

## **GridPACK™ Overview**

This document is designed to provide an in-depth description of the GridPACK™ framework and the software modules contained within it. In combination with the Doxygen-based documentation on the GridPACK webpage, users and application developers should have a complete description of the framework components and how to use them. The applications area in the source code directory, as well as the GridPACK modules and components, provide additional examples of how GridPACK can be used to create power grid applications. However, if there are still questions on GridPACK, users should feel free to contact the GridPACK development team.

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## How to read this document

Depending on how you are planning on using GridPACK™, there are a variety of different ways of approaching the documentation. If you are only planning on using existing applications as is, without modification, then you should focus on the sections for configuring and building GridPACK and the application module documentation. Users that interested in developing their own applications may want to scan the section “Developing Applications” before going to the beginning of the document to learn about individual functionality in depth. The “GridPACK Examples” section contains additional examples of simple applications that can be used to get a sense of how to build an application from the ground up. The “Contingency Analysis” section provides some information on how to build applications that are based on the existing GridPACK application modules. Users that are interested in modifying the core functionality in GridPACK can look at the Doxygen documentation online under the “GridPACK API Documentation” link on [www.gridpack.org](http://www.gridpack.org), in addition to the documentation in this document.

## Introduction

The objective of the GridPACK™ toolkit project is to provide a framework to support the rapid development of power grid applications capable of running on high performance computing architectures (HPC) with high levels of performance and scalability. The toolkit allows power system engineers to focus on developing working applications from their models without getting bogged down in the details of decomposing the computation across multiple processors, managing data transfers between processors, working out index transformations between power grid networks and the matrices generated by different power applications, and managing input and output. GridPACK is being designed to encapsulate as much of the book-keeping required to set up HPC applications as possible using high-level programming abstractions that allow developers to concentrate on the physics and mathematics of their problems.

This document summarizes the overall design of the GridPACK framework and provides a detailed description of its components. The remainder of this document will describe the functionality incorporated into the GridPACK framework to support multiple power grid applications. The framework will continue to evolve as more real-world experience can be incorporated into the design process but many base classes that have already been identified that are capable of supporting a range of applications.

During the initial stages of GridPACK development, four power grid applications were targeted for implementation. These included:

- 1) Powerflow simulations of the electric grid
- 2) Contingency analysis of the electric grid
- 3) State estimation based on electric grid measurements
- 4) Dynamic simulations of the electric grid

From these applications, several cross-cutting functionalities were identified that could be used to support multiple applications.

- 1) Network topology and behavior. The network topology is the starting point for any power grid analysis. The topology defines the initial network model and is the connection point between the physical problem definition in terms of buses and branches and the solution method, which is usually expressed in terms of matrices and vectors.
- 2) Network components and their properties (e.g. bus and branch models, measurements, etc.). Grid components are the objects associated with the buses and branches of the power grid network. Along with the network topology itself, these define the physical system being modeled and in some cases the analysis that is to be performed. Bus and branch components can be differentiated into things like generators, loads, grounds, lines, transformers, measurements, etc. and depending on how they are defined and the level of detail incorporated into them, they define different power grid systems and analyses. The behavior of buses and branches can depend on the properties of branches or buses that are directly attached to them, e.g. figuring out the contribution of a particular bus to the solution procedure may require that properties of the attached branches are made available to the bus. The necessity for exchanging this data is built into the framework. Furthermore, these data exchanges must also be accounted for in a parallel computing context, since the grid component from which data is required may be located on a different processor.
- 3) Linear algebra and solvers. Basic algebraic objects, such as distributed matrices and vectors, are a core part of the solution algorithms required by power grid analyses. Most solution algorithms are dominated by sparse matrices but a few, such as Kalman filter analyses, require dense matrices. Vectors are typically dense. There exists a rich set of libraries for constructing distributed matrices and vectors and these are coupled to preconditioner and solver libraries. GridPACK can leverage this work heavily by creating wrappers to these libraries that can be used in solution algorithms. Wrapping these libraries instead of using them directly will have the advantage that creating algebraic objects can be simplified somewhat for power grid applications but more importantly, it will allow framework developers to investigate new solver and algebraic libraries seamlessly, without disrupting other parts of the code.
- 4) Mapping between network and algebraic objects. The physical properties of power grid systems are defined by networks and the properties of the network components but the equations describing the networks are algebraic in nature. The mappings between the physical networks and the algebraic equations depend on the indexing scheme used to describe the network and the number of parameters in the network components that appear in the equations. Constructing a map between network parameters and their corresponding locations in a matrix or vector can be complicated and error prone. Fortunately, much of this work can be automated and developers can focus more on developing code to evaluate individual matrix elements without worrying about where to locate them in the matrix. This can simplify coding considerably.

The elements described above have all been incorporated into GridPACK modules. More details about these modules and their interactions are provided in the remainder of this document.

## Configuring and Building GridPACK

**A note about CMake:** The command for invoking CMake in this manual and the documentation in <https://gridpack.org> is usually of the form

```
cmake [OPTIONS] ..
```

This particular form assumes that the build directory is below the directory that contains the top-level **CMakeLists.txt** file for the build. For GridPACK, this is located in the **src** directory. If your build directory for GridPACK is below **src** and you invoke CMake from this directory, the “**..**” at the end of the **cmake** command is pointing to **src**. You could also use the absolute path to the **src** directory instead of “**..**” and this would work no matter where you locate the build directory.

Building GridPACK requires several external libraries that must be built prior trying to configure and build GridPACK itself. On some systems, these libraries may already be available but in many cases, users will need to build them by hand. An exception is MPI, which is usually available on parallel platforms, although users interested in running parallel jobs on a multi-core workstation may still need to build it themselves. In any case, the best way to guarantee that all libraries are compatible with each other is to build them all using a consistent environment and set of compilers. There is extensive documentation on how to build GridPACK and the libraries on which it depends on the website located at <https://gridpack.org>. We refer to the information on the website for most of the details on how to build GridPACK and will only discuss some general properties of the configure procedure in this document.

Example scripts for building the libraries used by GridPACK on different systems can be found under **\$GRIDPACK/src/scripts**. In most cases these need to be modified slightly before they will work on your system, but the changes are usually small and self-evident. The scripts contain some additional documentation at the top to help you with these modifications. Find a script for a platform that is similar to your system and use this as the starting point for your build.

GridPACK uses the CMake build system to create a set of make files that can then be used to compile the entire GridPACK framework. Most of the effort in building GridPACK is focused on getting the configure process to work, once configure has been successfully completed, compilation is usually straightforward. Builds of GridPACK should be done in their own directory and this also makes it possible to have multiple builds that use different configuration parameters associated with the same source tree. Typically, the build directories are under **\$GRIDPACK/src** directory but they can be put anywhere the user chooses. The user then needs to run CMake from the build directory to configure GridPACK and then **make** and **make install** to compile and install the GridPACK libraries. After running **make**, all applications



in the GridPACK source tree are also available for use. The application executables will be located in the build directory and not in the source tree.

GridPACK currently makes use of five different libraries. MPI and Global Arrays are used for communication, Boost provides several C++ extensions used throughout GridPACK, Parmetis is used to partition networks over multiple processors and PETSc provides parallel solvers and algebraic functionality. Except for MPI, which is usually available through compiler wrappers such as **mpicc** and **mpicxx**, the locations of the remaining libraries need to be specified in the CMake configure command.

Because the cmake command takes a large number of arguments, it is usually a good idea to put the entire command in a script. The script can then be edited as needed. Make sure that the script is executable by running the **chmod +x** command on it. A typical CMake configure script is

```
rm -rf CMake*

cmake -Wdev \
  -D BOOST_ROOT:STRING='$HOME/software_new/boost_1_55_0' \
  -D PETSC_DIR:STRING='$HOME/software_new/petsc-3.6.0' \
  -D PETSC_ARCH:STRING='linux-openmpi-gnu-cxx' \
  -D PARMETIS_DIR:STRING= \
    '$HOME/software_new/petsc-3.6.0/linux-openmpi-gnu-cxx/lib' \
  -D GA_DIR:STRING='$HOME/software_new/ga-5-4-ib' \
  -D USE_PROGRESS_RANKS:BOOL=FALSE \
  -D GA_EXTRA_LIBS='-lrt -libverbs' \
  -D MPI_CXX_COMPILER:STRING='mpicxx' \
  -D MPI_C_COMPILER:STRING='mpicc' \
  -D MPIEXEC:STRING='mpiexec' \
  -D CMAKE_INSTALL_PREFIX:PATH='$GRIDPACK/src/build/install' \
  -D CMAKE_BUILD_TYPE:STRING='RELWITHDEBINFO' \
  -D MPIEXEC_MAX_NUMPROCS:STRING="2" \
  -D CMAKE_VERBOSE_MAKEFILE:STRING=TRUE \
  ..
```

The first line removes any configuration files that may be left over from a previous configuration attempt. Removing these files is generally a good idea since parameters from a previous unsuccessful attempt may bleed over into the current configuration and either spoil the configuration itself or lead to problems when you try to compile the code. The Boost, PETSc, Parmetis and Global Array library locations are specified by the **BOOST\_ROOT**, **PETSC\_DIR**, **PARMETIS\_DIR** and **GA\_DIR** variables. The **PETSC\_ARCH** variable specifies the particular build within PETSc that you want GridPACK to use. It is usually possible when configuring and

building PETSc to have it download and build Parmetis as well. This was done in the example above and thus the Parmetis libraries are located within the PETSc source tree in the directory corresponding to the architecture specified in **PETSC\_ARCH**.

The Global Arrays library can be built using a number of different runtimes. The default runtime uses MPI two-sided communication. While it is very easy to use, this runtime does not scale well beyond a dozen or so processors. Users interested on running on large numbers of cores should look at configuring Global Arrays with other runtimes. A high performing GA runtime that is available on most platforms is called progress ranks. This runtime has a peculiarity in that it reserves one MPI process per SMP node to manage communication. Thus, if you request a total of 20 MPI processes on 4 nodes with 5 processes running on each node only 4 MPI process per node will actually be available to the application for a total of 16. In order to notify GridPACK that you are using this runtime, you need to set the parameter **USE\_PROGRESS\_RANKS** to true. In the example above, we are not using progress ranks so we set **USE\_PROGRESS\_RANKS** to false.

The **GA\_EXTRA\_LIBS** parameter can be used to include extra libraries in the link step that are not picked up as part of the configuration process. In this example, GA is configured to run on an Infiniband network so it is necessary to explicitly include **libibverbs** and **librt**. For most of the MPI-based runtimes, this variable is not needed.

The MPI wrappers for the C and C++ compilers can be specified by setting **MPI\_C\_COMPILER** and **MPI\_CXX\_COMPILER** and the MPI launch command can be specified using **MPIEXEC**. The **CMAKE\_INSTALL\_PREFIX** specifies the location of the installed build of GridPACK. This location should be used when linking external applications to GridPACK. The **CMAKE\_BUILD\_TYPE** can be used to control the level of debugging symbols in the library. **MPIEXEC\_NUM\_PROCS** should be set to a small number and controls the number of processors that will be used if running the parallel tests in the GridPACK test suite. Many of the application tests are small (9 or 14 buses) and will fail if you try and run on a large number of cores. Finally, **CMAKE\_VERBOSE\_MAKEFILE** controls the level of information generated during the compilation. It is mainly of interest for people doing development in GridPACK and most other users can safely set it to false.

The final argument of the cmake command is the location of the top level **CMakeLists.txt** file in the source tree. For GridPACK, this file is located in the **\$GRIDPACK/src** directory. The above example assumes that the build directory is located directly under **\$GRIDPACK/src** so the **..** at the end of the configure script is pointing to the directory containing the **CMakeLists.txt** file.

Once the GridPACK framework has been built, applications and framework tests can be run using standard MPI scripts for running jobs. A typical invocation to run a code `code.x` on some number of processors is

```
mpirun -n 2 code.x
```

In this case the code will run on 2 processors. Different platforms may use different scripts to run the parallel job. Consult your local system documentation for details. Applications may also have additional arguments that are processed inside the application itself. Most GridPACK applications will take an argument representing the input file for the application.

## Building GridPACK Applications

GridPACK comes with several applications that are included in the main distribution. These currently include power flow, contingency analysis, dynamic simulation, state estimation and Kalman filter applications as well as some non-power grid examples that illustrate features of the framework. These applications are automatically built whenever the full GridPACK distribution is built.

For applications developed outside the GridPACK distribution, the build process is fairly simple, provided you are using CMake (you will need to have CMake installed on your system to build GridPACK so using CMake for your application build should be a straightforward extension). For a CMake build, you need to create a CMakeLists.txt file in the same directory that includes your application files. A template for the CMakeLists.txt file is

```
1 cmake_minimum_required(VERSION 2.6.4)
2
3 if (NOT GRIDPACK_DIR)
4   set(GRIDPACK_DIR /HOME/gridpack-install
5       CACHE PATH "GridPACK installation directory")
6 endif()
7
8 include("${GRIDPACK_DIR}/lib/GridPACK.cmake")
9
10 project(MyProject)
11
12 enable_language(CXX)
13
14 gridpack_setup()
15
16 add_definitions(${GRIDPACK_DEFINITIONS})
17 include_directories(BEFORE ${CMAKE_CURRENT_SOURCE_DIR})
18 include_directories(BEFORE ${GRIDPACK_INCLUDE_DIRS})
```

```

19
20 add_executable(myapp.x
21     myapp_main.cpp
22     myapp_driver.cpp
23     myapp_file1.cpp
24     myapp_file2.cpp
25 )
26 target_link_libraries(myapp.x ${GRIDPACK_LIBS})
27
28 add_custom_target(myapp.input
29
30     COMMAND ${CMAKE_COMMAND} -E copy
31     ${CMAKE_CURRENT_SOURCE_DIR}/input.xml
32     ${CMAKE_CURRENT_BINARY_DIR}
33
34     COMMAND ${CMAKE_COMMAND} -E copy
35     ${CMAKE_CURRENT_SOURCE_DIR}/myapp_test.raw
36     ${CMAKE_CURRENT_BINARY_DIR}
37
38     DEPENDS
39     ${CMAKE_CURRENT_SOURCE_DIR}/input.xml
40     ${CMAKE_CURRENT_SOURCE_DIR}/myapp_test.raw
41 )
42 add_dependencies(myapp.x myapp.input)

```

Lines 1-6 check to see if the CMake installation is recent enough and also make sure that the **GRIDPACK\_DIR** variable has been defined in the configuration step. If it hasn't, then the CMake will try and use a default value and look for a build in **\$HOME/gridpack-install**. However, this is unlikely to be successful, so it is better to define **GRIDPACK\_DIR** when configuring your application. Line 8 picks up a file that is used by the application build to link to libraries and header files in the GridPACK build and line 10 can be used to assign a name to your application. Lines 12-18 can be included as is, if all application files are in the same directory as the CMakeLists.txt file. If other directories contain source and header files, then they can be included using the directives in lines 17 and 18.

Lines 20-25 define the name of the executable and all the source code files that are used in the application. The **add\_executable** command on line 26 adds the executable **myapp.x** to the build. The arguments to this command consist of the name of the executable followed by the executable source files. There can be an arbitrary number of source files associated with any one executable. Note that the source files just consist of the user application source files, the

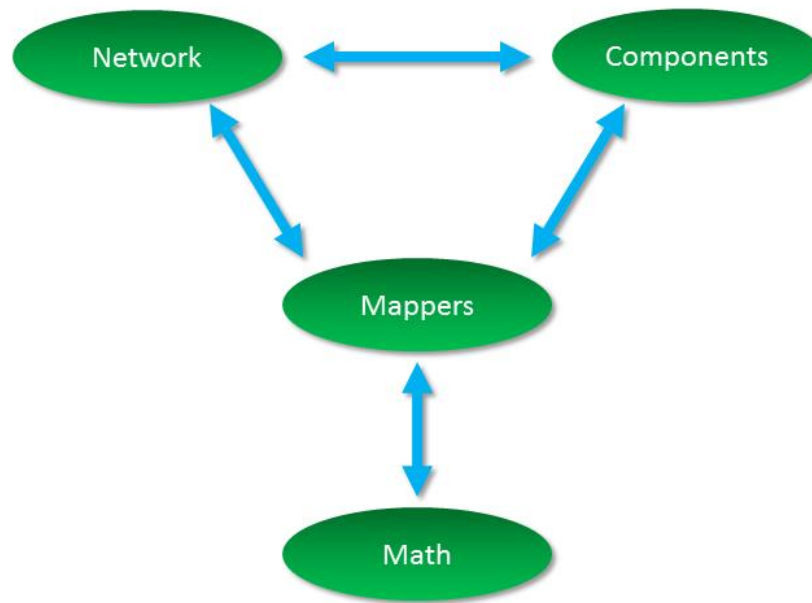
framework files are handled automatically. If some of the files are located in subdirectories, then the path relative to the directory where the CMakeLists.txt file is located should be included.

The remaining lines 28-42 are optional and can be used to automatically copy files from the application source file directory to the build directory. These could include example input files or external configuration files that are called by the code to set internal parameters. The **add\_custom\_target** command on line 28 defines a list of files and what should be done with them. In this example, the two files **input.xml** and **myapp\_test.raw** are the files to be copied. The **COMMAND** line specifies the action (copy) and the next two lines specify the location of the file to be copied and its destination. The **DEPENDS** keyword (line 38) indicates that any time the **input.xml** or **myapp\_test.raw** files are modified, they should be recopied to the build directory if make is invoked and the **add\_dependencies** command (line 42) binds the custom target to the build of the executable.

A template file for **CMakeLists.txt** can be found in the **src** directory under **CMakeLists.template.txt**. Users should copy this file to their application directory, modify the name to **CMakeLists.txt** and add their own source files and test input.

## GridPACK Framework Components

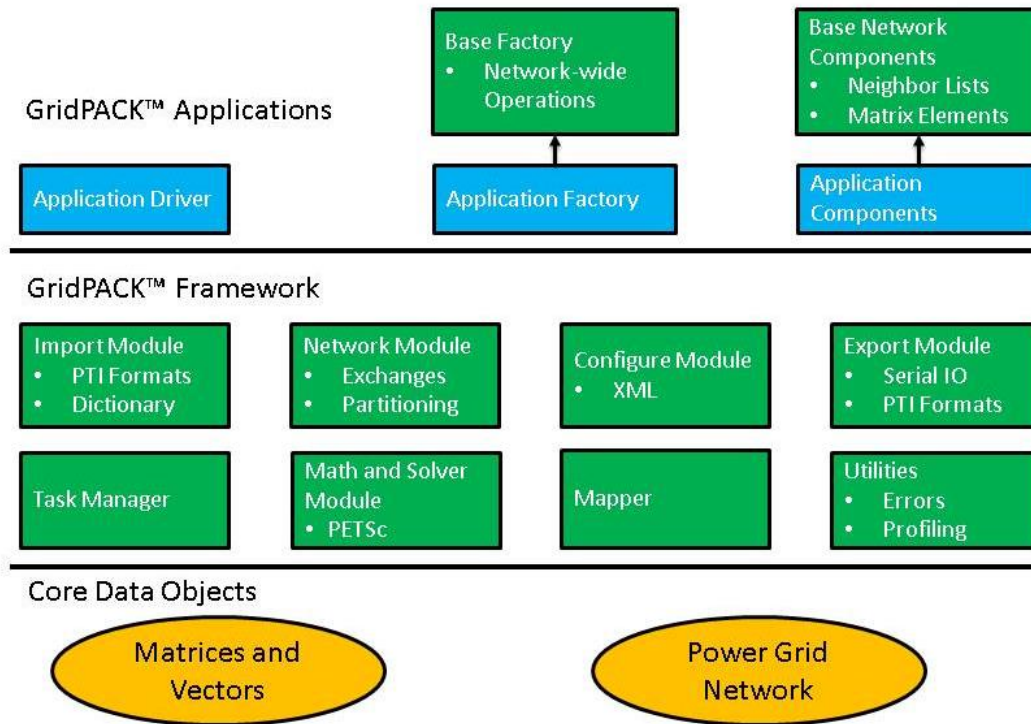
This section will describe the GridPACK components and the functionality they support. The four major GridPACK components are networks, bus and branch components, the mappers and the math module. The math module is relatively self-contained and can be used as a conventional library, but the other three are tightly coupled and need to be used together to do anything useful. A schematic that illustrates the relationship between these components is shown in Figure 1.



**Figure 1.** Relationship between major GridPACK components.

A full description of a power grid network requires specification of both the network topology and the physical properties of the bus and branch components. The combination of the models and the network generate algebraic equations that can be solved to get desired system properties. GridPACK supplies numerous modules to simplify the process of specifying the model and solving it. These include power grid components that describe the physics of the different network models or analyses, grid component factories that initialize the grid components, mappers that convert the current state of the grid components into matrices and vectors, solvers that supply the preconditioner and solver functionality necessary to implement solution algorithms, input and output modules that allow developers to import and export data, and other utility modules that support standard code develop operations like timing, event logging, and error handling.

Many of these modules are constructed using libraries developed elsewhere so as to minimize framework development time. However, by wrapping them in interfaces geared towards power grid applications these libraries can be made easier to use by power grid engineers. The interfaces also make it possible in the future to exchange libraries for new or improved implementations of specific functionality without requiring application developers to rewrite their codes. This can significantly reduce the cost of introducing new technology into the framework. The software layers in the GridPACK framework are shown schematically in Figure 2.



**Figure 2.** A schematic diagram of the GridPACK framework software data stack. Green represents components supplied by the framework and blue represents code that is developed by the user.

Core framework components are described below. Before discussing the components themselves, some of the coding conventions and libraries used in GridPACK will be described.

