KSP Residual norm 3.25952e-09
5 KSP Residual norm 7.274e-11
1 SNES Function norm 2.251e-10
N: 625 error l2 1.21529e-13 inf 9.53484e-12
Full Multigrid Experiment

V-cycle

./ex5 -mms 1 -par 0.0 -da_refine 3 -snes_type newtonls -snes_max_it 1 -ksp_rtol 1e-10 -pc_type mg -snes_monitor_short -ksp_monitor_short

and it changes little if we refine six more times

0 SNES Function norm 0.000455131
  0 KSP Residual norm 50.6842
  1 KSP Residual norm 0.00618427
  2 KSP Residual norm 9.87833e-07
  3 KSP Residual norm 2.99517e-09
1 SNES Function norm 2.83358e-09
N: 2362369 error l2 1.28677e-15 inf 7.68693e-12
Full Multigrid Experiment

FMG

./ex5 -mms 1 -par 0.0 -da_refine 3 -snes_type newtonls -snes_max_it 1 -ksp_rtol 1e-10 -pc_type mg -snes_monitor_short -ksp_monitor_short -pc_mg_type full

We do not seem to see the convergence acceleration

0 SNES Function norm 0.0287773
 0 KSP Residual norm 0.799687
1 KSP Residual norm 6.95292e-05
2 KSP Residual norm 1.50836e-06
3 KSP Residual norm 2.62524e-08
4 KSP Residual norm 6.184e-10
5 KSP Residual norm 1.275e-11
1 SNES Function norm 3.757e-11
N: 625 error l2 2.1428e-14 inf 1.80611e-12
Full Multigrid Experiment

FMG

./ex5 -mms 1 -par 0.0 -da_refine 3 -snes_type newtonls -snes_max_it 1 -ksp_rtol 1e-10 -pc_type mg -snes_monitor_short -ksp_monitor_short -pc_mg_type full

although its a little more apparent as we refine,

0 SNES Function norm 0.000455131
0 KSP Residual norm 51.2
1 KSP Residual norm 2.92416e-06
2 KSP Residual norm 3.76404e-09
1 SNES Function norm 8.50096e-09
N: 2362369 error l2 1.70304e-15 inf 6.22476e-11
# ! /usr/bin/env python
import argparse
import subprocess
import numpy as np

parser = argparse.ArgumentParser(
    description = 'CAAM 519 FMG',
    epilog = 'For more information, visit http://www.mcs.anl.gov/petsc',
    formatter_class = argparse.ArgumentDefaultsHelpFormatter)
parser.add_argument('--kmax', type=int, default=5,
    help='The number of doublings to test')
parser.add_argument('--save', action='store_true', default=False,
    help='Save the figures')
args = parser.parse_args()

sizesA = []
sizesB = []
errorA = []
errorB = []
```python
for k in range(args.kmax):
    options = ['-snes_type', 'newtonls', '-snes_max_it', '1', '-da_refine', '-par', '0.0', '-ksp_atol', '1e-1', '-mms', '1', '-pc_type', 'mg', '-pc_mg_type', 'multiplicative', '-mg_levels_ksp_max_it', '5']
    cmd = './ex5' + ' '.join(options)
    out = subprocess.check_output(['./ex5'] + options).split(' ')
    # This is l_2, out[6] is l_infty
    sizesA.append(int(out[1]))
    errorA.append(float(out[4]))

for k in range(args.kmax):
    options = ['-snes_type', 'newtonls', '-snes_max_it', '1', '-da_refine', '-par', '0.0', '-ksp_atol', '1e-1', '-mms', '1', '-pc_type', 'mg', '-pc_mg_type', 'full', '-mg_levels_ksp_max_it', '5']
    cmd = './ex5' + ' '.join(options)
    out = subprocess.check_output(['./ex5'] + options).split(' ')
    # This is l_2, out[6] is l_infty
    sizesB.append(int(out[1]))
    errorB.append(float(out[4]))
```
SA = np.array(sizesA)
SB = np.array(sizesB)

from pylab import legend, plot, loglog, show, title, xlabel, ylabel, savefig
loglog(SA, errorA, 'r', SA, 5e-6 * SA ** -0.5, 'r--',
       SB, errorB, 'g', SB, 5e-6 * SB ** -1., 'g--')
title('SNES ex5')
xlabel('Problem size $N$')
ylabel('Error $\|u - u^*\|$')
legend(['GMRES/GMG', '$h = N^{-1/2}$', 'GMRES/FMG', '$h^2 = N^{-1}$'],
       'upper right', shadow = True)

if args.save:
    savefig('fmg.png')
else:
    show()
Full Multigrid Experiment
Comparison

