

STRUCTURE OF GOMA

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DIRECTORY LAYOUT

```
| -- cmake                # CMake modules
| -- CMakeLists.txt       # CMake configuration
| -- docs                 # docs, has some undocumented things
| -- include              # Include files in here
|   | -- ac_conti.h
|   | -- ...
|   | -- brkfix
|   |   | -- bbb.h
|   |   | -- ...
| -- LICENSE              # Open source license
| -- scripts              # build scripts, some postproc utilities
| -- src                  # Source files (c/cpp/fortran files)
|   | -- ac_conti.c
|   | -- ...
|   | -- brkfix
|   |   | -- bbb.c
|   |   | -- ...
```

FILE NAMING CONVENTION

- **ac_** Augmenting condition
- **bc_** Boundary condition
- **dp_** Distributed Processing
- **el_** element (num nodes, gauss points, etc...)
- **loca_** builtin loca library
- **mm_** Moving mesh (more of a catch all)
- **rd_/wr_** I/O read/write
- **rf_** Reacting flow (historical name from `rf_salssa`, which Goma was forked)
- **sl_** Interfaces to solvers like Trilinos
- **user_** user routines, less used now

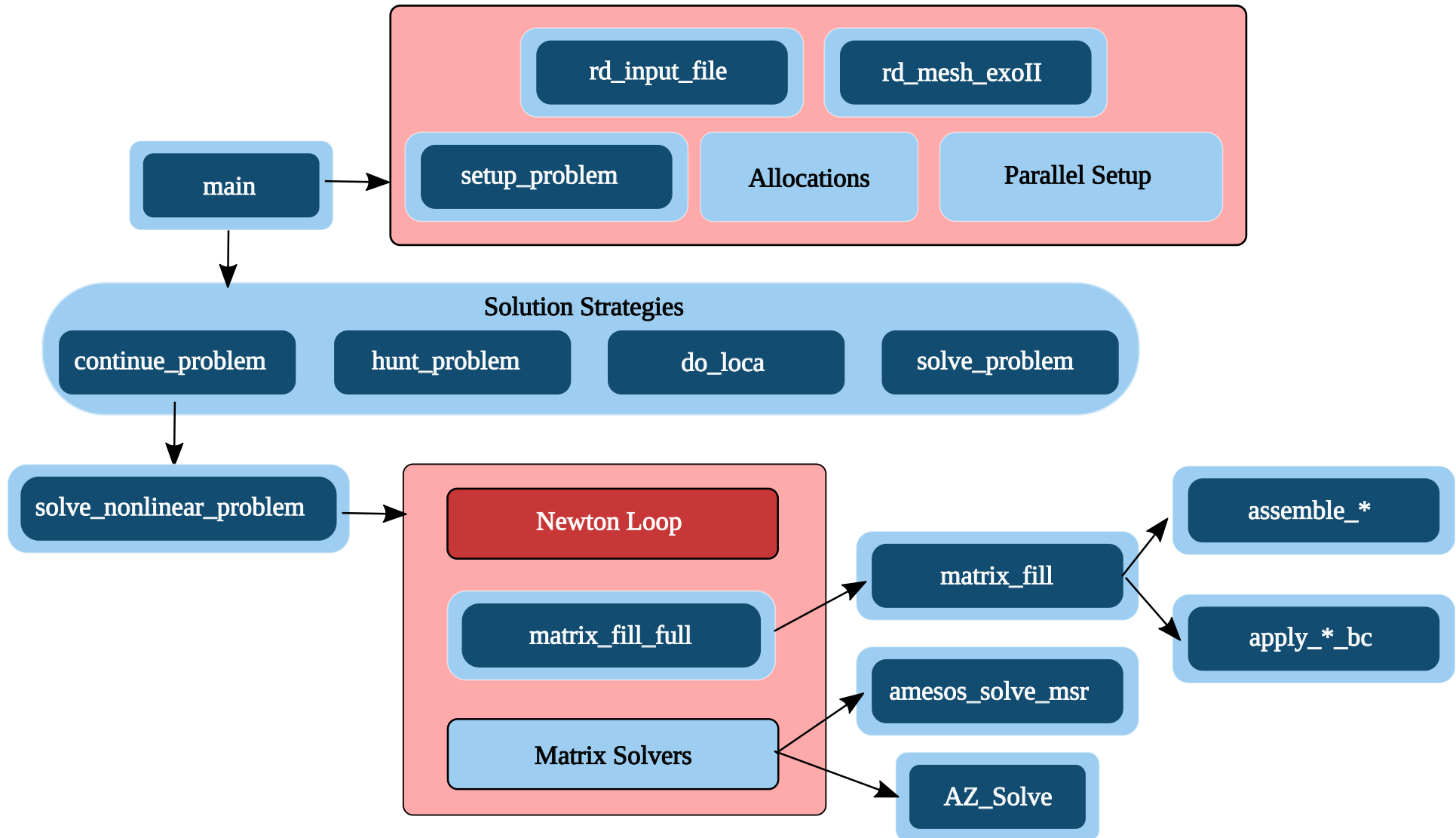
WHERE TO LOOK FOR FILES

- `mm_input*` Input for equations,BCs,materials
- `mm_fill_*` Assembly functions (Equations)
- `rf_*` Older basic finite element routines like shape functions
- `rf_solve.c` Steady and Transient solver

FILES WITH MODELS

- `mm_post_proc.c` Post Processing
- `mm_fill_terms.c` Many assembly functions
- `mm_viscosity.c` Viscosity models
- `mm_std_models.c` Many source models
- `mm_flux.c` Flux and Volume integration
- `mm_fill_species.c` Species BCs and equation
- `mm_ns_bc.c` Integrated BCs