**Preliminaries:** The GridPACK software uses a few coding conventions to help improve memory management and to minimize run-time errors. The first of these is to employ namespaces for all GridPACK modules. The entire GridPACK framework uses the **gridpack** namespace, individual modules within GridPACK are further delimited by their own namespaces. For example, the BaseNetwork class discussed in the next section resides in the **gridpack::network** namespace and other modules have similar delineations. The example applications included in the source code also have their own namespaces, but this is not a requirement for developing GridPACK-based applications.

To help with memory management, many GridPACK functions return boost shared pointers instead of conventional C++ pointers. These can be converted to a conventional pointer using the **get()** command. We also recommend that the type of pointers be converted using a **dynamic\_cast** instead of conventional C-style cast.

Application files should include the **gridpack.hpp** header file. This can be done by adding the line

```
#include "gridpack/include/gridpack.hpp"
```

at the top of the application .hpp and/or .cpp files. This file contains definitions of all the GridPACK modules and their associated functions.

Matrices and vectors in GridPACK were originally complex but now either complex or real matrices can be created using the library. Inside the GridPACK implementation, the underlying distributed matrices are either complex or real, but the framework adds a layer that supports both types of objects, even if the underlying math library does not. However, computations on complex matrices will perform better if the underlying math library is configured to use complex matrices directly. This should be kept in mind when choosing the math library to build GridPACK on. The underlying PETSc library can be configured to support either real or complex matrices. Complex numbers are represented in GridPACK as having type **ComplexType**. The real and imaginary parts of a complex number **x** can be obtained using the functions **real(x)** and **imag(x)**.

## Network Module

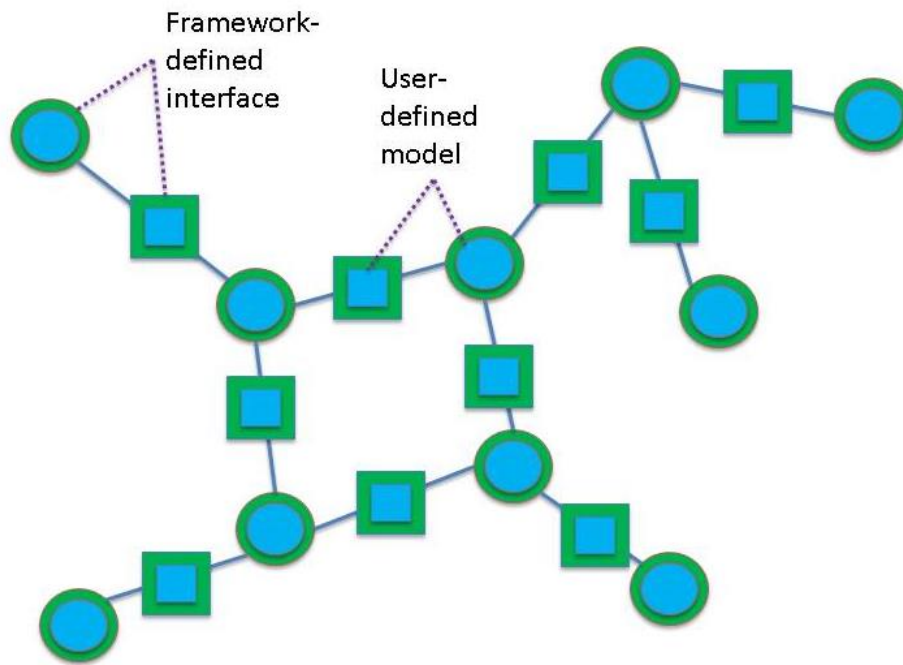
The network module is designed to represent the power grid and has four major functions

- 1) The network is a container for the grid topology. The connectivity of the network is maintained by the network object and can be made available through requests to the network. The network also maintains the “ghost” status of buses and branches and determines whether a bus or branch is owned by a particular processor or represents a ghost image of a bus or branch owned by a neighboring processor.
- 2) The network topology can be decorated with bus and branch objects that describe the properties of the particular physical system under investigation. Bus and branch objects are written by the application developer and incorporate the grid model and the analyses that need to be performed on it. Different applications will use different bus and branch implementations.
- 3) The network module is responsible for supplying update operations that can be used to fill in the value of ghost cell fields with current data from other processors. The updates of ghost buses and ghost branches have been split into separate operations to give users flexibility in optimizing performance by minimizing the amount of data that needs to be communicated in the code. Many applications do not require exchanges of branch data.
- 4) The network contains the partitioner. The partitioner is embedded in the network module but it is a substantial computational technology in its own right. Partitioning is a key part of parallel application development. It represents the act of dividing up the problem so that each processor is left with approximately equal amounts of work. At the same time,



the partition is chosen so that communication between processors (a major source of computational inefficiency in HPC programs) is minimized.

A network is illustrated schematically in Figure 3. Each bus and branch has an associated bus or branch object. The buses and branches are derived from base classes that specify certain functions that must be implemented by the application developer so that the network can interact with other GridPACK modules. In addition, the application can have functionality outside the base class that is unique to the particular application.



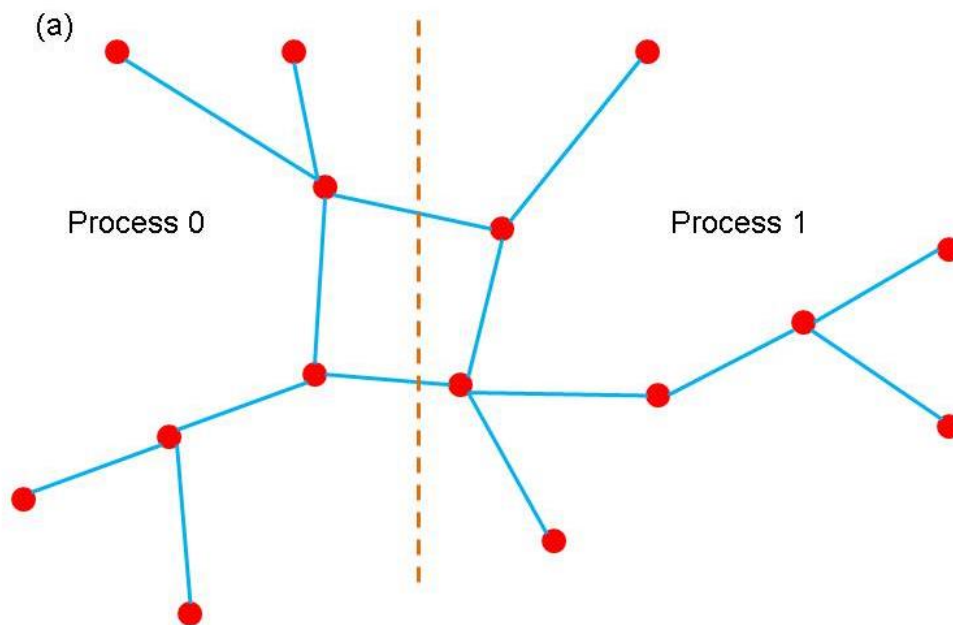
**Figure 3.** Schematic representation of a GridPACK network. The squares are branch objects and the circles are bus objects. Framework-specified interfaces are green and user supplied functionality is blue.

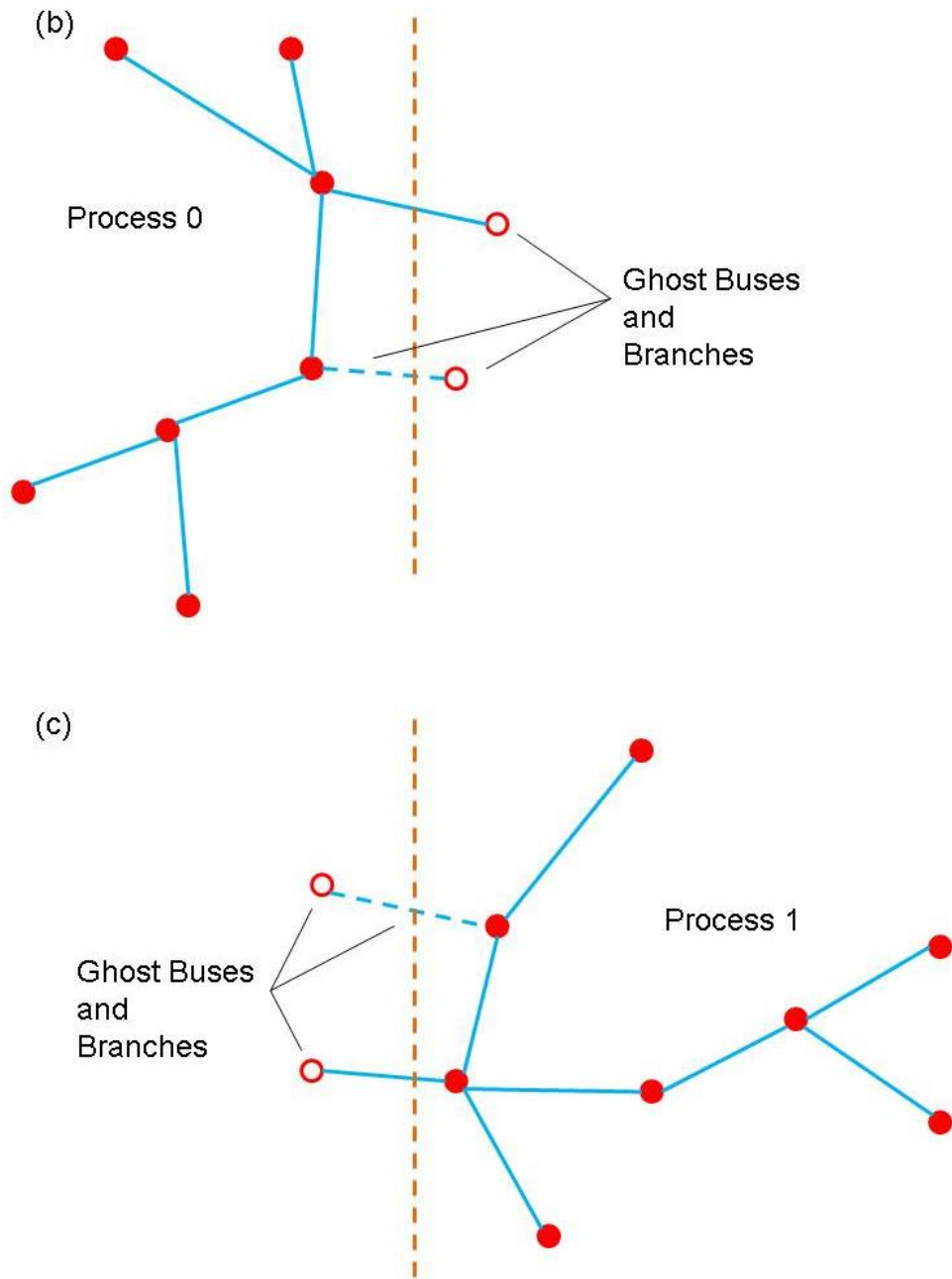
A major use of the partitioner is to rearrange the network in a form that is useful for computation immediately after it is read in from an external file. Typically, the information in the external file is not organized in a way that is necessarily optimal for computation, so the partitioner must redistribute data such that each processor contains at most a few large connected subsets of the network. The partitioner is also responsible for adding the ghost buses and branches to the system.

Ghost buses and branches in a parallel program represent images of buses and branches that are owned by other processes. In order to carry out operations on buses and branches it is frequently necessary to gain access to data associated with attached buses and branches. The most efficient way to do this is to create copies of the buses and branches from other processors that are

connected to locally owned objects. All local network components then have a complete set of attached neighbors. The ghost objects are updated collectively with current information from their home processors at points in the computation. Updating all ghosts at once is almost always more efficient than accessing data from one remote bus or branch at a time.

The use of the partitioner to distribute the network between different processors and create ghost nodes and branches is illustrated in Figure 4. Figure 4(a) shows a simple network and Figures 4(b) and 4(c) show the result of distributing the network between two processors.





**Figure 4.** (a) a simple network (b) partition of network on processor 0 (b) partition of network on processor 1. Open circles indicate ghost buses and dotted lines indicate ghost branches.

Networks can be created using the templated base class **BaseNetwork**<class **Bus**, class **Branch**>, where **Bus** and **Branch** are application-specific classes describing the properties of buses and branches in the network. The **BaseNetwork** class is defined within the **gridpack::network** namespace. In addition to the **Bus** and **Branch** classes, each bus and branch has an associated **DataCollection** object, which is described in more detail in the

network components section. The **DataCollection** object is a collection of key-value pairs that acts as an intermediary between data that is read in from external configuration files and the bus and branch classes that define the network.

The **BaseNetwork** class contains a large number of methods, but only a relatively small number will be of interest to application developers. The remaining methods are used primarily within other GridPACK modules to implement higher level capabilities. This document will focus on calls that are likely to be used by application developers.

The constructor for the network class is the function

**BaseNetwork(const parallel::Communicator &comm)**

The **Communicator** object is used to define the set of processors over which the network is distributed. Communicators are discussed in more detail below. The network constructor creates an empty shell that does not contain any information about an actual network. The remainder of the network must be built up by adding buses and branches to it. Typically, buses and branches are added by passing the network to a parser (see import module) which will create an initial version of the network. The constructor is paired with a corresponding destructor

**~BaseNetwork()**

that is called when the network object passes out of scope or is explicitly deleted by the user.

Two functions are available that return the number of buses or branches that are available on a process. This number includes both buses and branches that are held locally as well as any ghosts that may be located on the process.

**int numBuses()**

**int numBranches()**

There are also functions that will return the total number of buses or branches in the network. These numbers ignore ghost buses and ghost branches.

**int totalBuses()**

**int totalBranches()**

Buses and branches in the network can be identified using a local index that runs from 0 to the number of buses or branches on the process minus 1 (0-based indexing). For some calculations, it is necessary to identify one bus in the network as a reference bus. This bus is usually set when the network is created using an import parser. It can subsequently be identified using the function

**int getReferenceBus()**

If the reference bus is located on this processor (either as a local bus or a ghost) then this function returns the local index of the bus, otherwise it returns -1.

Ghost buses and branches are distinguished from locally owned buses and branches based on whether or not they are “active”. The two functions

```
bool getActiveBus(int idx)
```

```
bool getActiveBranch(int idx)
```

provide the active status of a bus or branch on a process. The index **idx** is a local index for the bus or branch.

Buses and branches are characterized by a number of different indices. One is the local index, already discussed above, but there are several others. Most of these are used internally by other parts of the framework but one index is of interest to application developers. This is the “original” bus index. When the network is described in the input file, the buses are labeled with a (usually) positive integer. There are no requirements that these integers be consecutive, only that each bus has its own unique index. The value of this index can be recovered using the function

```
int getOriginalBusIndex(int idx)
```

The variable **idx** is the local index of the bus. Branches are usually described in terms of the original bus indices for the two buses at each end of the branch, so there is no corresponding function for branches. Instead, the procedure is to get the local indices of the two buses at each end of the branch and then get the corresponding original indices of the buses. This information is usually used for output.

It is frequently necessary to gain access to the objects associated with each bus or branch. The following four methods can be used to access these objects

```
boost::shared_ptr<Bus> getBus(int idx)
```

```
boost::shared_ptr<Branch> getBranch(int idx)
```

```
boost::shared_ptr<DataCollection> getBusData(int idx)
```

```
boost::shared_ptr<DataCollection> getBranchData(int idx)
```

The first two methods can be used to get Boost shared pointers to individual bus or branch objects indexed by local indices **idx**. The second two functions return pointers to the **DataCollection** objects associated with each bus or branch. These **DataCollection** objects can be used to initialize the bus and branch objects at the start of a calculation but they

are also useful when converting a network of one type to a network of another type. This often happens when different computations are chained together.

The following functions can be useful for handling input that is directed at certain network components

```
std::vector<int> getLocalBusIndices(int idx)  
std::vector<int> getLocalBranchIndices(int idx1, int idx2)
```

These functions return a list of local indices that correspond to either the original bus index **idx** for a bus, or the pair of indices **idx1**, **idx2** for a branch. The reason that a list is returned instead of a single index is that in the case of ghost buses and branches, more than one copy of a network component may exist on a process. If no copies of a network component exist on a process then the returned vector has zero length. These functions can be used for applications such as contingency analysis, where modifications are made to a single network component and the modifications are specified in terms of the original bus indices. These functions can be used to find the local index of the component, if it exists.

The network partitioner can be accessed via the function

```
void partition()
```

The partition function distributes the buses and branches across processors such that the connectivity to branches and buses on other processors is minimized. It is also responsible for adding ghost buses and branches to the network. This function should be called after the network is read in but before any other operations, such as setting up exchange buffers or creating neighbor lists, have been performed.

Finally, two sets of functions are required in order to set up and execute data exchanges between buses and branches in a distributed network. These exchanges are used to move data from active components to ghost components residing on other processors. Before these functions can be called, the buffers in individual network components must be allocated. See the documentation below on network components and the network factory for more information on how to do this. Once the buffers are in place, bus and branch exchanges can be set up and executed with just a few calls. The functions

```
void initBusUpdate()
```

```
void initBranchUpdate()
```

are used to initialize the data structures inside the network object that manage data exchanges. Exchanges between buses and branches are handled separately, since not all applications will require exchanges between both sets of objects. The initialization routines are relatively complex

and allocate several large internal data structures, so they should not be called if there is no need to exchange data as part of the algorithm.

After the updates have been initialized, it is possible to execute a data exchange at any point in the code by calling the functions

```
void updateBuses()
```

```
void updateBranches()
```

These functions will cause data on ghost buses and branches to be updated with current values from active buses and branches located on other processors.

One additional network function that can be useful in certain circumstances is the capability for recovering the communicator on which the network is defined

```
const Communicator& communicator() const
```

This function can be used in implementing algorithms based on multilevel parallelism. Recovering the communicator is also needed for converting applications to modules that can be used to create higher level workflows that combine multiple different types of applications. This is discussed in more detail below.

The **BaseNetwork** methods described in this section are only a subset of the total functionality available but they represent most of the methods that a typical developer would use. The remaining functions are primarily used to implement other parts of the GridPACK framework but are generally not required by people writing applications. More information on how the functions described above are used in practice can be found in the section on GridPACK factories.

## **Math Module**

The math module provides support in GridPACK for distributed matrices and vectors, linear solvers, non-linear solvers, and preconditioners. Once created, matrices can be treated as opaque objects and manipulated using a high level syntax that is comparable to writing Matlab code. The distributed matrix and vector data structures themselves are based on external solver libraries and represent relatively lightweight wrappers on multipurpose HPC codes. The current math module is built on the PETSc library but other libraries, such as Hypre and Trilinos could potentially be used instead.

The main functionality associated with the math module is the ability to instantiate new matrices and vectors, add individual matrix and vector elements (and their values) to the matrix/vector objects, invoke the assemble operation on the object, perform basic algebraic operations, such as matrix-vector multiply, and solve systems of algebraic equations. The assemble operation is designed to give the library a chance to set up internal data structures and repartition the matrix

elements, etc. in a way that will optimize subsequent calculations. Inclusion of this operation follows the syntax of most solver libraries when they construct a matrix or vector.

In addition to basic matrix operations, the math module contains linear and non-linear solvers and preconditioners. The module provides a simple interface on top of the PETSc libraries that will allow users access to this functionality without having to be familiar with the libraries themselves. This should make it possible to construct solver routines that are comparable in complexity to Matlab scripts. The use of a wrapper instead of having users directly access the libraries will also make it simpler to switch the underlying library in an application. All that will be required is that developers link to an implementation of the math module interface that is built on a different library. There will not be a need to rewrite any application code. This has the advantage that if a different library is used for the math module in one application, it instantly becomes available for other applications.

The functionality in the math component is distributed between the classes **Matrix**, **RealMatrix**, **Vector**, **RealVector**, **LinearSolver**, **RealLinearSolver**, **NonlinearSolver** and **RealNonlinearSolver**. Each of these classes is in the **gridpack::math** namespace and is described below. Like the **BaseNetwork** class, there are a lot of functions in **Matrix** and **Vector** that do not need to be used by users. Most of the functions related to matrix/vector instantiation and creation are used inside the mapper classes described below, which eliminates the need for users to deal with them directly. However, users may be interested in creating functions not covered by existing library methods and in this case access to these functions is useful.

An additional note on the math module class names is in order. Originally, GridPACK only supported complex objects and used the names **Vector**, **Matrix**, etc. More recently, the capability for supporting real objects was added and hence the new names **RealVector**, etc. The original names continued to be used for complex objects to maintain backwards compatibility. Complex objects can also be accessed using the names **ComplexVector**, **ComplexMatrix**, etc., which are mapped to the original complex objects.

### *Matrices*

The **Matrix** and **RealMatrix** classes are designed to create distributed matrices. **Matrix** is used for complex matrices and **RealMatrix** is used for real matrices. The matrix classes support two types of matrix, **Dense** and **Sparse**. In most cases users will want to use the sparse matrix but some applications require dense matrices. The **Matrix** and **RealMatrix** classes are nearly identical in functionality, so in the following we will only outline operations on the **Matrix** class. In most cases, the **RealMatrix** class contains the same operations. The only point to note is that for any operations that involve multiple matrices or a matrix and a vector, all matrix and vector objects must be either all complex or all real. In the future, we plan on adding some operations that will allow users to convert between types.