SNES ex5

- GMRES/GMG
- $h = N^{-1/2}$
- GMRES/FMG
- $h^2 = N^{-1}$

Error vs Problem size $N$

- Error $|u - u^*|$
Outline

3 Advanced Solvers
- Fieldsplit
- Multigrid
- Nonlinear Preconditioning
Driven Cavity Problem

SNES ex19.c

./ex19  -lidvelocity 100 -grashof 1e2
   -da_grid_x 16 -da_grid_y 16 -da_refine 2
   -snes_monitor_short -snes_converged_reason -snes_view

\[ -\Delta U - \partial_y \Omega = 0 \]
\[ -\Delta V + \partial_x \Omega = 0 \]
\[ -\Delta \Omega + \nabla \cdot ([U \Omega, V \Omega]) - \text{Gr} \ \partial_x T = 0 \]
\[ -\Delta T + \text{Pr} \ \nabla \cdot ([U T, V T]) = 0 \]
Driven Cavity Problem

**SNES ex19.c**

```bash
./ex19 -lidvelocity 100 -grashof 1e2
    -da_grid_x 16 -da_grid_y 16 -da_refine 2
    -snes_monitor_short -snes_converged_reason -snes_view
```

lid velocity = 100, prandtl # = 1, grashof # = 100

0 SNES Function norm 768.116
1 SNES Function norm 658.288
2 SNES Function norm 529.404
3 SNES Function norm 377.51
4 SNES Function norm 304.723
5 SNES Function norm 2.59998
6 SNES Function norm 0.00942733
7 SNES Function norm 5.20667e-08

Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE iterations 7
Driven Cavity Problem

SNES ex19.c

./ex19 -lidvelocity 100 -grashof 1e4
   -da_grid_x 16 -da_grid_y 16 -da_refine 2
   -snes_monitor_short -snes_converged_reason -snes_view
Driven Cavity Problem

SNES ex19.c

./ex19 -lidvelocity 100 -grashof 1e4
     -da_grid_x 16 -da_grid_y 16 -da_refine 2
     -snes_monitor_short -snes_converged_reason -snes_view

lid velocity = 100, prandtl # = 1, grashof # = 10000
  0 SNES Function norm 785.404
  1 SNES Function norm 663.055
  2 SNES Function norm 519.583
  3 SNES Function norm 360.87
  4 SNES Function norm 245.893
  5 SNES Function norm 1.8117
  6 SNES Function norm 0.00468828
  7 SNES Function norm 4.417e-08
Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE iterations 7
Driven Cavity Problem

SNES ex19.c

./ex19 -lidvelocity 100 -grashof 1e5
-\texttt{da\_grid\_x} 16 -\texttt{da\_grid\_y} 16 -\texttt{da\_refine} 2
-\texttt{snes\_monitor\_short} -\texttt{snes\_converged\_reason} -\texttt{snes\_view}
Driven Cavity Problem

SNES ex19.c

./ex19 -lidvelocity 100 -grashof 1e5
   -da_grid_x 16 -da_grid_y 16 -da_refine 2
   -snes_monitor_short -snes_converged_reason -snes_view

lid velocity = 100, prandtl # = 1, grashof # = 100000
0 SNES Function norm 1809.96
Nonlinear solve did not converge due to DIVERGED_LINEAR_SOLVE iterations 0
Driven Cavity Problem