FLOW OF GOMA



TOOLS FOR EXPLORING GOMA

The number one tool you want to have is some kind of global search

- QtCreator / VSCode / IDE's generally will have this built-in
- Ripgrep is usually faster than grep and has nice defaults and output
(<https://github.com/BurntSushi/ripgrep>)
- If none of those are available `git-grep` is your friend
(<https://git-scm.com/docs/git-grep>)
- silver searcher, ack, other grep like tools

GREP EXAMPLE

```
$ grep -R "fluid_stress"  
$ git grep "fluid_stress"  
$ rg "fluid_stress"
```


FUNCTION REFERENCE OF GOMA

FUNCTION REFERENCE OF GOMA: I/O

- I/O and General Setup:
 - `translate_command_line()`: command line arguments like `-a`, `-i`, `-brk`
 - `pd_alloc(),...*_alloc`: global allocations....
 - `read_input_file()`: Reads the input file (problem setup, enabled equations, time stepping)
 - `rd_bc_specs()`: Read boundary conditions
 - `rd_mp_specs()`: Read material properties
 - `rd_post_process_specs()`: Read Post Processing
 - Quite a few more take a look at `read_input_file`

FUNCTION REFERENCE OF GOMA: PARALLEL

- `noahs_raven()` Share some initial information with other processors
- `raven_landing()` Using above allocate structures on all processors
- `noahs_ark()` Share information from Proc 0 to all other processors
- `ark_landing()` Allocate space for structures based on above
- `noahs_dove` send information from Proc 0 to all other processors

FUNCTION REFERENCE OF GOMA: SOLUTION PROCESS

- `solve_problem()` Transient and Steady state solution
- `continue_problem()` Generic Continuation
- `hunt_problem()` Hunting continuation
- `do_loca()` Solutions using built-in loca library

FUNCTION REFERENCE OF GOMA: SPARSE MATRICES

Usually setup in Solution Process functions like `solve_problem`

- `alloc_MSR_sparse_arrays()` MSR matrix
- `alloc_VBR_sparse_arrays()` VBR matrix
- `EpetraCreateRowMatrix()` Epetra CSR
- `EpetraCreateGomaProblemGraph()` Epetra problem graph

FUNCTION REFERENCE OF GOMA: SOLUTION VECTORS

Usually setup in Solution Process functions like `solve_problem`

- `x` Solution Vector
- `x_old` Previous solution
- `xdot` Time derivative
- `x_older` Previous Previous solution

FUNCTION REFERENCE OF GOMA: ADAPTIVE TIMESTEPPING

In `solve_problem()`

- `predict_solution()` Update initial guess for next time step with explicit step, setup `xdot`
- `time_step_control()` Restrictions on time step size based on predicted solution

FUNCTION REFERENCE OF GOMA: NONLINEAR SOLVE

`solve_nonlinear_problem` Solves the nonlinear problem using Newton's method, in here we have a newton loop and call assembly and solve functions

$$\mathbf{J}(x, \dot{x}) \Delta x = -\mathbf{R}(x, \dot{x})$$
$$x = x_{old} + \Delta x$$

FUNCTION REFERENCE OF GOMA: ASSEMBLY PROCESS

In `solve_nonlinear_problem()`

- `matrix_fill_full()` Element Loop
 - `matrix_fill()` Element contributions, contains almost everything used for the finite element assembly
 - loading basis functions
 - loading field variables
 - element residuals and Jacobians
 - boundary conditions

FUNCTION REFERENCE OF GOMA: FILL DEEP DIVE

- `matrix_fill()` Element contributions
 - `load_elem_dofptr()` loads `esp` from solution vectors
 - `load_ei`
loads `ei` which contains DOF info and element information

FUNCTION REFERENCE OF GOMA: FILL DEEP DIVE

In `matrix_fill()`

- Gauss Loop

- `load_basis_functions()` loads basis functions $\phi_i, \frac{d\phi_i}{d\xi_d}$
- `beer_belly()` loads element Jacobians, $J_{el}, |J_{el}|, B$
- `load_fv()` load field variables $u = \sum_j \phi_j u_j$
- `load_bf_grad()` loads basis gradients $\nabla \phi_i$
- `load_fv_grads()` load field gradients $\nabla u = \sum_j \nabla \phi_j u_j$

FUNCTION REFERENCE OF GOMA: FILL DEEP DIVE

In `matrix_fill()`

- Gauss Loop continued
 - `assemble_*`() Compute Element Residual and Jacobian for equation
 - `assemble_momentum`() NS Momentum
 - `assemble_continuity`() NS Continuity
 - `assemble_energy`() Energy equation
 - `assemble_mass_transport`() Species equations

FUNCTION REFERENCE OF GOMA: FILL DEEP DIVE

In `matrix_fill()`

- After Gauss loop (highlights)
 - `apply_embedded_bc()` LS boundary conditions
 - `apply_integrated_bc()` Strong and Weak integrated BCs
 - `apply_point_colloc_bc()` Collocated BCs (at nodes)
 - `put_dirichlet_in_matrix()` Zero rows and puts 1 on diagonal for Dirichlet

FUNCTION REFERENCE OF GOMA: FILL DEEP DIVE

In `matrix_fill()`

- After Gauss loop (highlights)
 - `load_lec()` Local element contributions into
global: $J(global_i, global_j) + = lec_J(i, j)$,
 $R(global_i) + = lec_R(i)$