The matrix constructor is

```
Matrix(const parallel::Communicator &comm,  
       const int &local_rows,  
       const int &cols,  
       const StorageType &storage_type=Sparse)
```

The communicator object **comm** specifies the set of processors that the matrix is defined on, the **local\_rows** parameter corresponds to the number of rows contributed to the matrix by the processor, the **cols** parameter indicates what the second dimension of the matrix is and the **storage\_type** parameter determines whether the matrix is sparse or dense. If the total dimension of the matrix is  $M \times N$ , then the sum of the **local\_rows** parameters over all processors must equal  $M$  and the **cols** parameter is equal to  $N$ . The matrix destructor is

```
~Matrix()
```

Once a matrix has been created some inquiry functions can be used to probe the matrix size and distribution. The following functions return information about the matrix.

```
int rows() const
```

```
int localRows() const
```

```
void localRowRange(int &lo, int &hi) const
```

```
int cols()
```

The function **rows** will return the total number of rows in the matrix, **localRows** returns the number of rows associated with the calling processor, **localRowRange** returns the **lo** and **hi** index of the rows associated with the calling processor and **cols** returns the number of columns in the matrix. Note that matrices are partitioned into row blocks on each processor.

Additional functions can be used to add matrix elements to the matrix, either one at a time or in blocks. The following two calls can be used to reset existing elements or insert new ones.

```
void setElement(const int &i, const int &j,  
               const ComplexType &x)
```

```
void setElements(const int &n, const int *i, const int *j,  
                const ComplexType *x)
```

For real matrices, all variables of type **ComplexType** should be switched to type **double**. The first function will set the matrix element at the index location **(i, j)** to the value **x**. If the

matrix element already exists, this function overwrites the value, if the element is not already part of the matrix, it gets added with the value **x**. Note that both **i** and **j** are zero-based indices. For the current PETSc based implementation of the math module, it is not required that the index **i** lie between the values of **lo** and **hi** obtained with **localRowRange** function, but for performance reasons it is desirable. Other implementations may require that **i** lie in this range. The second function can be used to add a collection of elements all at once. The variable **n** is the number of elements to be added, the arrays **i** and **j** contain the row and column indices of the matrix elements and the array **x** contains their values. Again, it is preferable that all values in **i** lie within the range [**lo**,**hi**].

Two functions that are similar to the set element functions above are the functions

```
void addElement(const int &i, const int &j,  
               const ComplexType &x)
```

```
void addElements(const int &n, const int *i, const int *j,  
                const ComplexType *x)
```

These differ from the set element functions only in that instead of overwriting the new values into the matrix, these functions will add the new values to whatever is already there. If no value is present in the matrix at that location the function inserts it.

In addition to setting or adding new elements, it is possible to retrieve matrix values using the functions

```
void getElement(const int &i, const int &j,  
               ComplexType &x) const
```

```
void getElements(const int &n, const int *i, const int *j,  
                ComplexType *x) const
```

These functions can only access elements that are local to the processor. This means that the index **i** must lie in the range [**lo**,**hi**] returned by the function **localRowRange**.

Finally, before a matrix can be used in computations, it must be assembled and internal data structures must be set up. This can be accomplished by calling the function

```
void ready()
```

After this function has been invoked, the matrix is read for use and can be used in computations. In general, the procedure for building a matrix is 1) create the matrix object 2) determine local

parameters such as **lo** and **hi** 3) set or add matrix elements and 4) assemble the matrix using the **ready** function. For most applications, users can avoid these operations by building matrices and vectors using the mapper functionality described below.

Some additional functions have been included in the matrix class that can be useful for creating matrices or writing out their values (e.g. for debugging purposes). It is often useful to create a copy of a matrix. This can be done using the clone method

**Matrix\* clone() const**

The new matrix is an exact replica of the matrix that invokes this function.

Two functions that can be used to write the contents of a matrix, either to standard output or to a file are

**void print (const char \*filename=NULL) const**  
**void save(const char \*filename) const**

The first function will write the contents of the matrix to standard output if no filename is specified, otherwise it writes to the specified file, the second function will write a file in MatLAB format. These functions can be used for debugging or to create matrices that can be fed into other programs.

Once a matrix has been created, a variety of methods can be applied to it. Most of these are applied after the **ready** call has been made by the matrix, but some operations can be used to actually build a matrix. These functions are listed below.

**void equate(const Matrix &A)**

This function sets the calling matrix equal to matrix **A**.

**void scale(const ComplexType &x)**

Multiply all matrix elements by the value **x** (use a value of type **double** for a real matrix).

**void multiplyDiagonal(const Vector &x)**

Multiply all elements on the diagonal of the calling matrix by the corresponding element of the vector **x**. The **Vector** class is described below.

**void addDiagonal(const Vector &x)**

Add elements of the vector **x** to the diagonal elements of the calling matrix.

**void add(const Matrix &A)**

Add the matrix **A** to the calling matrix. The two matrices must have the same number of rows and columns, but otherwise there are no restrictions on the data layout or the number and location of the non-zero entries.

#### **void identity()**

Create an identity matrix. This function assumes that the calling matrix has been created but no matrix elements have been assigned to it.

#### **void zero()**

Set all non-zero entries to zero.

#### **void conjugate(void)**

Set all entries to their complex conjugate value. This function only applies to complex matrices.

The following functions create a new matrix or vector.

#### **Matrix \*multiply(const Matrix &A, const Matrix& B)**

Multiply matrix **A** times matrix **B** to create a new matrix.

#### **Vector \*multiply(const Matrix &A, const Vector &x)**

Multiply matrix **A** times vector **x** to get a new vector.

#### **Matrix \*transpose(const Matrix &A)**

Take the transpose of matrix **A**.

### **Vectors**

The vector class operates in much the same way as the matrix class. As above, most functions apply to both the **Vector** and **RealVector** class so only the **Vector** operations are described here. The vector constructor is

#### **Vector(const parallel::Communicator& comm, const int& local\_length)**

The parameter **local\_length** is the number of contiguous elements in the vector that are held on the calling processor. The sum of **local\_length** over all processors must equal the total length of the vector. The functions

#### **int size(void) const**

#### **int localSize(void) const**

#### **void localIndexRange(int &lo, int &hi) const**

can be used to get the global size of the vector or the size of the vector segment held locally on the calling processor. The **localIndexRange** function can be used to find the indices of the vector elements that are held locally.

Vector elements can be set and accessed using the functions

```
void setElement(const int &i, const ComplexType &x)
void setElementRange(const int& lo, const int &hi, ComplexType *x)
void setElements(const int &n, const int *i, const ComplexType *x)
void addElement(const int &i, const ComplexType &x)
void addElements(const int& n, const int *i, const ComplexType *x)
void getElement(const int& i, ComplexType& x) const
void getElements(const int& n, const int *i, ComplexType *x) const
void getElementRange(const int& lo, const int& hi,
                    ComplexType *x) const
void ready(void)
```

These functions all operate in a similar way to the corresponding matrix operations. The **setElementRange** function, etc. are similar to the **setElements** function except that instead of specifying individual element indices in a separate vector, the low and high indices of the segment to which the values are assigned is specified (this assumes that the values in the array **x** represent a contiguous segment of the vector). Again, for real vectors, all values of type **ComplexType** should be replaced by values of type double. The utility functions

```
Vector *clone(void) const
void print(const char* filename = NULL) const
void save(const char *filename) const
```

also have similar behaviors to their matrix counterparts.

Additional operations that can be performed on the entire vector include

```
void zero(void)
void equate(const Vector &x)
void fill(const ComplexType& v)
ComplexType norm1(void) const
ComplexType norm2(void) const
ComplexType normInfinity(void) const
void scale(const ComplexType& x)
void add(const ComplexType& x)
void add(const Vector& x, const ComplexType& scale = 1.0)
void elementMultiply(const Vector& x)
```

**void elementDivide(const Vector& x)**

The **zero** function sets all vector elements to zero, the **equate** function copies all values of the vector **x** to the corresponding elements of the calling vector, **fill** sets all elements to the value **v**, **norm1** returns the  $L_1$  norm of the vector, **norm2** returns the  $L_2$  norm and **normInfinity** returns the  $L_\infty$  norm. The **scale** function can be used to multiply all vector elements by the value **x**, the first **add** function can be used to add the constant **x** to all vector elements and the second **add** function can be used to add the vector **x** to the calling vector after first multiplying it by the value **scale**. The final two functions multiply or divide each element of the calling vector by the value in the vector **x**.

The following methods modify the values of the vector elements using some function of the element value.

**void abs(void)**  
**void real(void)**  
**void imaginary(void)**  
**void conjugate(void)**  
**void exp(void)**  
**void reciprocal(void)**

The function **abs** replaces each element with its complex norm (absolute value), **real** and **imaginary** replace the elements with their real or imaginary values, **conjugate** replaces the vector elements with their conjugate values, **exp** replaces each vector element with the exponential of its original value and **reciprocal** replaces each element by its reciprocal. The **real**, **imaginary** and **conjugate** functions only apply to complex vectors.

### *Linear Solvers*

The math module also contains solvers. The **LinearSolver** class contains a constructor

**LinearSolver(const Matrix &A)**

that creates an instance of the solver. The matrix **A** defines the set of linear equations **Ax=b** that must be solved. If matrix **A** is a **RealMatrix** then the corresponding class and its constructor is

**RealLinearSolver(const RealMatrix &A)**

The properties of the solver can be modified by calling the function

**void configure(utility::Configuration::Cursor \*props)**

The **Configuration** module is described in more detail below. This function can be used to pass information from the input file to the solver to alter its properties. For the PETSc library, the solver algorithm can be controlled using PETSc's runtime options database. Different options can be passed to PETSc by including a block in the input deck (there is more documentation on input decks in the section on the **Configuration** module). An example of this type of input is

```
<LinearSolver>
  <PETScOptions>
    -ksp_view
    -ksp_type richardson
    -pc_type lu
    -pc_factor_mat_solver_package superlu_dist
    -ksp_max_it 1
  </PETScOptions>
</LinearSolver>
```

The **LinearSolver** block is where different solver parameters are defined and the **PETScOptions** block is where a string can be passed to the runtime options database. Additional parameters that can be passed to the solver include **SolutionTolerance**, **MaxIterations** and **FunctionTolerance**. Some solvers that are available in PETSc only run serially and will fail if run on more than one processor. However, for the problem size ranges frequently encountered in power grid analysis, the serial solvers may be the fastest options. Other parts of the code may be more scalable so it is desirable to run them in parallel. GridPACK has options that allow users to run the code in parallel while using a serial solver, without the need to modify any application source code. This can be done by including the options

```
<ForceSerial>true</ForceSerial>
<InitialGuessZero>true</InitialGuessZero>
<SerialMatrixConstant>true</SerialMatrixConstant>
```

in the **LinearSolver** block. The first option can be used to replicate the linear solver across all processors in the system and then distribute the answer to processors. The second option eliminates the need for obtaining an initial guess for the solution from all processors and provides additional performance gains. The final option can be used if the matrix does not change between function calls. Only new versions of the RHS vector need to be replicated on each processor after the first call. This can also result in performance gains.

After configuring the solver, it can be used to solve the set of linear equations by calling the method

```
void solve(const Vector &b, Vector &x) const
```

This function returns the solution  $\mathbf{x}$  based on the right hand side vector  $\mathbf{b}$ .

### *Non-linear Solvers*

The math module also supports non-linear solvers for systems of the type  $\mathbf{A}(\mathbf{x}) \cdot \mathbf{x} = \mathbf{b}(\mathbf{x})$  but the interface is more complicated than for the linear solvers. In order for the non-linear solver to work, two functions must be defined by the user. The first evaluates the Jacobian of the system for a given trial state  $\mathbf{x}$  of the system and the second computes the right hand side vector for a given trial state  $\mathbf{x}$ . The two functions are of type **JacobianBuilder** and **FunctionBuilder**. The **JacobianBuilder** function is a function with arguments

```
(const math::Vector &vec, math::Matrix &jacobian)
```

and **FunctionBuilder** is a function with arguments

```
(const math::Vector &xCurrent, math::Vector &newRHS)
```

These functions need to be added to the system somewhere. They can then be assigned to objects of type **JacobianBuilder** and **FunctionBuilder** and passed to the constructor of the non-linear solver. There are a number of ways to do this. In the following discussion, we will adopt the method used in the non-linear solver version of the power flow code that is distributed with GridPACK.

The first step is to define a struct that can be used to implement the functions needed by the non-linear solver (the actual implementation contains additional declarations and code, but the important features of this helper class are outlined here)

```
struct SolverHelper : private utility::Uncopyable
{
    //Constructor
    SolverHelper(/* Arguments to initialize helper */)
    {
        // Initialize non-linear calculation
    }

    :
    boost::shared_ptr<math::Matrix> matrix; // Jacobian matrix
    boost::shared_ptr<math::Vector> X; // Current state
    :
    void operator() (const math::Vector &xCurrent, math::vector &newRHS)
    {
        // Evaluate RHS vector from current state xCurrent
    }
    void operator() (const math::Vector &xCurrent,
```



```

        math::Matrix &Jacobian)
    {
        // Evaluate Jacobian from current state xCurrent
    }
}

```

The important functions for this discussion are the overloaded **operator()** functions. In the application code, this helper struct can be initialized and used to create two functions of type **JacobianBuilder** and **FunctionBuilder** using the syntax

```

SolverHelper helper(//Arguments to initialize helper //);
math::JacobianBuilder jbuild = boost::ref(helper);
math::FunctionBuilder fbuild = boost::ref(helper);

```

At this point **jbuild** and **fbuild** are pointing to the overloaded functions in **helper** that have the appropriate arguments for a function of type **JacobianBuilder** and type **FunctionBuilder**. The **boost::ref** command provides a reference to the appropriate function in **helper** instead of making a copy, this preserves any state that might be present in **helper** between invocations of the functions **jbuild** and **fbuild** by the solver.

For the power flow application using a non-linear solver, the creation of the solver is a two-step process. First, a pointer to a non-linear solver interface is created and then a particular solver instance is assigned to this interface. The power flow application can point to a hand-coded Newton-Raphson solver or a wrapper to the PETSc library of solvers. The code for this is the following

```

boost::scoped_ptr<math::NonlinearSolverInterface> solver;
if (useNewton) {
    math::NewtonRaphsonSolver *tmpsolver =
        new math::NewtonRaphsonSolver(*(helper.matrix), jbuild, fbuild);
    solver.reset(tmpsolver);
} else {
    solver.reset(new math::NonlinearSolver(*(helper.matrix), jbuild,
fbuild));
}

```

If you are only interested in using the **NonlinearSolver**, then it is possible to dispense with the **NonlinearSolverInterface** and just use the **NonlinearSolver** directly. The remaining call to invoke the solver is just

```

solver->solver(*helper.X);

```

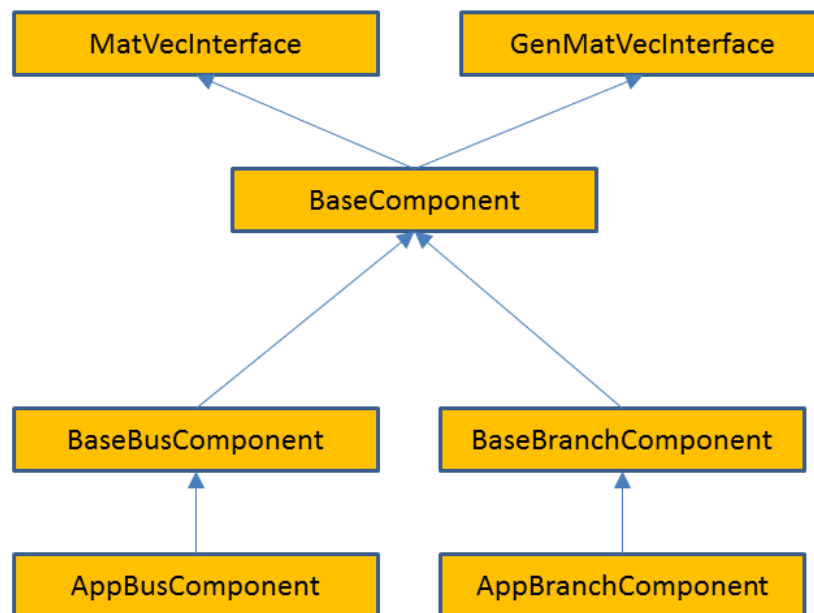
Additional calls are likely to be added to these to allow user-specified parameters from the input deck to be sent to the solver. In the case of the **NonlinearSolver**, these can be used to specify which PETSc solver should be used.

More details on how to use the non-linear solvers can be found by looking at the powerflow module in the GridPACK source code.

### Network Components

Network component is a generic term for objects representing buses and branches. These objects determine the behavior of the system and the type of analyses being done. Branch components can represent transmission lines and transformers while bus components could model loads, generators, or something else. Both kinds of components could represent measurements (e.g. for a state estimation analysis).

Network components cover a fairly broad range of behaviors and there is little that can be said about them outside the context of a specific problem. Each component inherits from a matrix-vector interface, which enables the framework to generate matrices and vectors from the network in a relatively straightforward way. In addition, buses inherit from a base bus interface and branches inherit from a base branch interface. The relationship between these interfaces is shown in Figure 5.



**Figure 5.** Schematic diagram showing the interface hierarchy for network components.

These base interfaces provide mechanisms for accessing the neighbors of a bus or branch and allow developers to specify what data is transferred in ghost exchanges. They do not define any physical properties of the bus or branch, it is up to application developers to do this.

Of these interfaces, the matrix-vector interfaces are the most important. The **MatVecInterface** is used for most calculations that directly model the physics of the power grid and described problems where the dependent and independent variables are associated with buses. The **GenMatVecInterface** is used for problems where variables are also associated with branches, such as state estimation or Kalman filter calculations. This section will describe the **MatVecInterface**, the **GenMatVecInterface** is described in more detail later in this document. The **MatVecInterface** is designed to answer the question of what block of data is contributed by a bus or branch to a matrix or vector and what the dimensions of the block are. For example, in constructing the Y-matrix for a power flow problem using a real-valued formulation, the grid components representing buses contribute a  $2 \times 2$  block to the diagonal of the matrix. Similarly, the grid components representing branches contribute a  $2 \times 2$  block to the off-diagonal elements. (Note that if the Y-matrix is expressed as a complex matrix, then the blocks are of size  $1 \times 1$ .) The location of these blocks in the matrix is determined by the location of the corresponding buses and branches in the network, but the indexing calculations required to determine how this location maps to a location in the matrix can be made completely transparent to the user via the mapper module.

Because the matrix-vector interface focuses on small blocks, it is relatively easy for power grid engineers to write the corresponding methods. The full matrices and vectors can then be generated from the network using simple calls to the mapper interface (see the discussion below on the mapper module). All of the base network component classes reside in the **gridpack::component** namespace.

The primary function of the **MatVecInterface** class is to enable developers to build the matrices and vectors used in the solution algorithms for the network. It eliminates a large number of tedious and error-prone index calculations that would otherwise need to be performed in order to determine where in a matrix a particular data element should be placed. The **MatVecInterface** includes basic constructors and destructors. The first set of non-trivial operations are implemented on buses and set the values of diagonal blocks in the matrix. Additional functions are implemented on branches and set values for off-diagonal elements. Vectors can be created by calling functions defined on buses. These functions are described in detail below.

The functions that are used to create diagonal matrix blocks are

```
virtual bool matrixDiagSize(int *isize, int *jsize) const
```

```
virtual bool matrixDiagValues(ComplexType *values)
```

**virtual bool matrixDiagValues(RealType \*values)**

These functions are virtual functions and are expected to be overwritten by application-specific bus and branch classes. Depending on whether the application should create real or complex matrices, either the real or complex versions of **matrixDiagValues** can be implemented. The default behavior is to return 0 for **isize** and **jsize** for **matrixDiagSize** and to return false for all functions. These functions will not build a matrix unless overwritten by the application. Not all functions need to be overwritten by a given bus or branch class. Generally, only a subset of functions may be needed by an application.

The **matrixDiagSize** function returns the size of the matrix block that is contributed by the bus to a matrix. If a single number is contributed by the bus, the **matrixDiagSize** function returns 1 for both **isize** and **jsize**. Similarly, for a 2×2 block then both **isize** and **jsize** are set to 2. The return value is true if the bus contributes to the matrix, otherwise it is false. Returning false can occur, for example, if the bus is the reference bus in a power flow calculation. For a more complicated calculation, such as a dynamic simulation with multiple generators on some buses, the size of the matrix blocks can differ from bus to bus. Note that the values returned by **matrixDiagSize** refer only to the particular bus on which the function is invoked. It does not say anything about other buses in the system.

The **matrixDiagValues** function returns the actual values for the matrix block associated with the bus for which the function is invoked. The values are returned as a linear array with values returned in column-major order. For a 2×2 block, this means the first value is at the (0,0) position, the second value is at the (1,0) position, the third value is at the (0,1) position and the fourth value is at the (1,1) position. This function also returns true if the bus contributes to the matrix and false otherwise. This may seem redundant, since the **matrixDiagSize** function has already returned this information but it turns out there are certain applications where it is desirable for the **matrixDiagSize** function to return true and the **matrixDiagValues** function to return false. The buffer **values** is supplied by the calling program and is expected to be big enough, based on the dimensions returned by the **matrixDiagSize** function, to contain all returned values.

The functions that are used to return values for off-diagonal matrix elements are listed below. These are usually only implemented for branches.

**virtual bool matrixForwardSize(int \*isize, int \*jsize) const**

**virtual bool matrixForwardValues(ComplexType \*values)**

**virtual bool matrixReverseSize(int \*isize, int \*jsize) const**

**virtual bool matrixReverseValues(ComplexType \*values)**

Only the complex versions of these functions are listed but equivalent functions for real matrices are available. These functions work in a similar way to the functions for creating blocks along the diagonal, except that they split off-diagonal matrix calculations into forward elements and reverse elements. The initial approximate location of an off-diagonal matrix element in a matrix is based in some internal indices assigned to the buses at either end of the branch. Suppose that these indices are **i**, corresponding to the “from” bus and **j**, corresponding to the “to” bus. The “forward” functions assume that the request is for the **ij** element while the “reverse” functions assume that the request is for the **ji** element. Another way of looking at this is the following: as discussed below, branches contain pointers to two buses. The first is the “from” bus and the second is the “to” bus. The forward functions assume that the “from” bus corresponds to the first index of the element, the reverse functions assume that the “from” bus corresponds to the second index of the element. Note that if a bus does not contribute to a matrix, then the branches that are connected to the bus should also not contribute to the matrix.