SNES ex19.c

./ex19 -lidvelocity 100 -grashof 1e5
   -da_grid_x 16 -da_grid_y 16 -da_refine 2 -pc_type lu
   -snes_monitor_short -snes_converged_reason -snes_view

lid velocity = 100, prandtl # = 1, grashof # = 100000
0  SNES Function norm 1809.96
1  SNES Function norm 1678.37
2  SNES Function norm 1643.76
3  SNES Function norm 1559.34
4  SNES Function norm 1557.6
5  SNES Function norm 1510.71
6  SNES Function norm 1500.47
7  SNES Function norm 1498.93
8  SNES Function norm 1498.44
9  SNES Function norm 1498.27
10 SNES Function norm 1498.18
11 SNES Function norm 1498.12
12 SNES Function norm 1498.11
13 SNES Function norm 1498.11
14 SNES Function norm 1498.11
...
Deceleration of Convergence

```
./ex19 -lidvelocity 100 -grashof 1.3372e2
    -da_grid_x 16 -da_grid_y 16 -da_refine 2
    -snes_type nrichardson -snes_linesearch_type cp -snes_max_it 10000
```
Stagnation of Newton

./ex19 -lidvelocity 100 -grashof 1.3372e4
-da_grid_x 16 -da_grid_y 16 -da_refine 2
-snes_type newtonls -snes_max_it 100 -pc_type lu
Stagnation of Newton

./ex19 -lidity 100 -grashof 1.3373e4
-da_grid_x 16 -da_grid_y 16 -da_refine 2
-snes_type newtonls -snes_max_it 100 -pc_type lu

![Newton Convergence for SNES ex19](image)
Preconditioning NRichardson with Newton

```
./ex19 -lidvelocity 100 -grashof 1.3373e2
-da_grid_x 16 -da_grid_y 16 -da_refine 2
-snes_type nrichardson -snes_max_it 200
-npc_snes_type newtonls -npc_snes_max_it 3 -npc_pc_type lu
```
Preconditioning NRichardson with Newton

./ex19 -lidvelocity 100 -grashof 1.3373e2
   -da_grid_x 16 -da_grid_y 16 -da_refine 2
   -snes_type nrichardson -snes_max_it 200
   -npc_snes_type newtonls -npc_snes_max_it 4 -npc_pc_type lu

NRichardson — Newton Convergence for SNES ex19

- 4 inner Newton its
- 3 inner Newton its
Preconditioning Newton with NRichardson

./ex19 -lidvelo city 100 -grashof 1.3373e2
   -da_grid_x 16 -da_grid_y 16 -da_refine 2
   -snes_type newtonls -snes_max_it 1000 -pc_type lu
   -npc_snes_type nrichardson -npc_snes_max_it 1
Preconditioning Newton with NRichardson

./ex19 -lidvelocity 100 -grashof 1.3373e2 -da_grid_x 16 -da_grid_y 16 -da_refine 2 -snes_type newtonls -snes_max_it 1000 -pc_type lu -npc_snes_type nrichardson -npc_snes_max_it 3
Preconditioning Newton with NRichardson

./ex19 -lidvelocity 100 -grashof 1.3373e2
   -da_grid_x 16 -da_grid_y 16 -da_refine 2
   -snes_type newtonls -snes_max_it 1000 -pc_type lu
   -npc_snes_type nrichardson -npc_snes_max_it 5
Preconditioning Newton with NRichardson

./ex19 -lidvelocity 100 -grashof 1.3373e2
-da_grid_x 16 -da_grid_y 16 -da_refine 2
-snes_type newtonls -snes_max_it 1000 -pc_type lu
-npc_snes_type nrichardson -npc_snes_max_it 6

![Newton - NRichardson Convergence for SNES ex19](image)
Nonlinear Preconditioning

./ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short -snes_type newtonls -snes_converged_reason -pc_type lu

lid velocity = 100, prandtl # = 1, grashof # = 50000

0 SNES Function norm 1228.95
1 SNES Function norm 1132.29
2 SNES Function norm 1026.17
3 SNES Function norm 925.717
4 SNES Function norm 924.778
5 SNES Function norm 836.867

21 SNES Function norm 585.143
22 SNES Function norm 585.142
23 SNES Function norm 585.142
24 SNES Function norm 585.142

...
Nonlinear Preconditioning

./ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short
   -snes_type fas -snes_converged_reason
   -fas_levels_snes_type gs -fas_levels_snes_max_it 6

lid velocity = 100, prandtl # = 1, grashof # = 50000
0 SNES Function norm 1228.95
1 SNES Function norm 574.793
2 SNES Function norm 513.02
3 SNES Function norm 216.721
4 SNES Function norm 85.949
Nonlinear solve did not converge due to DIVERGED_INNER iterations 4
Advanced Solvers  Nonlinear Preconditioning