The final set of functions in the **MatVecInterface** that are of interest to application developers are designed to set up vectors. These are usually implemented only for buses. These functions are analogous to the functions for creating matrix elements

**virtual bool vectorSize(int \*isize) const**

**virtual bool vectorValues(ComplexType \*values)**

The **vectorSize** function returns the number of elements contributed to the vector by a bus and the **vectorValues** returns the corresponding values. The **vectorValues** function expects the buffer values to be allocated by the calling program. In addition to functions that can be used to specify a vector, there is an additional function that can be used to push values from a vector back onto a bus. This function is

**virtual void setValues(ComplexType \*values)**

The buffer contains values from the vector corresponding to internal variables in the bus and this function can be used to set the bus variables. The **setValues** function could be used to assign bus variables so that they can be used to recalculate matrices and vectors for an iterative loop in a non-linear solver or so that the results of a calculation can be exported to an output file. Real versions of the **vectorValues** and **setValues** functions are available for real vectors.

The **BaseComponent** class contains additional functions that contribute to the base properties of a bus or branch. Again, most of the functions in this class are virtual and are expected to be overwritten by actual implementations. However, not all of them need to be overwritten by a

particular bus or branch class. Many of these functions are used in conjunction with the **BaseFactory** class, which defines methods that run over all buses and branches in the network and invokes the functions defined below.

The **load** function

```
virtual void load(const boost::shared_ptr<DataCollection> &data)
```

is used to instantiate components based on data in the network configuration file that is used to create the network. It is used in conjunction with the **DataCollection** object, which is described in more detail below. Networks are generally created by first instantiating a network parser. The parser is used to read in an external network file and create the network topology. The next step is to invoke the partition function on the network to get all network elements properly distributed between processors. At this point, the network, including ghost buses and branches, is complete and each bus and branch has a **DataCollection** object containing all the data in the network configuration file that pertains to that particular bus or branch. The data in the **DataCollection** object is stored as simple key-value pairs. This data is used to initialize the corresponding bus or branch by invoking the load function on all buses and branches in the system. The bus and branch classes must implement the **load** function to extract the correct parameters from the **DataCollection** object and use them to assign internal component parameters.

Only one type of bus and one type of branch is associated with each network but many different types of equations can be generated by the network. To allow developers to embed many different behaviors into a single network and to control at what points in the simulation those behaviors can be manifested, the concept of modes is used. The function

```
virtual void setMode(int mode)
```

can be used to set an internal variable in the component that tells it how to behave. The variable “**mode**” usually corresponds to an enumerated constant that is part of the application definition. For example, in a power flow calculation it might be necessary to calculate both the Y-matrix and the equations for the power flow solution containing the Jacobian matrix and the right-hand side vector. To control which matrix gets created, two modes are defined: “**YBus**” and “**Jacobian**”. Inside the matrix functions in the **MatVecInterface**, there is a condition

```
if (p_mode == YBus) {  
    // Return values for Y-matrix calculation  
} else if (p_mode == Jacobian) {  
    // Return values for power flow calculation  
}
```

The variable “**p\_mode**” is an internal variable in the bus or branch that is set using the **setMode** function.

The function

```
virtual bool serialWrite(char *string, const int bufsize,  
                        const char *signal = NULL)
```

is used in the serial IO modules described below to write out properties of buses or branches to standard output. The character buffer “**string**” contains a formatted line of text representing the properties of the bus or branch that is written to standard output, the variable “**bufsize**” gives the number of characters that “**string**” can hold, and the variable “**signal**” can be used to control what data is written out. The return value is true if the bus or branch is writing out data and false otherwise. For example, if the application is writing out the properties of all buses with generators, then the signal “**generator**” might be passed to this subroutine. If a bus has generators, then a string is copied into the buffer “**string**” and the function returns true, otherwise it returns false. The buffer “**string**” is allocated by the calling program. The variable “**bufsize**” is provided so that the bus or branch can determine if it is overwriting the buffer. Returning to the generator example, if this call returns a separate line for each generator, then it is possible that a bus with too many generators might exceed the buffer size. This could be detected by the implementation if the buffer size is known. More information on how this function is used can be found in the discussion of the serial IO modules.

The **BaseComponent** class also contains two functions that must be implemented if buses and/or branches need to exchange data with other processors. Data that must be exchanged needs to be placed in buffers that have been allocated by the network. The bus and branch objects specify how large the buffers need to be by implementing the function

```
virtual int getXCBufSize()
```

This function must return the same value for all buses and all branches in the same bus or branch classes. Buses can return a different value than branches. For example, in a power flow calculation, it is necessary that ghost buses get new values of the phase angle and voltage magnitude increments. These are both real numbers so the **getXCBusSize** routine needs to return the value **2\*sizeof(double)**. Note that all buses must return this value even if the bus is a reference bus and does not participate in the calculation.

This function is queried by the network and used to allocate a buffer of the appropriate size. The network then informs the bus and branch objects where the location of the buffer is by invoking the function

```
virtual void setXCBuf(void *buf)
```

The bus or branch can use this function to set internal pointers to this buffer that can be used to assign values to the buffer (which is done before a ghost exchange) or to collect values from the buffer (which is done after a ghost exchange). Continuing with the powerflow example, the bus implementation of the **setXCBuf** function would look like

```
setXCBuf(void *buf)
{
    p_Ang_ptr = static_cast<double*>(buf);
    p_Mag_ptr = p_Ang_ptr+1;
}
```

The pointers **p\_Ang\_ptr** and **p\_Mag\_ptr** of type **double** are internal variables of the bus implementation and can be used elsewhere in the bus whenever the voltage angle and voltage magnitude variables are needed. After a network update operation, ghost buses will contain values for these variables that were calculated on the home processor that owns the corresponding bus.

The **BaseBusComponent** and **BaseBranchComponent** classes contain a few additional functions that are specific to whether or not a component is a bus or a branch. The **BaseBusComponent** class contains functions that can be used to identify attached buses or branches, determine if the bus is a reference bus, and recover the original indices of the bus. Other functions are included in the **BaseBusClass** but these are not usually required by application developers and are used primarily to implement other GridPACK functions.

To get a list of pointers to all branches connected to a bus, the function

```
void getNeighborBranches(
    std::vector<boost::shared_ptr<BaseComponent> > &nghbrs) const
```

can be called. This provides a list of pointers to all branches that have the calling bus as one of its endpoints. This function can be used inside a bus method to loop over attached branches, which is a common motif in matrix calculations. For example, to evaluate the contribution to a diagonal element of the Y-matrix coming from transmission lines, it is necessary to perform the sum

$$Y_{ii} = -\sum_{j \neq i} Y_{ij}$$

where the  $Y_{ij}$  are the contribution due to transmission lines from the branch connecting i and j. The code inside a bus component that evaluates this sum can be written as

```
std::vector<boost::shared_ptr<BaseComponent> > branches;
getNeighborBranches(branches);
ComplexType y_diag(0.0,0.0);
```



```

for (int i=0; i<branches.size(); i++) {
    YBranch *branch = dynamic_cast<YBranch*>(branches[i].get());
    y_diag += branch->getYContribution();
}

```

The function **getYContribution** evaluates the quantity  $Y_{ij}$  using parameters that are local to the branch. The return value is then accumulated into the bus variable **y\_diag**, which is eventually returned through the **matrixDiagValues** function. The **dynamic\_cast** is necessary to convert the pointer from a **BaseComponent** object to the application class **YBranch**. The **BaseComponent** class has no knowledge of the **getYContribution** function, this is only implemented in the class **YBranch**.

A function that is similar to **getNeighborBranches** is

```

void getNeighborBuses(
    std::vector<boost::shared_ptr<BaseComponent> > &nghbrs) const

```

which can be used to get a list of the buses that are connected to the calling bus via a single branch.

Many power grid problems require the specification of a special bus as a reference bus. This designation can be handled by the two functions

```

void setReferenceBus(bool status)

```

```

bool getReferenceBus() const

```

The first function can be used (if called with the argument true) to designate a bus as the reference bus and the second function can be called to inquire whether a bus is the reference bus. A reference bus is usually set when the network configuration file is read in and does not need to be set explicitly by the application.

Finally, it is often useful for exporting results if the original index of the bus is available. This can be recovered using the function

```

int getOriginalIndex() const

```

The functions in the **BaseBusComponent** class only work correctly after a call to the base factory method **setComponents**, which is described below. Other functions in the **BaseBusComponent** class are needed within the framework but are not usually required by application developers.

The **BaseBranchComponent** class is similar to the **BaseBusComponent** class and provides basic information about branches and the buses at either end of the branch. To retrieve pointers to the buses at the ends of the branch, the following two functions are available

```
boost::shared_ptr<BaseComponent> getBus1() const
```

```
boost::shared_ptr<BaseComponent> getBus2() const
```

The **getBus1** function returns a pointer to the “from” bus, the **getBus2** function returns a pointer to the “to” bus.

Two other functions in the **BaseBranchComponent** class that are useful for writing output are

```
int getBus1OriginalIndex() const
```

```
int getBus2OriginalIndex() const
```

These functions get the original index of “from” and “to” buses. Unlike buses, the branches are not characterized by a single index. Similar to the functions in the **BaseBusComponent** class, the functions in the **BaseBranchComponent** class will not work correctly until the **setComponents** method has been called in the base factory class.

Finally, a separate network component class that is associated with all buses and branches (including ghost buses and branches) is the **DataCollection** class. This class is a simple container that can be used to store key-value pairs. It also resides in the **gridpack::component** namespace. When the network is created using a standard parser to read a network configuration file (see more on parsers below), each bus and branch in the network, including the ghosts, has an associated **DataCollection** object that contains all parameters from the configuration file that are associated with that particular bus or branch. These can be retrieved from the **DataCollection** object using some simple accessors. Data can be stored in two ways inside the **DataCollection** object. The first method assumes that there is only a single instance of the key-value pair, the second assumes there are multiple instances. This second case can occur, for example, if there are multiple generators on a bus. Generators are characterized by a collection of parameters and each generator has its own set of parameters. The generator parameters can be indexed so that they can be matched with a specific generator.

Assuming that a parameter only appears once in the data collection, the contents of a **DataCollection** object can be accessed using the functions

```
bool getValue(const char *name, int *value)  
bool getValue(const char *name, long *value)
```

```

bool getValue(const char *name, bool *value)
bool getValue(const char *name, std::string *value)
bool getValue(const char *name, float *value)
bool getValue(const char *name, double *value)
bool getValue(const char *name, ComplexType *value)

```

These functions return true if a variable of the correct type is stored in the **DataCollection** object with the key “**name**”, otherwise it returns false.

If the variable is stored multiple times in the **DataCollection**, then it can be accessed with the functions

```

bool getValue(const char *name, int *value, const int idx)
bool getValue(const char *name, long *value, const int idx)
bool getValue(const char *name, bool *value, const int idx)
bool getValue(const char *name, std::string *value, const int idx)
bool getValue(const char *name, float *value, const int idx)
bool getValue(const char *name, double *value, const int idx)
bool getValue(const char *name, ComplexType *value, const int idx)

```

where **idx** is an index that identifies a particular instance of the key. These functions are used primarily to implement the network component **load** method, described above for the **BaseComponent** class.

The data collection is generally filled with values after the parser is called to create the network. The nomenclature for these values can be found in the **dictionary.hpp** file located under **src/parser** under the main GridPACK directory.

The **DataCollection** objects can also be used to transfer data between different networks. This is important for chaining different types of calculations together. For example, a powerflow or state estimation calculation might be used to initialize a dynamic simulation and the **DataCollection** object can be used as a mechanism for transferring data between the two different networks. Because of this, the functions for adding more data to the **DataCollection** and the functions for overwriting the values of existing data are useful. New key value pairs can be added to a data collection object using the functions

```

void addValue(const char *name, int value)
void addValue(const char *name, long value)
void addValue(const char *name, bool value)
void addValue(const char *name, char *value)
void addValue(const char *name, float value)
void addValue(const char *name, double value)

```

```

void addValue(const char *name, ComplexType value)

void addValue(const char *name, int value, const int idx)
void addValue(const char *name, long value, const int idx)
void addValue(const char *name, bool value, const int idx)
void addValue(const char *name, char *value, const int idx)
void addValue(const char *name, float value, const int idx)
void addValue(const char *name, double value, const int idx)
void addValue(const char *name, ComplexType value, const int idx)

```

Existing values can be overwritten with the functions

```

bool setValue(const char *name, int value)
bool setValue(const char *name, long value)
bool setValue(const char *name, bool value)
bool setValue(const char *name, char *value)
bool setValue(const char *name, float value)
bool setValue(const char *name, double value)
bool setValue(const char *name, ComplexType value)

bool setValue(const char *name, int value, const int idx)
bool setValue(const char *name, long value, const int idx)
bool setValue(const char *name, bool value, const int idx)
bool setValue(const char *name, char *value, const int idx)
bool setValue(const char *name, float value, const int idx)
bool setValue(const char *name, double value, const int idx)
bool setValue(const char *name, ComplexType value, const int idx)

```

The functions return false if the value is not already in the **DataCollection** object.

## Factories

The network component factory is an application-dependent piece of software that is designed to manage interactions between the network and the network component objects. Most operations in the factory run over all buses and all branches and invoke some operation on each component. An example is the “**load**” operation. After the network is read in from an external file, it consists of a topology and a set of simple data collection objects containing key-value pairs associated with each bus and branch. The **load** operation then runs over all buses and branches and instantiates the appropriate objects by invoking a local **load** method that takes the values from the data collection object and uses it to instantiate the bus or branch. The application network factory is derived from a base network factory class that contains some additional routines that set up indices, assign neighbors to individual buses and branches and assign buffers.

The neighbors are originally only known to the network, so a separate operation is needed to push this information down into the bus and branch components. The network component factory may also execute other routines that contribute to setting up the network and creating a well-defined state.

Factories can be derived from the **BaseFactory** class, which is a templated class that is based on the network type. It resides in the **gridpack::factory** namespace. The constructor for a **BaseFactory** object has the form

```
BaseFactory<MyNetwork>(boost::shared_ptr<MyNetwork> network)
```

The **BaseFactory** class supplies some basic functions that can be used to help instantiate the components in a network. Other methods can be added for particular applications by inheriting from the **BaseFactory** class. The two most important functions in **BaseFactory** are

```
virtual void setComponents()
```

```
virtual void setExchange()
```

The **setComponents** method pushes topology information available from the network into the individual buses and branches using methods in the base component classes. This operation ensures that operations such as **getNeighborBuses**, etc. work correctly. The topology information is originally only available in the network and it requires additional operations to imbed it in individual buses and branches. Being able to access this information directly from the buses and branches can simplify application programming substantially.

The **setExchange** function allocates buffers and sets up pointers in the components so that exchange of data between buses and branches can occur and ghost buses and branches can receive updated values of the exchanged parameters. This function loops over the **getXCBusSize** and **setXCBuf** commands defined in the network component classes and guarantees that buffers are properly allocated and exposed to the network components.

Two other functions are defined in the **BaseFactory** class as convenience functions. The first is

```
virtual void load()
```

This function loops over all buses and branches and invokes the individual bus and branch **load** methods. This moves information from the **DataCollection** objects that are instantiated when the network is created from a network configuration file to the bus and branch objects themselves. The second convenience function is

```
virtual void setMode(int mode)
```

This function loops over all buses and branches in the network and invokes each bus and branch **setMode** method. It can be used to set the behavior of the entire network in single function call.

Some utility functions in the **BaseFactory** class that are occasionally useful are

**bool checkTrue(bool flag)**

**bool checkTrueSomewhere(bool flag)**

The **checkTrue** function returns true if the variable **flag** is true on all processors, otherwise it returns false. This function can be used to check if a condition has been violated somewhere in the network. The **checkTrueSomewhere** function returns true if **flag** is true on at least one processor. This function can be used to check if a condition is true anywhere in the system.

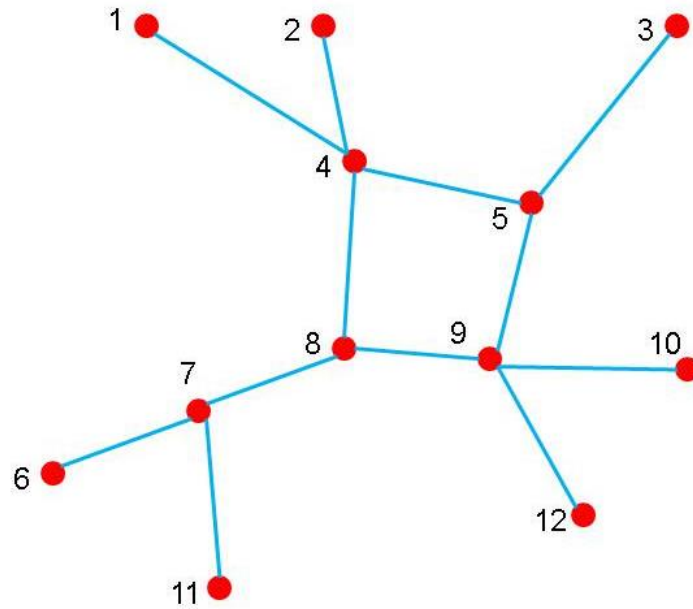
### Mapper Module

The mappers are a collection of generic capabilities that can be used to generate a matrices or vectors from the network components. This is done by running over all the network components and invoking methods in the matrix-vector interface. The mapper is basically a transformation that converts a set of network components into a matrix or vector based on the behavior of their matrix-vector interfaces. It has no explicit dependencies on either the network components or the type of analyses being performed so this capability is applicable across a wide range of problems. At present there are three types of mapper, the standard mapper described here that is implemented on top of the **MatVecInterface**, a more generalized mapper that utilizes the **GenMatVecInterface** and a mapper for generating matrices resembling “fat” vectors. These are dense matrices that are basically a collection of column vectors. The generalized mapper and its corresponding interface are described in a later section below, along with the mapper for generating fat vectors. The mapper discussed in this section is used for problems where both dependent and independent variables are associated with buses, which is the case for problems such as power flow calculations and dynamic simulation. Other problems, such as state estimation, have variables associated with both buses and branches and require the more general interface.

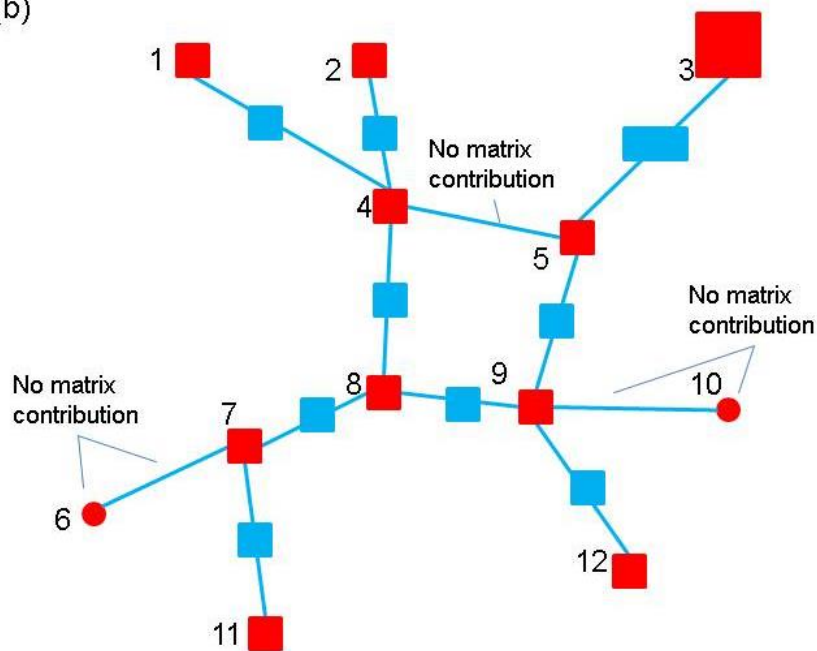
The basic matrix-vector interface contains functions that provide two pieces of information about each network component. The first is the size of the matrix block that is contributed by the component and the second is the values in that block. Using this information, the mapper can calculate the dimensions of the matrix and where individual elements in the matrix are located. The construction of a matrix by the mapper is illustrated in Figure 6 for a small network. Figure 6(a) shows a hypothetical network. The contributions from each network component are shown in Figure 6(b). Note that some buses and branches do not contribute to the matrix. This could occur in real systems because the transmission line corresponding to the branch has failed or because a bus represents the reference bus. In addition, it is possible that buses and branches can contribute different size blocks to the matrix. The mapping of the individual contributions from

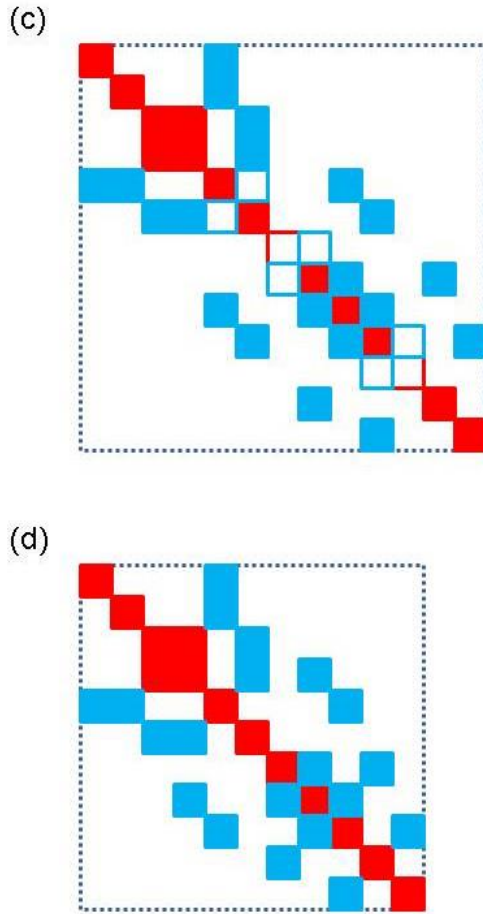
the network in Figure 6(b) to initial matrix locations is shown in Figure 6(c). This is followed by elimination of gaps in the matrix in Figure 6(d).

(a)



(b)





**Figure 6.** A schematic diagram of the matrix map function. The bus numbers in (a) and (b) map to approximate row and column locations in (c). (a) a small network (b) matrix blocks associated with branches and buses. Not that not all blocks are the same size and not all buses and branches contribute (c) initial construction of matrix based on network indices (d) final matrix after eliminating gaps

The most complex part of generating matrices and vectors is implementing the functions in the **MatVecInterface**. Once this has been done, actually creating matrices and vectors using the mappers is quite simple. The **MatVecInterface** is associated with two mappers, one that creates matrices from buses and branches and a second that can create vectors from buses. Both mappers are templated objects based on the type of network being used and use the **gridpack::mapper** namespace. The **FullMatrixMap** object runs over both buses and branches to set up a matrix. The constructor is

```
FullMatrixMap<MyNetwork>(boost::shared_ptr<MyNetwork> network)
```

The network is passed in to the object via the constructor. The constructor sets up a number of internal data structures based on what mode has been set in the network components. For



example, for a power flow application where it might be necessary to create both a Y-matrix and a Jacobian matrix, it would be necessary to create two mappers. If the first mapper is created while the mode is set to construct the Y-matrix, then it will be necessary to instantiate a second mapper to create the Jacobian for a power flow calculation. Before instantiating the second matrix, the mode should be set to Jacobian.

Once the mapper has been created, a matrix can be generated using the call

```
boost::shared_ptr<gridpack::math::Matrix> mapToMatrix()
```

This function creates a new matrix and returns a pointer to it. If a matrix already exists and it is only necessary to update the values, then the functions

```
void mapToMatrix(  
    boost::shared_ptr<gridpack::math::Matrix> &matrix)
```

```
void mapToMatrix(gridpack::math::Matrix &matrix)
```

can be used. These functions use the existing matrix data structures and overwrite the values of individual elements. For these to work, it is necessary to use the same mapper that was used to create the original matrix and to have the same mode set in the network components.

Additional operations that can be used on existing matrices include

```
void overwriteMatrix(boost::shared_ptr<gridpack::math::Matrix> matrix)  
void overwriteMatrix(gridpack::math::Matrix &matrix)  
void incrementMatrix(boost::shared_ptr<gridpack::math::Matrix> matrix)  
void incrementMatrix(gridpack::math::Matrix &matrix)
```

These operations are designed to support making small changes in an existing matrix instead of reconstructing the full matrix from scratch. This can happen in contingency calculations or simulations of faults where a single grid element goes out or changes value. Instead of rebuilding the entire matrix, it is possible to modify only a small portion if it. To use these functions, it is necessary to define at least two modes in the network components. The first mode is used to build the original matrix, the second is used to make changes. All **MatVecInterface** functions that return true using the second mode (the one making changes) must return true for the first mode (used to build the original matrix). Furthermore, all block sizes for the second mode must match the block sizes in the first mode. The **overwriteMatrix** functions replace the values in the matrix with the values returned by the **MatVecInterface** functions, the **incrementMatrix** functions add these values to whatever is already in the matrix.

The vector mapper works in an entirely analogous way to the matrix mapper. The constructor for the **BusVectorMap** class is

```
BusVectorMap<MyNetwork>(boost::shared_ptr<MyNetwork> network)
```

and the function for building a new vector is

```
boost::share_ptr<gridpack::math::Vector> mapToVector()
```

The functions for overwriting the values of an existing vector are

```
void mapToVector(  
    boost::shared_ptr<gridpack::math::Vector> &vector)
```

```
void mapToVector(gridpack::math::Vector &vector)
```

The vector map can also be used to write values back to buses using the function

```
void mapToBus(const gridpack::math::Vector &vector)
```

This function will copy values from the vector into the bus using the **setValues** function in the **MatVecInterface**.

### Parser Module

The parser modules are designed to read an external network file, set up the network topology and assign any parameter fields in the file to simple fields. The parsers do not partition the network, they are only responsible for reading in the network and distributing the different network elements in a way that guarantees that not too much data ends up on any one processor. The parsers are also not responsible for determining if the input is compatible with the analysis being performed. This can be handled, if desired, by building checks into the network factory. The parsers are only responsible for determining if they can read the file.

Currently, GridPACK only supports two file formats. Files based on the PSS/E PTI version 23 and version 33 formats can be read in using the classes **PTI23\_parser** and **PTI33\_parser**. Both parsers can also read PSS/E formatted .dyr files that are used to read in extra parameters used in dynamic simulation. The parsers are templated classes that again use the network type as a template argument. Both **PTI23\_parser** and **PTI33\_parser** are located in the **gridpack::parser** namespace. These classes have only a few important functions. The first are the constructors

```
PTI23_parser<MyNetwork>(boost::shared_ptr<MyNetwork> network)
```

```
PTI33_parser<MyNetwork>(boost::shared_ptr<MyNetwork> network)
```

The remaining functions are common to both parsers. To read a PSS/E PTI file containing a network configuration and generate a network, the parser calls the method

```
void parse(const std::string &filename)
```

where filename refers to the location of the network configuration file. To use this parser, the network object with the appropriate bus and branch classes is instantiated and then passed to the constructor of the **PTI23\_parser** or **PTI33\_parser** object. The parse method is then invoked with the name of the network configuration file passed in as an argument and the network is filled out with buses and branches. The parameters in the network configuration file are stored as key-value pairs in the **DataCollection** object associated with each bus and branch. Once the network partition method has been called, the network is fully distributed and ghost buses and branches have been created. Other operations can then be performed. A variant on parse is the command

```
void externalParse(const std::string &filename)
```

This command can be used to parse .dyr files containing dynamic simulation parameters. The difference between this function and **parse** is that **externalParse** assumes that the network already exists and that the parameters that are read in will be added to it. This command should therefore only be called after a network has been created using **parse**.

Another key part of the parsing capability is the **dictionary.hpp** file, which is designed to provide a common nomenclature for parameters associated with power grid components. This file is located in the **src/parser** directory under the main GridPACK directory. The definitions in the **dictionary.hpp** are used to extract parameters from the **DataCollection** objects created by the parser. For example, the parameter describing the resistance of a transmission element is given the common name **BRANCH\_R**. This string is defined as a macro in the **dictionary.hpp** file as

```
#define BRANCH_R "BRANCH_R"
```

The use of a macro provides compile time error checking on the name. The goal of using the dictionary is that all parsers will eventually store the branch resistance parameter in the **DataCollection** object using this common name. Applications can then switch easily between different network configuration file formats by simply exchanging parsers, which will all store corresponding parameters using a common naming convention that can be used within the code to access data.

### Serial IO Module

The serial IO module is designed to provide a simple mechanism for writing information from selected buses and/or branches to standard output or a file using a consistent ordering scheme. Individual buses and/or branches implement a write method that will write bus/branch information to a single string. This information usually consists of bus or branch identifiers plus some parameters that are desired in the output. The serial IO module then gathers this information, moves it to the head node, and writes it out in a consistent order. An example of this type of output is shown below.

### Bus Voltages and Phase Angles

Bus Number	Phase Angle	Voltage Magnitude
1	0.000000	1.060000
2	-4.982589	1.045000
3	-12.725100	1.010000
4	-10.312901	1.017671
5	-8.773854	1.019514
6	-14.220946	1.070000
7	-13.359627	1.061520
8	-13.359627	1.090000
9	-14.938521	1.055932
10	-15.097288	1.050985
11	-14.790622	1.056907
12	-15.075585	1.055189
13	-15.156276	1.050382
14	-16.033645	1.035530

**Figure 7.** Example output from buses in a 14 bus problem.

Note that the output is ordered by bus index (which matches the ordering of the buses in the original network configuration file). This ordering would be preserved regardless of the number of processors used in the calculation.

Like the mapper, the serial IO classes are relatively easy to use. Most of the complexity is associated with implementing the **serialWrite** methods in the buses and branches. Data can be written out for buses and/or branches using either the **SerialBusIO** class or the **SerialBranchIO** class. These are again templated classes that take the network as an argument in the constructor. Both classes reside in the **gridpack::serial\_io** namespace. The **SerialBusIO** constructor has the form

```
SerialBusIO<MyNetwork>(int max_str_len,  
    boost::shared_ptr<MyNetwork> network)
```

The variable **max\_str\_len** is the length, in bytes, of the maximum size string you would want to write out using this class and **network** is a pointer to the network that is used to generate output. The value of **max\_str\_len** is used to allocate internal memory and also determines how much data needs to be moved around each time data from the entire network is written out. As the value of this parameter increases, the amount of memory needed and the amount of data that needs to move increases, so this value should be kept to a minimum, if possible.

Two additional functions can be used to actually generate output. They are

```
void header(const char *string) const
```

and

```
void write(const char *signal = NULL)
```

The **header** method is a convenience function that will only write the buffer string from the head processor (process 0) and can be used for creating the headings above an output listing. The **write** function traverses all the buses in the network and writes out the strings generated by the **serialWrite** methods in the buses. The **SerialBusIO** object is responsible for reordering these strings in a consistent manner, even if the buses are distributed over many processors. The optional variable “**signal**” is passed to the **serialWrite** methods and can be used to control what output is listed. For example, in one part of a simulation it might be desirable to list the voltage magnitude and phase angle from a powerflow calculation and in another part of the calculation to list the rotor angle for a generator. These two outputs could be distinguished from each other in the **serialWrite** function using the **signal** variable.

To generate the output in Figure 7, the following calls are used

```
gridpack::serial_io::SerialBusIO<MyNetwork> busIO(128,network);  
busIO.header("\n  Bus Voltages and Phase Angles\n");  
busIO.header(  
    "\n  Bus Number      Phase Angle      Voltage Magnitude\n");  
busIO.write();
```

The first call creates the **SerialIOBus** object and specifies the internal buffers size (128 bytes). This buffer must be large enough to incorporate the output from any call to one of the **serialWrite** calls in the bus components. The next two lines print out the header on top of the bus listing and the last line generates the listing itself. The **serialWrite** implementation looks like

```
bool gridpack::myapp::MyBus::serialWrite(char *string,  
    const int bufsize, const char *signal)  
{  
    double pi = 4.0*atan(1.0);  
    double angle = p_a*180.0/pi;  
    sprintf(string, "      %6d      %12.6f      %12.6f\n",  
        getOriginalIndex(),angle,p_v);  
}
```

For this simple case, **signal** is ignored as well as the variable **bufsize**. If more than one type of bus listing was desired, additional conditions based on the value of **signal** could be included. For the case of generators, the length of the output may vary from one bus to the next since buses can have different numbers of generators associated with them. In this case it may be important to check the length of the output string being generated against the size of the buffer to make sure there is no overwrite and to take some kind of appropriate action if there is.

If you wish to direct the output to a file, then calling the function

```
void open(const char *filename)
```

will direct all output from the serial IO object to the file specified in the variable filename. Similarly, calling the function

```
void close(void)
```

will close the file and all subsequent writes are directed back to standard output. The same **SerialBusIO** object can be used to write data to multiple different files, as long as the files are opened and closed sequentially. If two files need to be used at the same time, then two **SerialBusIO** objects need to be created. Two additional methods can be used to further control where output goes. If a file already exists, you can use the function

```
boost::shared_ptr<std::ofstream> getStream()
```

to recover a pointer to the file stream currently being used by the **SerialBusIO** object. This can then be used to redirect output from some other part of the code to the same file. The function

```
void setStream(boost::shared_ptr<std::ofstream> stream)
```

can be used to redirect the output from the **SerialIOBus** object to an already existing file. The main use of these two functions is to direct the output from both buses and branches to the same file instead of standard output.

The **SerialBranchIO** module is similar to the **SerialBusIO** module but works by creating listings for branches. The constructor is

```
SerialBranchIO<MyNetwork>(int max_str_len,  
    boost::shared_ptr<MyNetwork> network)
```

and the header and write methods are

```
void header(const char *string) const
```

```
void write(const char *signal = NULL)
```

These have exactly the same behavior as in the **SerialBusIO** class. Similarly, the methods

```
void open(const char *filename)
```

```
void close(void)
```

```
boost::shared_ptr<std::ofstream> getStream()
```

```
void setStream(boost::shared_ptr<std::ofstream> stream)
```

can be used to redirect output to a file instead of standard output.

The usual method for directing the output from both a `SerialBusIO` object and `SerialBranchIO` object to the same file is to use the calling sequence

```
SerialBusIO<MyNetwork> busIO(max_str_len, network);
SerialBranchIO<MyNetwork> branchIO(max_str_len, network);
busIO.open("file.dat");
branchIO.setStream(busIO.getStream());
```

The file can be closed by calling `close` from either `busIO` or `branchIO`.

### Configuration Module

The configuration module is designed to provide a central mechanism for directing information from the input file to the components making up a given application. For example, information about convergence thresholds and maximum numbers of iterations might need to be picked up by the solver module from an external configuration file. The configuration module is designed to read files using a simple XML format that supports a hierarchical input. This can be used to control which input gets directed to individual objects in the application, even if the object is a framework component and cannot be modified by the application developer.

The `Configuration` class is in the namespace `gridpack::utility`. This class follows the C++ singleton pattern and does not have a public constructor. The static method `configuration()` returns a pointer to the shared instance of this class and guarantees that the same instance is used by all modules in an application. The shared instance can be initialized with data from an external file using the code

```
gridpack::utility::Configuration * c =
    gridpack::utility::Configuration::configuration() ;
c->open(input_file, MPI_COMM_WORLD);
```

The input file uses XML syntax. The single top-level element must be named “Configuration” but other elements may have module- and application-specific names. Refer elsewhere in this document for details. For illustration purposes, an example configuration file might look like:

```
<?xml version="1.0" encoding="utf-8"?>
<Configuration>
  <PowerFlow>
    <networkConfiguration> IEEE118.raw </networkConfiguration>
  </PowerFlow>
  <DynamicSimulation>
    <StartTime> 0.0 </StartTime>
    <EndTime> 0.1 </EndTime>
    <TimeStep> 0.001 </TimeStep>
    <Faults>
      <Fault>
        <StartFault> 0.03 </StartFault>
```

```

        <EndFault> 0.06 </EndFault>
        <Branch> 3 7 </Branch>
    </Fault>
    <Fault>
        <StartFault> 0.07 </StartFault>
        <EndFault> 0.06=8 </EndFault>
        <Branch> 4 8 </Branch>
    </Fault>
</Faults>
</DynamicSimulation>
</Configuration>

```

A value in this configuration file is accessed with a call to the overloaded method **get()**. The following line will return the value of the input file corresponding to the XML field **"networkConfiguration"**

```

std::string s =
    c->get("Configuration.PowerFlow.networkConfiguration",
        "IEEE.raw");

```

The first argument has type **Configuration::KeyType** which is a **typedef** of **std::string**. Values are selected by hierarchically named “keys” using “.” as a separator. In the example input file, “PowerFlow” is a block under “Configuration” and “networkConfiguration” is, in turn, a block under “PowerFlow”. The second argument in **get** is a default value that is used if the field corresponding to the key can’t be found. There are overloaded versions of **get()** for accessing standard C++ data types: **bool**, **int**, **long**, **float**, **double**, **ComplexType** and **std::string**. For each type there are two variants. For integers these look like

```

int get(const KeyType &, int default_value) const ;
bool get(const KeyType &, int *) const;

```

The first variant takes a key name and a default value and returns either the value in the configuration file or the default value when none is specified. In the second variant, a Boolean value is returned indicating whether or not the value was in the configuration file and the second argument points to an object that is updated with the configuration value when it is present. For strings, the second argument is passed by reference.

The method **getCursor(KeyType)** returns a pointer to an internal element in the hierarchy. This “cursor” supports the same **get()** methods as above but the names are now relative to the name of the cursor. Thus, we might modify the previous example to:

```

Configuration::CursorPtr p =
    c->getCursor("Configuration.PowerFlow");

std::string s = p->get("networkConfiguration",

```



```
"IEEE14.raw");
```

An additional use of such cursors is to access lists of values. The method

```
typedef std::vector<CursorPtr> ChildCursors;
```

```
void children(ChildCursors &);
```

can be used to get a vector of all the elements that are children in the name hierarchy of some element. These children need not have unique names, as illustrated by the children of the “Faults” element shown above. In this example, each of the children is a cursor that can be used to access “StartFault”, “EndFault”, and “Branch” parameters for each of the “Fault” blocks. Again, returning to the sample input above, the following code will return a list of faults

```
Configuration::CursorPtr p =  
    c->getCursor("Configuration.DynamicSimulation.Faults");  
ChildCursors faults;  
p->children(faults);
```

The cursor `p` is set so that it is pointing at the `Faults` block in the input. The `children` function then picks up all XML blocks one level and returns a list of cursor pointers to those blocks. The individual data elements in `faults` can be accessed using the following loop

```
int nfaults = faults.size()  
for (int i=0; i<nfaults; i++) {  
    double start = faults[i]->get("StartFault", 0.0);  
    double end = faults[i]->get("EndFault", 0.0);  
    std::string indices = faults[i]->get("Branch", "-1 -1");  
    // Do something with these parameters  
}
```

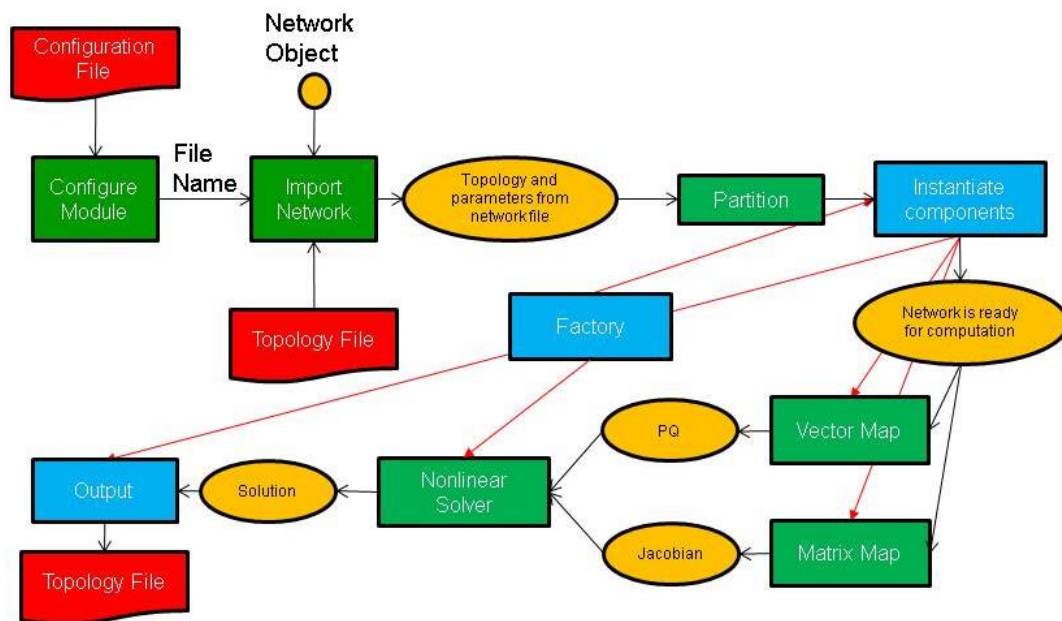
Note that this method does not have any way of distinguishing between different blocks below `Faults` and if two types of blocks were listed within the `Faults` block, the `children` method would pick up both of them.

## Developing Applications

The previous section outlined most of the basic modules in the GridPACK framework. In this section, we provide an overview of how to use these modules to create actual applications by discussing the development of a power flow simulation application in detail. Actual examples of a power flow application can be found by looking at an example code located under the top-level GridPACK directory in `src/applications/examples/powerflow`. Users can also look at the power flow module located in the `src/applications/modules/powerflow` directory. The main difference between the power flow example code and the power flow module is that the module breaks up

the power flow calculation into more separate function calls and the module also has options for using a non-linear solver. The power flow bus and branch classes are located in the directory `src/applications/components/pf_matrix`.

A schematic of a power flow code based on GridPACK is shown in Figure 8. For different power grid problems, the details of the code will be different, but most of these motifs will appear at some point or other. The main differences will probably be in feedback loops as results from one part of the calculation are fed back into other parts. For example, an iterative solver will need to update the current values of the network components, which can then be used to generate new matrices and vectors that are fed back into the next iteration of the solver. The diagram in Figure 8 is not complete, but gives an overall view of code structure and data movement.



**Figure 8.** Schematic of program flow for a power flow simulation. The yellow ovals are distributed data objects, the green blocks are GridPACK framework components and the blue blocks are application specific code. External files are red.

As shown in the figure, application developers will need to focus on writing two or three sets of modules. The first is the network components. These are the descriptions of the physics and/or measurements that are associated with buses and branches in the power grid network. The network factory is a module that initializes the grid components on the network after the network is originally created by the import module. The power flow problem is simple enough that it can use a non-linear solver directly from the math module but even a straightforward solution such as this requires the developer to overwrite some functions in the factory that are used in the non-linear solver iterations.

Most of the work involved in creating a new application is centered on creating the bus and branch classes. This discussion will describe in some detail the routines that need to be written in order to develop a working power flow simulation. Additional application modules for dynamic simulation and contingency analysis have also been included in the distribution and users are encouraged to look at these modules for additional coding examples on how to use GridPACK. The discussion below is designed to illustrate how to build an application and for brevity has left out some code compared to the working implementation. The source code contains more comment lines as well as some additional diagnostics that may not appear here. However, the overall design is the same and readers who have a good understanding of the following text should have no difficulty understanding the power flow source code.