Nonlinear Preconditioning

./ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short
-snes_type fas -snes_converged_reason
-fas_levels_snes_type gs -fas_levels_snes_max_it 6
-fas_coarse_snes_converged_reason

lid velocity = 100, prandtl # = 1, grashof # = 50000
  0 SNES Function norm 1228.95
    Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 12
  1 SNES Function norm 574.793
    Nonlinear solve did not converge due to DIVERGED_MAX_IT its 50
  2 SNES Function norm 513.02
    Nonlinear solve did not converge due to DIVERGED_MAX_IT its 50
  3 SNES Function norm 216.721
    Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 22
  4 SNES Function norm 85.949
    Nonlinear solve did not converge due to DIVERGED_LINE_SEARCH its 42
Nonlinear solve did not converge due to DIVERGED_INNER iterations 4
Nonlinear Preconditioning

./ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short
-snes_type fas -snes_converged_reason
-fas_levels_snes_type gs -fas_levels_snes_max_it 6
-fas_coarse_snes_linesearch_type basic
-fas_coarse_snes_converged_reason

lid velocity = 100, prandtl # = 1, grashof # = 50000
0 SNES Function norm 1228.95
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 6
47 SNES Function norm 78.8401
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 5
48 SNES Function norm 73.1185
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 6
49 SNES Function norm 78.834
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 5
50 SNES Function norm 73.1176
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 6

...
./ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short -snes_type nrichardson -npc_snes_max_it 1 -snes_converged_reason -npc_snes_type fas -npc_fas_coarse_snes_converged_reason -npc_fas_levels_snes_type gs -npc_fas_levels_snes_max_it 6 -npc_fas_coarse_snes_linesearch_type basic

lid velocity = 100, prandtl # = 1, grashof # = 50000
0 SNES Function norm 1228.95
   Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 6
1 SNES Function norm 552.271
   Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 27
2 SNES Function norm 173.45
   Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 45
   ...
43 SNES Function norm 3.45407e-05
   Nonlinear solve converged due to CONVERGED_SNORM_RELATIVE its 2
44 SNES Function norm 1.6141e-05
   Nonlinear solve converged due to CONVERGED_SNORM_RELATIVE its 2
45 SNES Function norm 9.13386e-06
   Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE iterations 45

Matt (ANL)
Nonlinear Preconditioning

./ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short -snes_type ngmres -npc_snes_max_it 1 -snes_converged_reason -npc_snes_type fas -npc_fas_coarse_snes_converged_reason -npc_fas_levels_snes_type gs -npc_fas_levels_snes_max_it 6 -npc_fas_coarse_snes_linesearch_type basic

lid velocity = 100, prandtl # = 1, grashof # = 50000
0 SNES Function norm 1228.95
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 6
1 SNES Function norm 538.605
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 13
2 SNES Function norm 178.005
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 24
...
27 SNES Function norm 0.000102487
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE its 2
28 SNES Function norm 4.2744e-05
  Nonlinear solve converged due to CONVERGED_SNORM_RELATIVE its 2
29 SNES Function norm 1.01621e-05
  Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE iterations 29
./ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short -snes_type ngmres -npc_snes_max_it 1 -snes_converged_reason -npc_snes_type fas -npc_fas_coarse_snes_converged_reason -npc_fas_levels_snes_type newtonls -npc_fas_levels_snes_max_it 6 -npc_fas_levels_snes_linesearch_type basic -npc_fas_levels_snes_max_linear_solve_fail 30 -npc_fas_levels_ksp_max_it 20 -npc_fas_levels_snes_converged_reason -npc_fas_coarse_snes_linesearch_type basic

lid velocity = 100, prandtl # = 1, grashof # = 50000

0 SNES Function norm 1228.95
  Nonlinear solve did not converge due to DIVERGED_MAX_IT its 6

  Nonlinear solve converged due to CONVERGED_SNORM_RELATIVE its 1

  1 SNES Function norm 0.1935
  2 SNES Function norm 0.0179938
  3 SNES Function norm 0.00223698
  4 SNES Function norm 0.000190461
  5 SNES Function norm 1.6946e-06

Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE iterations 5
Nonlinear Preconditioning

/ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short
  -snes_type composite -snes_composite_type additiveoptimal
  -snes_composite_sneses fas,newtonls -snes_converged_reason
  -sub_0_fas_levels_snes_type gs -sub_0_fas_levels_snes_max_it 6
    -sub_0_fas_coarse_snes_linesearch_type basic
  -sub_1_snes_linesearch_type basic -sub_1_pc_type mg

lid velocity = 100, prandtl # = 1, grashof # = 50000
  0 SNES Function norm 1228.95
  1 SNES Function norm 541.462
  2 SNES Function norm 162.92
  3 SNES Function norm 48.8138
  4 SNES Function norm 11.1822
  5 SNES Function norm 0.181469
  6 SNES Function norm 0.00170909
  7 SNES Function norm 3.24991e-08
Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE iterations 7
Nonlinear Preconditioning

 ./ex19 -lidvelocity 100 -grashof 5e4 -da_refine 4 -snes_monitor_short -snes_type composite -snesCompositeType multiplicative -snesCompositeSneses fas,newtonls -snes_converged_reason -sub_0_fas_levels_snes_type gs -sub_0_fas_levels_snes_max_it 6 -sub_0_fas_coarse_snes_linesearch_type basic -sub_1_snes_linesearch_type basic -sub_1_pc_type mg

lid velocity = 100, prandtl # = 1, grashof # = 50000
0 SNES Function norm 1228.95
1 SNES Function norm 544.404
2 SNES Function norm 18.2513
3 SNES Function norm 0.488689
4 SNES Function norm 0.000108712
5 SNES Function norm 5.68497e-08
Nonlinear solve converged due to CONVERGED_FNORM_RELATIVE iterations 5
## Nonlinear Preconditioning

<table>
<thead>
<tr>
<th>Solver</th>
<th>T</th>
<th>N. It</th>
<th>L. It</th>
<th>Func</th>
<th>Jac</th>
<th>PC</th>
<th>NPC</th>
</tr>
</thead>
<tbody>
<tr>
<td>$(\mathcal{K} - \text{MG})$</td>
<td>9.83</td>
<td>17</td>
<td>352</td>
<td>34</td>
<td>85</td>
<td>370</td>
<td>–</td>
</tr>
<tr>
<td>NGMRES $- R$</td>
<td>7.48</td>
<td>10</td>
<td>220</td>
<td>21</td>
<td>50</td>
<td>231</td>
<td>10</td>
</tr>
<tr>
<td>$(\mathcal{K} - \text{MG})$</td>
<td>6.23</td>
<td>162</td>
<td>0</td>
<td>2382</td>
<td>377</td>
<td>754</td>
<td>–</td>
</tr>
<tr>
<td>FAS</td>
<td>8.07</td>
<td>10</td>
<td>197</td>
<td>232</td>
<td>90</td>
<td>288</td>
<td>–</td>
</tr>
<tr>
<td>FAS + $(\mathcal{K} - \text{MG})$</td>
<td>4.01</td>
<td>5</td>
<td>80</td>
<td>103</td>
<td>45</td>
<td>125</td>
<td>–</td>
</tr>
<tr>
<td>FAS * $(\mathcal{K} - \text{MG})$</td>
<td>3.20</td>
<td>50</td>
<td>0</td>
<td>1180</td>
<td>192</td>
<td>384</td>
<td>50</td>
</tr>
<tr>
<td>NRICH $- L$ FAS</td>
<td>1.91</td>
<td>24</td>
<td>0</td>
<td>447</td>
<td>83</td>
<td>166</td>
<td>24</td>
</tr>
<tr>
<td>NGMRES $- R$ FAS</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
</tbody>
</table>
Nonlinear Preconditioning

See discussion in:


Outline

1. Getting Started with PETSc
2. PETSc Integration
3. Advanced Solvers
4. More Stuff
Communication abstractions
- PetscSF and VecScatter

Meshing abstractions
- DMDA, DMStag, DMPlex, DMForest

Mesh adaptivity
- Interfaces to p4est and Pragmatic

Finite elements and finite volumes
- PetscFE and PetscFV