For the power flow calculation, the buses and branches will be represented by the classes **PFBus** and **PFBranch**. **PFBus** inherits from the **BaseBusComponent** class, so it automatically inherits the **BaseComponent** and **MatVecInterface** classes as well. The first thing that must be done in creating the **PFBus** component is to overwrite the load function in the **BaseComponent** class. The original function is just a placeholder that performs no action. The **load** function should take parameters from the **DataCollection** object associated with each bus and use them to initialize the bus component itself. For the **PFBus** component, a simplified **load** function is

```
void gridpack::powerflow::PFBus::load(
    const boost::shared_ptr<gridpack::component
    ::DataCollection> &data)
{
    data->getValue(CASE_SBASE, &p_sbase);
    data->getValue(BUS_VOLTAGE_ANG, &p_angle);
    data->getValue(BUS_VOLTAGE_MAG, &p_voltage);
    p_v = p_voltage;
    double pi = 4.0*atan(1.0); p_angle = p_angle*pi/180.0;
    p_a = p_angle;
    int itype; data->getValue(BUS_TYPE, &itype);
    if (itype == 3) {
        setReferenceBus(true);
    }
    bool lgen;
    int i, ngen, gstatus;
    double pg, qg;
    if (data->getValue(GENERATOR_NUMBER, &ngen)) {
        for (i=0; i<ngen; i++) {
            lgen = true;
```

```

    lgen = lgen && data->getValue(GENERATOR_PG, &pg,i);
    lgen = lgen && data->getValue(GENERATOR_QG, &qg,i);
    lgen = lgen && data->getValue(GENERATOR_STAT, &gstatus,i);
    if (lgen) {
        p_pg.push_back(pg);
        p_qg.push_back(qg);
        p_gstatus.push_back(gstatus);
    }
}
}
}

```

This version of the **load** function has left off additional properties, such as shunts and loads and some transmission parameters, but it serves to illustrate how **load** is suppose to work. The **load** method in the base factory class will run over all buses, get the **DataCollection** object associated with each bus and then call the **PFBus::load** method using the **DataCollection** object as the argument. The parameters **p\_sbase**, **p\_angle**, **p\_voltage** are private members of **PFBus**. The variables corresponding to the keys **CASE\_SBASE**, **BUS\_VOLTAGE\_ANG**, **BUS\_VOLTAGE\_MAG** were stored in the **DataCollection** object when the network configuration file was parsed. They are retrieved from this object using the **getValue** functions and assigned to **p\_sbase**, **p\_angle**, **p\_voltage**. Additional internal variables are also assigned in a similar manner. The value of the **BUS\_TYPE** variable can be used to determine whether the bus is a reference bus. As mentioned previously, the **CASE\_SBASE** etc. are just preprocessor symbols that are defined in the **dictionary.hpp** file, which must be included in the file defining the **load** function. The **dictionary.hpp** file can be found in the **src/parser** directory of the GridPACK distribution.

The variables referring to generators have a different behavior than the other variables. A bus can have multiple generators and these are stored in the **DataCollection** object with an index. The total number of generators on the bus is also stored in the **DataCollection** object with the key **GENERATOR\_NUMBER**. First the number of generators is retrieved and then a loop is set up so that all the generator variables can be accessed. The generator parameters are stored in local private arrays. The loop shows how the return value of the **getValue** function can be used to verify that all three parameters for a generator were found. If they aren't found, then the generator is incomplete and the generator is not added to the local data. The boolean return value can also be used to determine if the bus has other properties and to set internal flags and parameters accordingly. The load function for the **PFBBranch** is constructed in a similar way, except that the focus is on extracting branch related parameters from the **DataCollection** object.

Both the **PFBus** and **PFBBranch** classes contain an application-specific function called **setYBus** that is used to set up values in the Y-matrix. There is also a function in the powerflow factory class that runs over all buses and branches and calls this function. The **setYBus** function in **PFBus** is

```
void gridpack::powerflow::PFBus::setYBus(void)
{
    gridpack::ComplexType ret(0.0,0.0);
    std::vector<boost::shared_ptr<BaseComponent> > branches;
    getNeighborBranches(branches);
    int size = branches.size();
    int i;
    for (i=0; i<size; i++) {
        gridpack::powerflow::PFBBranch *branch
            = dynamic_cast<gridpack::powerflow::PFBBranch*>
              (branches[i].get());
        ret -= branch->getAdmittance();
        ret -= branch->getTransformer(this);
        ret += branch->getShunt(this);
    }
    if (p_shunt) {
        gridpack::ComplexType shunt(p_shunt_gs,p_shunt_bs);
        ret += shunt;
    }
    p_ybusr = real(ret);
    p_ybusi = imag(ret);
}
```

This function evaluates the contributions to the Y-Matrix associated with buses. The real and imaginary parts of this number are stored in the internal variables **p\_ybusr** and **p\_ybusi**. The subroutine first creates the local variable **ret** and then gets a list of pointers to neighboring branches from the **BaseBusComponent** function **getNeighborBranches**. The function then loops over each of the branches and uses the dynamic cast function in C++ to convert the **BaseComponent** pointer to a **PFBBranch** pointer. Note that the cast is necessary since the **getNeighborBranches** function only returns a list of **BaseComponent** object pointers. The **BaseComponent** class does not contain application-specific functions such as **getAdmittance**. The **getAdmittance**, **getTransformer** and **getShunt** methods return the contributions from transmission elements, transformers and shunts associated with the branch. These are accumulated into the **ret** variable.

The reason that the **getAdmittance** variable has no argument while both **getTransformer** and **getShunt** take the pointer “**this**” as an argument is that the admittance contribution from simple transmission elements is symmetric with respect to whether or not the calling bus is the “from” or “to” buses while the transformer and shunt contributions are not. This can be seen by examining the **getTransformer** function.

```
gridpack::ComplexType
gridpack::powerflow::PFBranch::getTransformer(
    gridpack::powerflow::PFBUS *bus)
{
    gridpack::ComplexType ret(p_resistance,p_reactance);
    if (p_xform) {
        ret = -1.0/ret;
        gridpack::ComplexType a(cos(p_phase_shift),sin(p_phase_shift));
        a = p_tap_ratio*a;
        if (bus == getBus1().get()) {
            ret = ret/(conj(a)*a);
        } else if (bus == getBus2().get()) {
            // ret is unchanged
        }
    } else {
        ret = gridpack::ComplexType(0.0,0.0);
    }
    return ret;
}
```

The variables **p\_resistance**, **p\_reactance**, **p\_phase\_shift**, and **p\_tap\_ratio** are all internal variables that are set based on the variables read in from using the **load** method or are set in other initialization steps. The boolean variable **p\_xform** variable is set to true in the **PFBranch::load** method if transformer-related variables are detected in the **DataCollection** objects associated with the branch, otherwise it is false.

The **PFBranch** version of the **setYBus** function is

```
void gridpack::powerflow::PFBranch::setYBus(void)
{
    gridpack::ComplexType ret(p_resistance,p_reactance);
    ret = -1.0/ret;
    gridpack::ComplexType a(cos(p_phase_shift),sin(p_phase_shift));
    a = p_tap_ratio*a;
    if (p_xform) {
```

```

    p_ybusr_frwd = real(ret/conj(a));
    p_ybusi_frwd = imag(ret/conj(a));
    p_ybusr_rvrs = real(ret/a);
    p_ybusi_rvrs = imag(ret/a);
} else {
    p_ybusr_frwd = real(ret);
    p_ybusi_frwd = imag(ret);
    p_ybusr_rvrs = real(ret);
    p_ybusi_rvrs = imag(ret);
}
gridpack::powerflow::PFBUS *bus1 =
    dynamic_cast<gridpack::powerflow::PFBUS*>(getBus1().get());
gridpack::powerflow::PFBUS *bus2 =
    dynamic_cast<gridpack::powerflow::PFBUS*>(getBus2().get());
p_theta = (bus1->getPhase() - bus2->getPhase());
}

```

Note that the branch version of the **setYBus** function calculates different values for the Y-matrix contribution depending on whether the first index in the Y-matrix element corresponds to bus 1 (the forward direction) or bus 2 (the reverse direction). These are stored in the separate variables **p\_ybusr\_frwd** and **p\_ybusi\_frwd** for the forward directions and **p\_ybusr\_rvrs** and **p\_ybusi\_rvrs** for the reverse direction. This routine also calculates the variable **p\_theta** which is equal to the difference in the phase angle variable associated with the two buses at either end of the branch. This last variable provides an example of calculating a branch parameter based on the values of parameters located on the terminal buses.

The **setYBus** functions described above are used in the power flow components to set some basic parameters. These are eventually incorporated into the Jacobian matrix and PQ vector that constitute the matrix and right hand side vector of the power flow equations. To build the matrix, it is necessary to implement the matrix size and matrix values functions in the **MatVecInterface**. The functions for setting up the matrix are discussed in detail in the following, the vector functions are simpler but follow the same pattern. The mode used for setting up the Jacobian matrix is “**Jacobian**”. The corresponding **matrixDiagSize** routine is

```

bool gridpack::powerflow::PFBUS::matrixDiagSize(int *isize,
    int *jsize) const
{
    if (p_mode == Jacobian) {
        *isize = 2;
        *jsize = 2;
    }
}

```

```

        return true;
    } else if (p_mode == YBus) {
        *isize = 1;
        *jsize = 1;
        return true;
    }
}

```

This function handles two modes, stored in the internal variable **p\_mode**. If the mode equals **Jacobian**, then the function returns a contribution to a 2×2 matrix. In the case that the mode is “**YBus**” the function would return a contribution to a 1×1 matrix. (The Jacobian is treated as a real matrix where the real and complex parts of the problem are treated as separate variables, the Y-matrix is handle as a regular complex matrix). The corresponding code for returning the diagonal values is

```

bool gridpack::powerflow::PFBus::matrixDiagValues(ComplexType *values)
{
    if (p_mode == YBus) {
        gridpack::ComplexType ret(p_ybusr,p_ybusi);
        values[0] = ret;
        return true;
    } else if (p_mode == Jacobian) {
        if (!getReferenceBus()) {
            values[0] = -p_Qinj - p_ybusi * p_v *p_v;
            values[1] = p_Pinj - p_ybusr * p_v *p_v;
            values[2] = p_Pinj / p_v + p_ybusr * p_v;
            values[3] = p_Qinj / p_v - p_ybusi * p_v;
            if (p_isPV) {
                values[1] = 0.0;
                values[2] = 0.0;
                values[3] = 1.0;
            }
            return true;
        } else {
            values[0] = 1.0;
            values[1] = 0.0;
            values[2] = 0.0;
            values[3] = 1.0;
            return true;
        }
    }
}

```



```

    }
}

```

In this implementation, the return values are of type **ComplexType**, even if they are real. For real values, the complex part is set to zero. If the mode is “**YBus**”, the function returns a single complex value. If the mode is “**Jacobian**”, the function checks first to see if the bus is a reference bus or not. If the bus is not a reference bus, then the function returns a 2×2 block corresponding to the contributions to the Jacobian matrix coming from a bus element. If the bus is a reference bus, the function returns a 2×2 identity matrix. This is a result of the fact that the variables associated with a reference bus are fixed. In fact, the variables contributed by the reference bus could be eliminated from the matrix entirely by returning false if the mode is “**Jacobian**” and the bus is a reference bus for both the matrix size and matrix values routines. This would also require some adjustments to the off-diagonal routines as well. There is an additional condition for the case where the bus is a “PV” bus. In this case one of the independent variables is eliminated by setting the off-diagonal elements of the block to zero and the second diagonal element equal to 1. The values are returned in column-major order, so **values[0]** corresponds to the (0,0) location in the 2×2 block, **values[1]** is the (1,0) location, **values[2]** is the (0,1) location and **values[3]** is the (1,1) location.

The **matrixForwardSize** and **matrixForwardValues** routines, as well as the corresponding Reverse routines, are implemented in the **PFBranch** class. These functions determine the off-diagonal blocks of the Jacobian and Y-matrix. The **matrixForwardSize** routine is given by

```

bool gridpack::powerflow::PFBranch::matrixForwardSize(int *isize,
    int *jsize) const
{
    if (p_mode == Jacobian) {
        gridpack::powerflow::PFBUS *bus1
            = dynamic_cast<gridpack::powerflow::PFBUS*>(getBus1().get());
        gridpack::powerflow::PFBUS *bus2
            = dynamic_cast<gridpack::powerflow::PFBUS*>(getBus2().get());
        bool ok = !bus1->getReferenceBus();
        ok = ok && !bus2->getReferenceBus();
        if (ok) {
            *isize = 2;
            *jsize = 2;
            return true;
        } else {
            return false;
        }
    }
}

```

```

    } else if (p_mode == YBus) {
        *isize = 1;
        *jsize = 1;
        return true;
    }
}

```

If the mode is “**YBus**”, the size function returns a  $1 \times 1$  block for the off-diagonal matrix block. For the Jacobian, this function first checks to see if either end of the branch is a reference bus by evaluating the Boolean variable “ok”. If neither end is the reference bus then the function returns a  $2 \times 2$  block, if one end is the reference bus then the function returns false. The false value indicates that this branch does not contribute to the matrix. For this system, the **matrixReverseSize** function is the same. For applications that return a non-square block, the reverse function will transpose the block dimensions relative to the forward direction.

The **matrixForwardValues** function is

```

bool gridpack::powerflow::PFBranch::matrixForwardValues(
    ComplexType *values)
{
    if (p_mode == Jacobian) {
        gridpack::powerflow::PFBus *bus1
            = dynamic_cast<gridpack::powerflow::PFBus*>(getBus1().get());
        gridpack::powerflow::PFBus *bus2
            = dynamic_cast<gridpack::powerflow::PFBus*>(getBus2().get());
        bool ok = !bus1->getReferenceBus();
        ok = ok && !bus2->getReferenceBus();
        if (ok) {
            double cs = cos(p_theta);
            double sn = sin(p_theta);
            values[0] = (p_ybusr_frwd*sn - p_ybusi_frwd*cs);
            values[1] = (p_ybusr_frwd*cs + p_ybusi_frwd*sn);
            values[2] = (p_ybusr_frwd*cs + p_ybusi_frwd*sn);
            values[3] = (p_ybusr_frwd*sn - p_ybusi_frwd*cs);
            values[0] *= ((bus1->getVoltage())*(bus2->getVoltage()));
            values[1] *= -((bus1->getVoltage())*(bus2->getVoltage()));
            values[2] *= bus1->getVoltage();
            values[3] *= bus1->getVoltage();
            bool bus1PV = bus1->isPV();
            bool bus2PV = bus2->isPV();
            if (bus1PV & bus2PV) {

```

```

        values[1] = 0.0;
        values[2] = 0.0;
        values[3] = 0.0;
    } else if (bus1PV) {
        values[1] = 0.0;
        values[3] = 0.0;
    } else if (bus2PV) {
        values[2] = 0.0;
        values[3] = 0.0;
    }
    return true;
} else {
    return false;
}
} else if (p_mode == YBus) {
    values[0] = gridpack::ComplexType(p_ybusr_frwd,p_ybusi_frwd);
    return true;
}
}

```

For the “**YBus**” mode, the function simply returns the complex contribution to the Y-matrix. For the “**Jacobian**” mode, the function first determines if either end of the branch is connected to the reference bus. If it is, then the function returns false and there is no contribution to the Jacobian. If neither end of the branch is the reference bus then the function evaluates the 4 elements of the 2×2 contribution to the Jacobian coming from the branch. To do this, the branch needs to get the current values of the voltages on the buses at either end by using the **getVoltage** accessor functions that have been defined in the **PFBus** class. Finally, if one end or the other of the branch is a PV bus, then some variables need to be eliminated from the equations. This can be done by setting appropriate values in the 2×2 block equal to zero.

The **matrixReverseValues** function is similar to the **matrixForwardValues** functions with a few key differences. 1) the variables **p\_ybusr\_rvrs** and **p\_ybusi\_rvrs** are used instead of **p\_ybusr\_frwd** and **p\_ybusi\_frwd** 2) instead of using **cos(p\_theta)** and **sin(p\_theta)** the function uses **cos(-p\_theta)** and **sin(-p\_theta)** since **p\_theta** is defined as difference in phase angle on bus 1 minus the difference in phase angle on bus 2 and 3) the values that are set to zero in the conditions for PV buses are transposed. The PV conditions are the same as the forward case if both bus 1 and bus 2 are PV buses, if only bus 1 is a PV bus then **values[2]** and **values[3]** are zero and if only bus2 is a PV bus then **values[1]** and **values[3]** are zero.

The functions for setting up vectors are similar to the corresponding matrix functions, although they are a bit simpler. The vector part of the **MatVecInterface** contains one function that does not have a counterpart in the set of matrix functions and that is the **setValues** function. This function can be used to push values in a vector object back into the buses that were used to generate the vector. For the Newton-Raphson method used to solve the power flow equations, it is necessary, at each iteration, to push the current solution back into the buses so they can be used to evaluate new Jacobian and right hand side vectors. The solution vector contains the current increments to the voltage and phase angle. These are written back to the buses using the function

```
void gridpack::powerflow::PFBus::setValues(
    gridpack::ComplexType *values)
{
    p_a -= real(values[0]);
    p_v -= real(values[1]);
    *p_vAng_ptr = p_a;
    *p_vMag_ptr = p_v;
}
```

This function is paired with a mapper that is used to create a vector with the same pattern of contributions. If for example, the matrix equation  $\mathbf{Ax} = \mathbf{b}$  is being solved, then the mapper used to create the right hand side vector  $\mathbf{b}$  should be used to push results back onto the buses using the **mapToBus** method. The **setValues** method above takes the contributions from the solution vector and uses them to decrement the internal variables **p\_a** (voltage angle) and **p\_v** (voltage magnitude). The new values of **p\_a** and **p\_v** are then assigned to the buffers **p\_vAng\_ptr** and **p\_vMag\_ptr** so that they can be exchanged with other buses. This is discussed below.

The two routines that need to be created in the **PFBus** class to copy data to ghost buses are both simple. There is no need to create corresponding routines in the **PFBBranch** class since branches do not exchange data in the power flow calculation. Two values need to be exchanged between buses, the current voltage angle and the current voltage magnitude. This requires a buffer that is the size of two doubles so the **getXCBufSize** function is written as

```
int gridpack::powerflow::PFBus::getXCBufSize(void)
{
    return 2*sizeof(double);
}
```

The **setXCBuf** assigns the buffer created in the base factory **setExchange** function to internal variables used within the **PFBus** component. It has the form

```
void gridpack::powerflow::PFBus::setXCBuf(void *buf)
```

```

{
    p_vAng_ptr = static_cast<double*>(buf);
    p_vMag_ptr = p_vAng_ptr+1;
    *p_vAng_ptr = p_a;
    *p_vMag_ptr = p_v;
}

```

The buffer created in the **setExchange** routine is split between the two internal pointers **p\_vAng\_ptr** and **p\_vMag\_ptr**. These are then initialized to the current values of **p\_a** and **p\_v**. Whenever the **updateBuses** routine is called, the buffers on the ghost buses are refreshed with the current values of the variables from the processes that own the corresponding buses. Note that both the **getXCBufSize** and the **setXCBuf** routines are only called during the **setExchange** routine. They are not called during the actual bus updates.

One final function in the **PFBus** and **PFBranch** class that is worth taking a brief look at is the set mode function. This function is used to set the internal **p\_mode** variable that is defined in both classes. The **PFMode** enumeration, which contains both the “YBus” and “Jacobian” modes, is defined within the `gridpack::powerflow` namespace. The **setMode** function for both buses and branches has the form

```

void gridpack::powerflow::PFBus::setMode(int mode)
{
    p_mode = mode;
}

```

This function is triggered on all buses and branches if the **setMode** method in the factory class is called.

Once the **PFBus** and **PFBranch** classes have been defined, it is possible to define a **PFNetwork** using a **typedef** statement. This can be done using the line

```

typedef network::BaseNetwork<PFBus, PFBranch > PFNetwork;

```

in the header file declaring the **PFBus** and **PFBranch** classes. This type can then be used in other powerflow files that need to create objects from templated classes.

The discussion above summarizes many of the important functions in the **PFBus** and **PFBranch** classes. Additional functions are included in these classes that are not discussed here, but the basic principles involved in implementing the remaining functions have been covered.

The first part of creating a new application is writing the network component classes. The second part is implementing the application-specific factory. For the power flow application, this is the

**PFFactory** class, which inherits from the **BaseFactory** class. Most of the important functionality in **PFFactory** is derived from the **BaseFactory** class and is used without modification, but several application-specific functions have been added to **PFFactory** that are used to set internal parameters in the network components. As an example, consider the **setYBus** function

```
void gridpack::powerflow::PFFactory::setYBus(void)
{
    int numBus = p_network->numBuses();
    int numBranch = p_network->numBranches();
    int i;
    for (i=0; i<numBus; i++) {
        p_network->getBus(i).get()->setYBus();
    }
    for (i=0; i<numBranch; i++) {
        p_network->getBranch(i).get()->setYBus();
    }
}
```

This function loops over all buses and branches and invokes the **setYBus** method in the individual **PFBus** and **PFBranh** objects. The first two lines in the factory **setYBus** method get the total number of buses and branches on the process. A loop over all buses on the process is initiated and a pointer to the bus object is obtained via the **getBus** bus method in the **BaseNetwork** class. This pointer is returned as a pointer of type **PFBus**, so it is not necessary to do a dynamic cast on it and the **setYBus** method, which is not part of the base class, can be invoked. The same set of steps is then repeated for the branches. The factory can be used to create other methods that invoke functions on buses and/or branches. Most of these functions follow the same general form as the **setYBus** method just described.

The last part of building an application is creating the top level application driver that actually instantiates all the objects used in the calculation and controls the program flow. Running the code is broken up into two parts. The first is creating a main program and the second is creating the application driver. The main routine is primarily responsible for initializing the communication libraries and creating the application object, the application object then controls the application itself. The main program for the powerflow application is

```
main(int argc, char **argv)
{
    gridpack::parallel::Environment env(argc,argv);
    gridpack::math::Initialize();
```

```

gridpack::powerflow::PFApp app;
app.execute();

gridpack::math::Finalize();
}

```

The first line in this program creates a variable of type **Environment** that initializes the MPI and GA communication libraries (the initialization happens in the constructor, so all that is necessary is to create the variable). The second line initializes the math library, which, in turn, calls the initialization routines of whatever library the math module is built on. The code then instantiates a power flow application object and calls the execute method for this object. The remainder of the power flow application is contained in the **PFApp::execute** method. Finally, when the application has finished running, the main program cleans up the communication and math libraries. The communication libraries are handled when the **env** variable goes out of scope and calls the **Environment** destructor. The main reason for breaking the code up in this way instead of including the execute function as part of **main** is to force the invocation of all the destructors in the GridPACK objects used to implement the application. Otherwise, these destructors get called after the communication libraries have been finalized and the program will fail to exit cleanly.

The top level control of the application is embedded in the power flow **execute** method. The **execute** method starts off with the code

```

gridpack::parallel::Communicator world;
boost::shared_ptr<PFNetwork> network(new PFNetwork(world));

gridpack::utility::Configuration *config
    = gridpack::utility::Configuration::configuration();
config->open("input.xml",world);
gridpack::utility::Configuration::Cursor *cursor;
cursor = config->getCursor("Configuration.Powerflow");
std::string filename;
if (!cursor->get("networkConfiguration", &filename)) {
    printf("No network configuration specified\n");
    return;
}
gridpack::parser::PTI23_parser<PFNetwork> parser(network);
parser.parse(filename.c_str());

network->partition();

```

The first two lines create a communicator for this application and use it to instantiate a **PFNetwork** object (note that this is really a **BaseNetwork** template class that is instantiated using the **PFBus** and **PFBranch** classes as template arguments). The network object exists but has no buses or branches associated with it. The next few lines get an instance of the configuration object and use this to open the **input.xml** file. This filename has been hardwired into this implementation but it could be passed in as a runtime argument, if desired. The code then creates a **Cursor** object and initializes this to point into the **Configuration.Powerflow** block of the **input.xml** file. The cursor can then be used to get the contents of the **networkConfiguration** block in **input.xml**, which corresponds to the name of the network configuration file containing the power grid network. This file is assumed to use the PSS/E version 23 format. After getting the file name, the code creates a **PTI23\_parser** object and passes in the current network object as an argument. When the parse method is called, the parser reads in the file specified in **filename** and uses that to add buses and branches to the network object. At this point, the network has all the bus and branches from the configuration file, but no ghost buses or branches exist and buses and branches are not distributed in an optimal way. Calling the partition method on the network then distributes the buses and branches and adds appropriate ghost buses and branches.

The next set of calls initialize the network components and prepare the network for computation.

```
gridpack::powerflow::PFFactory factory(network);
factory.load();

factory.setComponents();
factory.setExchange();

network->initBusUpdate();

factory.setYBus();
```

The first call creates a **PFFactory** object and instantiates it with a reference to the current network. **PFFactory** is defined as an instance of the **BaseFactory** class using a **PFNetwork** as the template argument. The next line calls the **BaseFactory load** method which invokes the component **load** method on all buses and branches. These use data from the **DataCollection** objects to initialize the corresponding bus and branch objects. Note that when the partition function creates the ghost bus and branch objects, it copies the associated **DataCollection** objects to these ghosts so the parameters from the network configuration file are available to instantiate all objects in the network. There is no need to do a data exchange at this point in the code in order to get current values on the ghost objects.



The next two calls are also implemented as **BaseFactory** methods. The **setComponents** method sets up pointers in the network components that point to neighboring branches and buses (in the case of buses) and terminal buses (in the case of branches). It is also responsible for setting up internal indices that are used by the mapper functions to create matrices and vectors. The **setExchange** method sets up the buffers that are used to exchange data between locally owned buses and branches and their corresponding ghost images on other processors. The call to **initBusUpdate** creates internal data structures that are used to exchange bus data between processors and the final factory call to **setYBus** evaluates the Y-matrix contributions from all network components. The network is fully initialized at this point and ready for computation.

The next calls create the Y-matrix and the matrices used in the Newton-Raphson iteration loop.

```
factory.setMode(YBus);
gridpack::mapper::FullMatrixMap<PFNetwork> mMap(network);
boost::shared_ptr<gridpack::math::Matrix> Y = mMap.mapToMatrix();

factory.setSBus();
factory.setMode(RHS);
gridpack::mapper::BusVectorMap<PFNetwork> vMap(network);
boost::shared_ptr<gridpack::math::Vector> PQ = vMap.mapToVector();

factory.setMode(Jacobian);
gridpack::mapper::FullMatrixMap<PFNetwork> jMap(network);
boost::shared_ptr<gridpack::math::Matrix> J = jMap.mapToMatrix();
boost::shared_ptr<gridpack::math::Vector> X(PQ->clone());
```

The first call sets the internal `p_mode` variable in all network components to “**YBus**”. The second call constructs a **FullMatrixMap** object **mMap** and the third call uses the **mapToMatrix** method to generate a Y-matrix based on the “**YBus**” mode. The factory then calls the **setSBus** method that sets some additional network component parameters (again, by looping over all buses and invoking a **setSBus** method on each bus). The next three lines set the mode to “**RHS**”, create a **BusVectorMap** object and create the right hand side vector in the powerflow equations using the **mapToVector** method. This builds the vector based on the “**RHS**” mode. The next three lines create the Jacobian using the same pattern as for the Y-matrix. The mode gets set to “**Jacobian**”, another **FullMatrixMap** object is created and this is used to create the Jacobian using the **mapToMatrix** method. Two separate mappers are used to create the Y-matrix and the Jacobian. This is required unless there is some reason to believe that the “**YBus**” and “**Jacobian**” modes generate matrices with the same dimensions and *exactly* the same fill pattern. This is not generally the case, so different mappers should be created for

each matrix in the problem. The last line creates a new vector by cloning the PQ vector. The X vector has the same dimension and data layout as PQ so it could be used with the **vMap** object.

Once the vectors and matrices for the problem have been created and set to their initial values, it is possible to start the Newton-Raphson iterations. The code to set up the first Newton-Raphson iteration is

```
double tolerance = 1.0e-6;
int max_iteration = 100;
ComplexType tol;

gridpack::math::LinearSolver solver(*J);
solver.configure(cursor);

int iter = 0;

X->zero();
solver.solve(*PQ, *X);
tol = PQ->normInfinity();
```

The first three lines define some parameters used in the Newton-Raphson loop. The tolerance and maximum number of iterations are hardwired in this example but could be made configurable via the input deck using the **Configuration** class. The next line creates a linear solver based on the current value of the Jacobian, **J**. The call to the configure method allows configuration parameters in the input file to be passed directly to the newly created solver. The iteration counter is set to zero and the value of **X** is also set to zero. The linear solver is called with **PQ** as the right hand side vector and **X** as the solution. An initial value of the tolerance is set by evaluating the infinity norm of **PQ**. The calculation can now enter the Newton-Raphson iteration loop

```
while (real(tol) > tolerance && iter < max_iteration) {
    factory.setMode(RHS);
    vMap.mapToBus(X);
    network->updateBuses();

    vMap.mapToVector(PQ);
    factory.setMode(Jacobian);
    jMap.mapToMatrix(J);

    X->zero();
    solver.solve(*PQ, *X);
```

```

    tol = PQ->normInfinity();
    iter++;
}

```

This code starts by pushing the values of the solution vector back on to the buses using the same mapper that was used to create **PQ**. The network then calls the **updateBus** routine so that the ghost buses have new values of the voltage angle and magnitude parameters from the solution vector. New values of the Jacobian and right hand side vector are created based on the solution values from the previous iteration. Note that since **J** and **PQ** already exist, the mappers are just overwriting the old values instead of creating new data objects. The linear solver is already pointing to the Jacobian matrix so it automatically uses the new Jacobian values when calculating the solution vector **x**. If the norm of the new **PQ** vector is still larger than the tolerance, the loop goes through another iteration. This continues until the tolerance condition is satisfied or the number of iterations reaches the value of **max\_iteration**.

If the Newton-Raphson loop converges, then the calculation is essentially done. The last part of the calculation is to write out the results. This can be accomplished using the code

```

gridpack::serial_io::SerialBusIO<PFNetwork> busIO(128,network);
busIO.header("\n  Bus Voltages and Phase Angles\n");
busIO.header("\n  Bus Number      Phase Angle");
busIO.header("      Voltage Magnitude\n");
busIO.write();

```

The first line creates a serial bus IO object that assumes that no line of output will exceed more than 128 characters. The next three lines write out the header for the output data and the last line writes a listing of data from all buses. The data from each bus is generated by the **serialWrite** method defined in the **PFBus** class. A similar set of calls can be used to write out data from the branches. This completes the execute method and the overview of the power flow application.

## Advanced Functionality

The core operations supported by GridPACK have been described above and these can be used in to create many different kinds of power grid applications. This section will describe features that are more advanced but can be extremely useful in certain cases. Additional capabilities of the GridPACK framework include

- Communicators and task managers that can be used to create multiple levels of parallelism and implement dynamic load balancing schemes
- A generalized matrix-vector interface to handle applications where the dependent and independent variables are associated with both buses and branches. The

**MatVecInterface** described above can only be used for systems where the dependent and independent variables are associated solely with buses

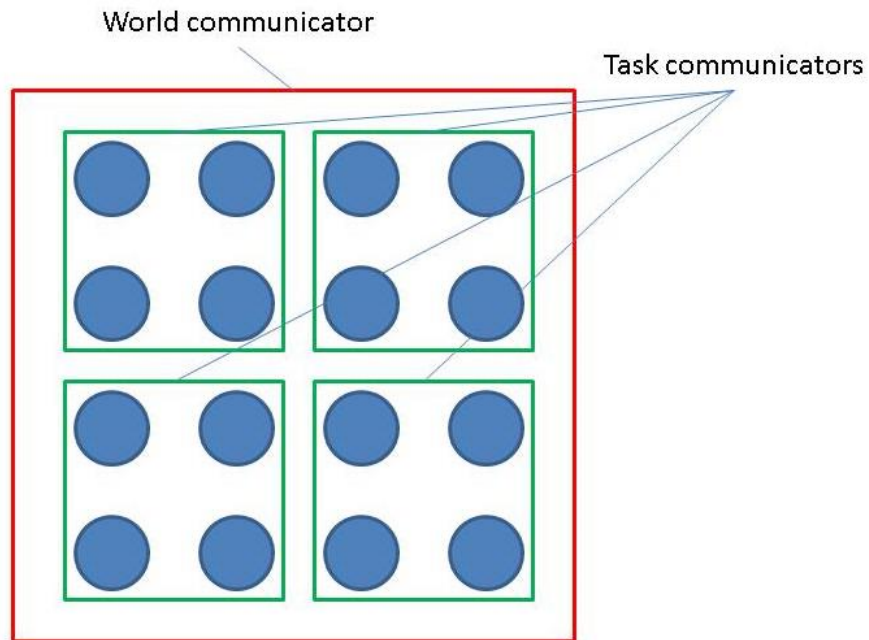
- A “slab” matrix-vector interface for creating matrices based on multiple values on each of the network components. This can be used for algorithms such as Kalman filter
- Profiling and error handling capabilities
- A hashed data distribution capability that can be used to direct network data to the processors that own the corresponding network components

This functionality is described in more detail in the following sections.

## Communicators

The subject of communicators has already been mentioned in the context of the constructor for the **BaseNetwork** class. This section will describe communicators in more detail and will show how the GridPACK communicators can be used to partition a large calculation into separate pieces that all run concurrently. A communicator can crudely be thought of as a communication link between a group of processors. Whenever a process needs to communicate with another process it needs to specify the communicator over which that communication will occur. When a parallel job is started, it creates a “world” communicator to which all processes implicitly belong. Any process can communicate with any other process via the world communicator. Other communicators can be created by an application and it is possible for a process to belong to multiple communicators. The concept of communicators is particularly important for restricting the scope of “global” operations. These are operations that require every process in the communicator to participate. Failure of a process to participate in the operation usually results in the calculation stalling because multiple processors are waiting for a communication from a process that is not part of the global operation. A program can remain in this state indefinitely. Many of the module functions in GridPACK represent global operations and contain imbedded calls that act collectively on a communicator. In order for two separate calculations to proceed concurrently, they must be run on disjoint sets of processors using separate communicators.

The use of communicators to create multiple concurrent parallel tasks within an application is usually straightforward to implement but it is frequently much more confusing to understand. A diagram of a set of 16 processes that are divided into 4 groups each containing 4 processes is shown schematically in Figure 9. In this example, each subgroup could potentially execute a separate parallel task within the larger parallel calculation.



**Figure 9.** Schematic diagram illustrating the use of multiple communicators

Global operations on the world communicator involve all 16 processes, global operations on one of the task communicators just involve the 4 processes in the group used to define the task communicator. If a network object is created on one of the task communicators, then a global operation such as the bus update only occurs between the 4 processes in the task communicator. The network object is, in a certain sense, “invisible” to the processes outside that communicator. If a network is created on a sub-communicator, then all objects derived from the network, such as factories, parsers, serial IO objects, etc. are also associated with the same sub-communicator.

The communicator supports some basic operations that are commonly used in parallel programming. GridPACK has been designed to minimize the amount of explicit communication that must be handled by application developers, but it is occasionally useful to have access to standard communication protocols in applications. In particular, it is useful to be able to divide a given communicator into a set of non-overlapping sub-communicators. The basic operations supported by the GridPACK communicator class are described below.

The GridPACK **Communicator** class is in the **gridpack::parallel** namespace. The constructor for this class creates a copy of the world communicator. The constructor has the form

**Communicator(void)**

and takes no arguments. Two basic functions associated with communicators are

```
int size(void) const
```

and

```
int rank(void) const
```

The first function returns the number of processors in the communicator and the second returns the index of the processor within the communicator. If the communicator contains  $N$  processes, then the rank will be an integer ranging from 0 to  $N-1$ . The process corresponding to rank 0 is often referred to as the head process or head node for the communicator. Note that if a process belongs to more than one communicator, its rank may differ depending on which communicator is being referred to. Information on size and rank is used extensively when explicitly programming in parallel. GridPACK has tried to abstract much of this programming so that developers do not need to pay attention to it, but it is still occasionally useful to be able to access these numbers. For example, the header function in the SerialIO classes is essentially equivalent to the following code fragment

```
Communicator comm;  
char buf[128];  
sprint(buf, "My message\n");  
if (comm.rank() == 0) {  
    printf("%s", buf);  
}
```

This code creates some output. If the condition was not there, the code would print out the message from all  $N$  processors in the world communicator and  $N$  copies of “My message” would appear in the output. The condition restricting the print statement to process 0 guarantees that the message appears only once.

A more important use of communicators is to divide up the world communicator into separate communicators that can be used to run independent parallel calculations. This is known as multi-level parallelism. Two functions can be used to split up an existing communicator into sub-communicators. The first is **split**

```
Communicator split(int color) const
```

This function divides the calling communicator into sub-communicators based on the **color** variable. All processors with the same value of the **color** variable end up in the same communicator. Thus, if 16 processors on the world communicator are divided up such that processes 0-3 are color 0, processes 4-7 are color 1, processes 8-11 are color 2 and processes 12-15 are color 3, then **split** will generate 4 sub-communicators from the world communicator such that 0-3 are on one communicator, 4-7 are on another communicator and so on. Note that this function divides the communicator completely into complementary pieces with all processes in

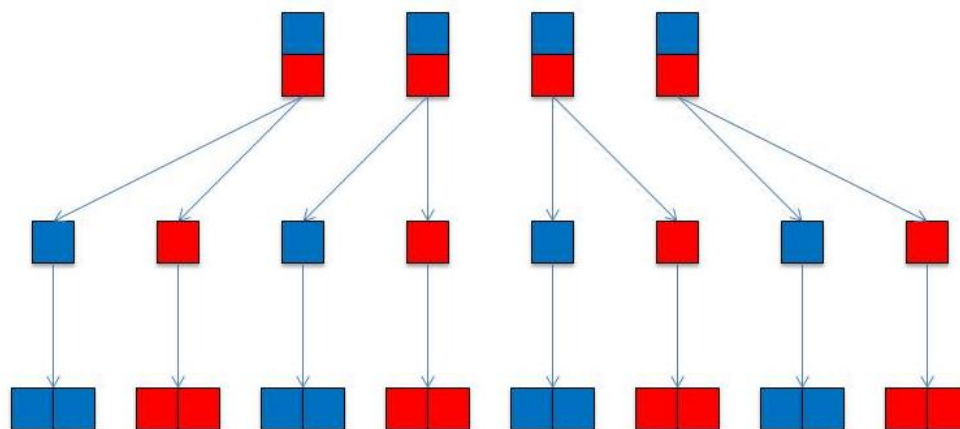
the old communicator ending up in a new communicator and no process ending up in more than one new communicator.

A second function that can be used to decompose a communicator into sub-communicators is **divide**. This function has the form

**Communicator divide(int nsize) const**

Each sub-communicator returned by this function contains at most **nsize** processes. The function will try and create as many communicators of size **nsize** as possible. For example, if the calling communicator contains 10 processes and **nsize** is set to 4, then this function will create 3 sub-communicators, two of which contain 4 processors and one containing 2 processors.

An example of how communicators can be used to create multiple levels of parallelism is illustrated in Figure 10. The example has 8 tasks that can be evaluated independently. The first row in Figure 10 shows four processors. Two of the 8 tasks are run on each processor so if each task has been parallelized then it needs to run on a communicator with only 1 processor in it. The second row shows the same calculation running on 8 processors. In this case, each processor only has 1 task associated with it but each task is still running on a single processor. If the tasks have not been parallelized, then this is as far as you can go. However, if the tasks *have* been parallelized, then you can move on to the configuration shown in the third line using 16 processors. In this case, the system has been divided into 8 groups, each containing two processors. Each group has its own separate subcommunicator and each task can be run in parallel on two processors. This gives additional speed-up over what can be achieved by simply distributing tasks to separate processors.



**Figure 10.** Schematic diagram of 8 tasks evaluated using multiple levels of parallelism. The first row represents 8 tasks on 4 processors, the second row is 8 tasks on 8 processors and the third row is 8 tasks running on 16 processors.

Additional functions are available for communicators that support basic parallel computing tasks. The objective of GridPACK is to abstract most aspects of parallel computing so that users do not need to deal with them directly, but there are some tasks, particularly those associated with collecting and organizing data, that are not difficult to program but are difficult to generalize. Some support for simple parallel operations is useful in these cases. The following operations can be used to sum data across all processors

```
void sum(float *x, int nvals) const
void sum(double *x, int nvals) const
void sum(int *x, int nvals) const
void sum(long *x, int nvals) const
void sum(ComplexType *x, int nvals) const
```

The array **x** holds the values to be summed and **nvals** is the number of values in **x**. This operation can be used to compute the total of some quantity after partial sums have been calculated on each processor. It can also be used to collect an array of values across a collection of processors by having each processor compute a portion of an array and then using the sum operation to create a complete copy of the array on all processors.

Maximum and minimum values can be calculated using the functions

```
void max(float *x, int nvals) const
void max(double *x, int nvals) const
void max(int *x, int nvals) const
void max(long *x, int nvals) const
```

```
void min(float *x, int nvals) const
void min(double *x, int nvals) const
void min(int *x, int nvals) const
void min(long *x, int nvals) const
```

Again, a global maximum or minimum can be calculated by first computing the local maximum or minim on each processor and then evaluating it across processors.

One other common parallel construct that may be useful in some contexts is the barrier or synchronization function. In GridPACK, this is available as the function

```
void sync() const
```



The **sync** function does not allow any processor in the communicator to proceed beyond this call until all processors in the communicator have reached the call. This is used in some parallel programming situations to guarantee a consistent state across all processors. In general, there should be relatively little need for this call in GridPACK. See, however, the comment below at the end of the section on **GlobalStore**.

### Task Manager

The task manager functionality is designed to parcel out tasks on a first come, first serve basis to processes in a parallel application. Each processor can request a task ID from the task manager and based on the value it receives, it will execute a block of work corresponding to the ID. The task manager guarantees that all IDs are sent out once and only once. The unique feature of the task manager is that if the tasks take unequal amounts of time, then processes with longer tasks will make fewer requests to the task manager than processes that have relatively short tasks. This leads to an automatic dynamic load balancing of the application that can substantially improve performance. The task manager also supports multi-level parallelism and can be used in conjunction with the sub-communicators described above to implement parallel tasks within a parallel application. An example of the use of communicators and task managers to create a code that uses multiple levels of parallelism can be found in the contingency analysis application that is part of the GridPACK distribution.

Task managers use the **gridpack::parallel** namespace. Task managers can be created either on the world communicator or on a subcommunicator. Two constructors are available.

#### TaskManager(void)

#### TaskManager(Communicator comm)

The first constructor must be called on all processors in the system and creates a task manager on the world communicator, the second is called on all processors within the communicator **comm**. Once the task manager has been created, the number of tasks must be set. This can be done with the function

#### void set(int ntask)

where the variable **ntask** corresponds to the total number of tasks to be performed. This call is collective on all processes in the communicator and each process must use the same value of **ntask**. The task IDs returned by the task manager will range from 0 to **ntask-1**.

Once the task manager has been created, task IDs can be retrieved from the task manager using one of the functions

#### bool nextTask(int \*next)

```
bool nextTask(Communicator &comm, int *next)
```

The first function is called on a single processor and returns the task ID in the variable **next**. The second is called on the communicator **comm** by all processors in **comm** and returns the same task ID on all processors (note that if all processors in **comm** called the first **nextTask** function, each processor in **comm** would end up with a different task ID). The communicator argument in the second **nextTask** call should be a subcommunicator relative to the communicator that was used to create the task manager. Both functions return true if the task manager has not run out of tasks, otherwise they return false and the value of **next** is set to -1.

The task manager also has a function

```
void printStats(void)
```

that can be used to print out information to standard out about how many tasks were assigned to each process.

A simple code fragment shows how communicators and task managers can be combined to create an application exhibiting multi-level parallelism.

```
gridpack::parallel::Communicator world  
int grp_size = 4;  
gridpack::parallel::Communicator task_comm = world.divide(grp_size);  
App app(task_comm);  
gridpack::parallel::TaskManager taskmgr;  
taskmgr.set(ntasks);  
int task_id;  
while(taskmgr.nextTask(task_comm, &task_id) {  
    app.execute(task_data[task_id]);  
}
```

This code divides the world communicator into sub-communicators containing at most 4 processes. An application is created on each task communicator and a task manager is created on the world group. The task manager is set to execute **ntasks** tasks and a while loop is created to execute each task. Each call to **nextTask** returns the same value of **task\_id** to the processors in **task\_comm**. This ID is used to index into an array **task\_data** of data structures that contain the input data necessary to execute the task. The size of **task\_data** corresponds to the value of **ntasks**. When the task manager runs out of tasks, the loop terminates. Note that this structure does not guarantee that tasks are mapped to processors in any fixed order. There is no guarantee that task 0 is executed on process 0 or that some process will execute a given number of tasks. If one task takes significantly longer than other tasks then it is likely that other processors will pick up work from the processors executing the longer task. This balances the

workload if each process is involved in multiple tasks. Once the workload drops to 1 task per process, this advantage is lost.

## Timers

Profiling applications is an important part of characterizing performance, identifying bottlenecks and proposing remedies. Profiling in a parallel context is also extremely tricky. Unbalanced applications can lead to incorrect conclusions about performance when load imbalance in one part of the application appears as poor performance in another part of the application. This occurs because the part of the application that appears slow has a global operation that acts as an accumulation point for load imbalance. Nevertheless, the first step in analyzing performance is to be able to time different parts of the code. GridPACK provides a timer functionality that can help users do this. These modules are designed to do relatively coarse-grained profiling, they should not be used to time the inside of computationally intensive loops.

GridPACK contains two different types of timers. The first is a global timer that can be used anywhere in the code and accumulates all results back to the same place for eventual display. Users can get a copy of this timer from any point in the calculation. The second timer is created locally and is designed to only time portions of the code. The second class of timers was created to support task based parallelism where there was an interest in profiling individual tasks instead of getting timing results averaged over all tasks. Both timers can be found in the **gridpack::utility** namespace.

The **CoarseTimer** class represents a timer that is globally accessible from any point in the code. A pointer to this timer can be obtained by calling the function

```
static CoarseTimer *instance()
```

A category within the timer corresponds to a set of things that are to be timed. A new category in the timer can be created using the command

```
int createCategory(const std::string title)
```

This command creates a category that is labeled by the name in the string **title**. The function returns an integer handle that can be used in subsequent timing calls. For example, suppose that all calls to **function1** within a code need to be timed. The first step is to get an instance of the timer and create the category “**Function1**”

```
gridpack::utility::CoarseTimer *timer =  
    gridpack::utility::CoarseTimer::instance();  
int t_func1 = timer->createCategory(“Function1”);
```

This code gets a copy of the timer and returns an integer handle **t\_func1** corresponding to this category. If the category has already been created, then **createCategory** returns a handle to the existing category, otherwise it adds the new category to the timer.

Time increments can be accumulated to this category using the functions

```
void start(const int idx)
void stop(const int idx)
```

The **start** command begins the timer for the category represented by the handle **idx** and **stop** turns the timer off and accumulates the increment.

At the end of the program, the timing results for all categories can be printed out using the command

```
void dump(void) const
```

The results for each category are printed to standard out. An example of a portion of the output from **dump** for a power flow code is shown below.

```
Timing statistics for: Total Application
  Average time:          14.7864
  Maximum time:          14.7864
  Minimum time:          14.7863
  RMS deviation:         0.0000
Timing statistics for: PTI Parser
  Average time:          0.1553
  Maximum time:          1.2420
  Minimum time:          0.0000
  RMS deviation:         0.4391
Timing statistics for: Partition
  Average time:          2.8026
  Maximum time:          2.9668
  Minimum time:          1.7142
  RMS deviation:         0.4398
Timing statistics for: Factory
  Average time:          1.2424
  Maximum time:          1.2540
  Minimum time:          1.2336
  RMS deviation:         0.0056
Timing statistics for: Bus Update
  Average time:          0.0019
  Maximum time:          0.0025
  Minimum time:          0.0016
  RMS deviation:         0.0003
```

For each category, the dump command prints out the average time spent in that category across all processors, the minimum and maximum times spent on a single processor and the RMS standard deviation from the mean across all processors. It is also possible to get more detailed output from a single category. The commands

```
void dumpProfile(const int idx) const  
void dumpProfile(const std::string title)
```

can both be used to print out how much time was spent in a single category across all processors. The first command identifies the category through its integer handle, the second via its name.

Some other timer commands also can be useful. The function

```
double currrentTime()
```

returns the current time in seconds (if you want to do timing on your own). If you want control profiling in different sections of the code the command

```
void configureTimer(bool flag)
```

can be used to turn timing off (**flag = false**) or on (**flag = true**). This can be used to restrict timing to a particular section of code and can be used for debugging and performance tuning.

In addition to the **CoarseTimer** class, there is a second class of timers called **LocalTimer**. **LocalTimer** supports the same functionality as **CoarseTimer** but differs from the **CoarseTimer** class in that **LocalTimer** has a conventional constructor. When an instance of a local timer goes out of scope, the information associated with it is destroyed. Apart from this, all functionality in **LocalTimer** is the same as **CoarseTimer**. The **LocalTimer** class was created to profile individual tasks in applications such as contingency analysis. Each contingency can be profiled separately and the results printed to a separate file. The only functions that are different from the **CoarseTimer** functions are the functions that print out results. The **dumpProfile** functions are not currently supported and the **dump** command has been modified to

```
void dump(boost::shared_ptr<ofstream> stream) const
```

This function requires a stream pointer that signifies which file the data is written to.

## Exceptions

The math module has been implemented so that failures throw exceptions. These can be caught by other parts of code and managed so that code does something more graceful than simply crash when an error is encountered. For example, a calculation that fails because the solver throws an exception might try to run again using a different solver. In a contingency analysis calculation, a

contingency that fails because the solver did not converge can be marked as a failed calculation and the code can proceed to the next contingency. This allows the code to evaluate all contingencies even if some do not converge.

A solver exception can be handled using the following construct

```
LinearSolver solver(*A);  
// User code...  
try {  
    solver.solve(*B,*X);  
} catch (const gridpack::Exception e) {  
    // Do something to manage exception  
}
```

If the solve command fails, it throws a **gridpack::Exception** that can then be managed by the code. This could consist of simply exiting cleanly or the code could try and take corrective action by using a different algorithm.

Exceptions can also be added to error conditions that are detected in user written code so that the error can be picked up in some other part of the application and managed there. Exceptions have two constructors that can be used in applications

```
Exception(const std::string msg)  
Exception(const char* msg)
```

where **msg** is a text string describing the error that was encountered. This message can be read later using the function

```
const char* what()
```

Exceptions are usually created in user code using the following syntax

```
if (...some_condition_is_violated...) {  
    throw gridpack::Exception("Describe error condition");  
}
```

The error message can be printed out to standard out (or standard error) by catching the exception and calling **what**

```
try {  
    // Some action  
} catch (const gridpack::Exception e) {  
    std::cout << e.what() << std::endl;  
    // After printing error take some action  
}
```

```
}
```

### Hash Distribution Module

The hash distribution functionality provides a simple mechanism for quickly distributing data associated with individual buses and branches to the processors that own those buses and branches. This situation can come up in several contexts, particularly when network data is distributed across multiple files. For example, the information on measurements in the state estimation calculation is contained in a file that is distinct from the file that holds the network configuration. The program starts by reading in the network configuration and partitioning it. The program next reads in the measurements, but there is no simple map between the measurements, each of which is associated with either a branch or a bus, and the distributed network. Even if the measurements are read in before the network is distributed, there is still no simple map between measurements and their corresponding buses and branches, since some components may have no measurements associated with them and other components may have multiple measurements. Moving this data to the right processor and providing a simple way of mapping it to the correct bus or branch on that processor is a non-trivial task.

The **HashDistribution** module is a templated class that assumes that the data that is to be sent to the buses and branches are held in user-defined structs. It is contained in the **gridpack::hash\_distr** namespace. The structs used for branches can be different from the structs used for buses. If we designate the bus and branch structs by the names **BusData** and **BranchData** then the constructor for the **HashDistribution** class has the form

```
HashDistribution<MyNetwork, BusData, BranchData>  
(const boost::shared_ptr<MyNetwork> network)
```

Both the **BusData** and **BranchData** structs must be specified when creating a new **HashDistribution** object, even if only bus or branch data is actually being used. If you are just using bus data you can simply repeat the **BusData** type in the branch slot without causing any problems. Similarly, you can also use **BranchData** in both slots if you are only interested in moving data to branches.

The following command can be used to move bus data to the processors that actually hold the corresponding buses

```
void distributeBusValues(std::vector<int> &keys,  
                        std::vector<BusData> &values)
```

The integer array “**keys**” holds the original indices of the buses that the data in the vector “**values**” is supposed to map to. The **keys** and **values** vectors should be the same and the data in the **values** array at index  $n$  should be mapped to the bus indicated by the original index stored at the same location in the **keys** array. This function is collective and must be called on

all processors. The amount of data on each processor does not need to be the same and some processors, or even most of them, can have no data (it is still necessary to call the **distributeBusValues** function across all processors even if some processors contain no data). It is also possible that the same original index can appear multiple times in the keys array, i.e. multiple pieces of data can map to the same bus. On output, the values array contains all the data objects that map to buses on that processor and the keys array contains the *local* indices of the bus. This will include data that maps to ghost buses so a piece of data may map to more than one processor in a distributed system.

An analogous command can be used to distribute data to branches. It has the form

```
void distributeBranchValues(std::vector<std::pair<int,int> > &keys,  
                           std::vector<int> &branch_ids,  
                           std::vector<BranchData> &values)
```

Branches are uniquely identified by the buses at each end of the branch, so the **keys** array in this case is a vector consisting of index pairs representing the original indices of these buses. The **values** array contains the data to be distributed to the branches and the **branch\_ids** array contains the *local* index of the branch on output. Unlike the command to distribute bus values, the **keys** array cannot be reused to store the destination index of the data. Similar to buses, multiple data items can be mapped to the same branch.

### String Utilities

At some point, users may want to develop their own parsers for reading in information in external files. The **StringUtils** class is contained in the **gridpack::utility** namespace and is designed to provide some useful string manipulation routines that can be used to parse individual lines of a file. Other capabilities are available in standard C routines such as **strcmp** and the Boost libraries also have many useful routines. The **StringUtils** class is just a convenient container for different string manipulation methods; it has no internal state.

Some basic routines for modifying strings so that they can be compared with other strings are

```
void trim(std::string &str)
```

which can be used to remove white space at the beginning and end of a string. This function will also convert all tabs and carriage returns to white space before trimming the white space at the ends of the string. The functions

```
void toUpper(std::string &str)  
void toLower(std::string &str)
```

can be used to convert all characters in the string to either upper or lower case.



Many devices in power grid applications are characterized by a one or two character alphanumeric string. It is useful to get these strings into a standard form so that they can be compared with other strings. The function

**std::string clean2Char(std::string &str)**

returns a two character string that is right justified. It will also remove any quotes that may or may not be around the original string. The strings C1, 'C1', "C1" and " C1" will all return a string containing the two characters C1. A single character string will return a two character string with a blank as the first character.

The function

**std::string trimQuotes(std::string &string)**

can be used to remove either single or double quotation marks from around a string and remove any remaining white space at the beginning and end of the string.

Finally, the function

**std::vector<std::string> blankTokenizer(std::string &str)**

will take a string in which individual elements are delimited by blank spaces and return a vector in which each element is a separate string (token). This function treats anything inside the original string that may be delimited by quotes as a single token, even if there are additional blank spaces between the quotes. Thus, the string

**1 5 "ATLANTA 001" 0.00056 1.02**

is broken up into a vector containing the strings

**1  
5  
"ATLANTA 001"  
0.00056  
1.02**

Both single and double quotes can be used as delimiters for internal strings.

### Global Store

The **GlobalStore** class was created to make large amounts of data globally accessible to any processor when replicating the data would be inefficient in terms of the amount of memory required. The premise of the **GlobalStore** class is that processors generate vectors of data and this data is added to a **GlobalStore** object. After all processors have completed adding data, the data is "uploaded" to the **GlobalStore** object so that it is visible to all processors in

the system. Prior to the upload operation, the data is held locally on the processor that generated it. The original motivation for creating this class was to save system state variables that represent the results of individual simulations in a contingency analysis context. These variables could then be used to initialize additional calculations.

The module **GlobalStore** is a templated class that is located in the **gridpack::parallel** namespace. The **GlobalStore** constructor is

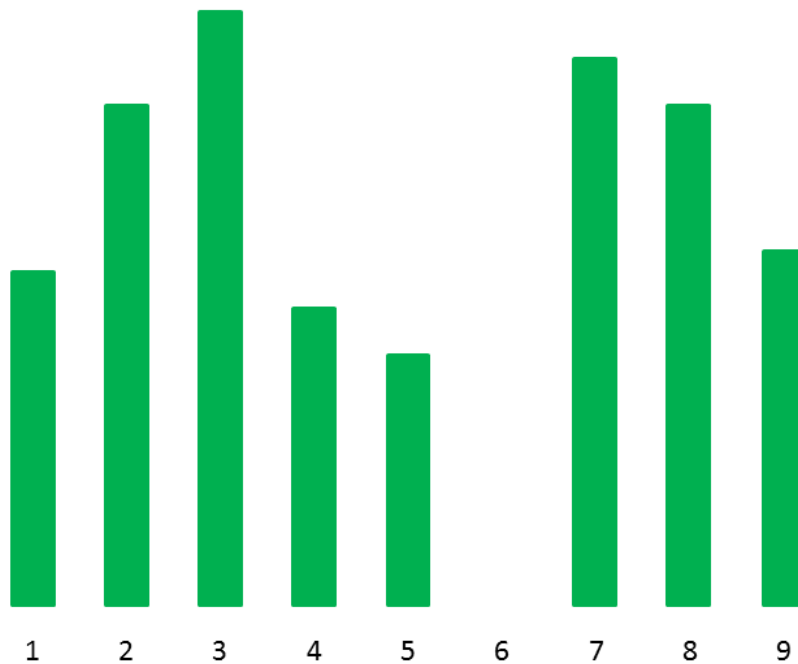
```
GlobalStore<data_type>(const gridpack::parallel::Communicator &comm)
```

The constructor takes a communicator as an argument so data in the **GlobalStore** object will only be visible to processors in the communicator. It also takes the template argument **data\_type** that can be any fixed-sized data type. This includes standard data types such as **int**, **float**, **double**, etc. but could also represent user-defined structs.

Data can be added to the **GlobalStore** object using the command

```
void addVector(const int idx, const std::vector<data_type> &vec)
```

This command assumes that the user has some way of uniquely identifying each contributed vector by an index **idx**. The indices do not have to be complete, i.e. not all indices in some interval  $[0, \dots, N-1]$  need to be added to the storage object, although large gaps between contributed indices are potentially wasteful. The length of the vectors can differ for different indices and there are also no restrictions on which processor contributes which index, so contributions can be made in any order from any processor. The only restriction on indices is that they are not used more than once, i.e. **addVector** is not call more than once on any processor for a given index. This behavior maps fairly well to contingency calculations where the index represents the index of a particular contingency. The data layout in the **GlobalStore** object is illustrated schematically in Figure 11.



**Figure 11.** Schematic diagram of data storage in a GlobalStore object. Vectors can have any length and some indices can be missing data.

Once the processors have completed adding vectors to the storage object, the data is still only stored locally. To make it globally accessible, it is necessary to move it from local buffers to a globally accessible data structure. This is accomplished by calling the function

```
void upload()
```

This function takes no arguments. After calling **upload**, it is no longer possible to continue adding data to the storage object using the **addVector** function.

Once data has been uploaded to the storage object, any processor can retrieve the data associated with a particular index using the function

```
void getVector(const int idx, std::vector<data_type> &vec)
```

This function retrieves the data corresponding to index **idx** from global storage and stores it in a local vector. The **getVector** function can be called an arbitrary number of times after the data has been uploaded. If no data is found, the return vector will have length zero.

One note about using the **getVector** function is worth mentioning. The implementation of the **GlobalStore** class uses some Global Array calls that can potentially interfere with MPI calls in a subsequent function call, resulting in the code hanging. If this occurs, it is usually possible to prevent the hang by calling **Communicator::sync** on the communicator that was used to

define the **GlobalStore** object. This should be done after completing all **getVector** calls but before making calls to other parallel functions.

### Bus Tables

The bus table module was created to allow applications to update the properties of buses over multiple scenarios. This module is designed to read files of the form

```
11002 BL    0.0011 0.0009 0.0018 0.0023
11003 BL    0.2232 0.2113 0.2202 0.2317
11005 BL    0.1188 0.1076 0.1211 0.1197
11008 BL    0.0053 0.0045 0.0067 0.0072
```

The first column is a bus ID, the second column is a one- or two-character tag identifying some device on the bus (e.g. a generator) and the remaining columns are properties of the bus for different scenarios. The columns are delimited by white space. If there are  $N$  columns of properties for the buses then the total number of columns in the file is  $N+2$ , where the extra two columns represent the bus indices and the device tags. The columns containing data are indexed from 0 to  $N-1$ . If the properties apply to the bus as a whole and not some device on the bus, then the tags can be ignored but some arbitrary one- or two-character string still needs to be included in the file for the second column. The scenarios themselves can represent different times, different parameter sets, different loads etc. The properties are assumed to be double precision values. Integer values can be used as properties by storing them as double precision values and then casting them back to integers inside the application. Not all buses need to be included in the table and in many cases, where a device is not present on a bus, it is undesirable to require that each bus be represented.

The **BusTable** module is a templated class that takes the network type as a parameter. It is located in the **gridpack::bus\_table** namespace. The constructor has the form

```
BusTable<MyNetwork>(const boost::shared_ptr<MyNetwork> network)
```

An external file with the format described above can be read in using the function

```
bool readTable(std::string filename)
```

where **filename** points to the appropriate file. This function will ingest the file and store the contents in a distributed form that can be readily access by the application. This function is collective and must be called by all processes over which the network is defined.

Accessing the data in the table can be accomplished using the following three functions

```
void getLocalIndices(std::vector<int> &indices)
void getTags(std::vector<std::string> &tags)
void getValues(int idx, std::vector<double> &values)
```

The first function returns a list of the local bus indices to which the data applies, the second function returns a list of the corresponding device tags and the third function returns the values from column **idx** in the table.

After calling the functions, the data can be applied to the appropriate buses using a loop of the form

```
MyBus *bus;
For (i=0; i<indices.size(); i++) {
    bus = network->getBus(indices[i]).get();
    bus->setProperty(tags[i], values[i]);
}
```

where **setProperty** is a user-defined function in the **MyBus** class that does something useful with the data. This example assumes that the **getLocalIndices** and **getTags** functions have already been called outside the loop.

The number of columns of properties can be accessed using the function

```
int getNumColumns()
```

This function is provided as a method for accessing the total number of scenarios directly from the bus table input, instead of having to include it as a separate parameter.

## Generalize Matrix-Vector Interface

The matrix-vector interface described earlier is suitable for problems where the independent and dependent variables are both associated with buses. However, it does not work for systems where some variables are associated with branches. This can occur in optimization problems such as state estimation, where measurements are made on both buses and branches. Every measurement contributes an equation to the state-estimation optimization, which results in dependent variables associated with branches. To handle these types of problems, a more general approach to creating matrices and vectors is required. This is implemented via the **GenMatVecInterface** class. As illustrated in Figure 5, the **BaseComponent** class directly inherits from this interface, along with the **MatVecInterface**.

Unlike the **MatVecInterface** class, there is no definitive way to map which elements are contributed by a branch or bus, and the number of elements contributed by a branch or bus does not reduce to simple blocks. Thus, the idea that buses and branches contribute simple blocks of data must be abandoned. The **GenMatVecInterface** just assumes that buses and branches contribute some number of equations (dependent variables) to the matrix and that they also contribute some number of independent variables to the matrix. This information is embedded in the function calls

```
virtual int matrixNumRows(void)  
virtual int matrixNumCols(void)
```

These two functions specify how many dependent variables (rows) and how many independent variables (columns) are associated with a bus or branch. For the state estimation module that is currently available in the GridPACK release, the dependent variables are the number of measurements that are associated with the bus or branch and the independent variables are the voltage magnitude and phase angle, which are only associated with buses. Thus, if the state estimation Jacobian is being built, the **matrixNumRows** function returns the number of measurements on each bus and branch. The **matrixNumCols** only returns a non-zero value for buses since the branches have no independent variables. This value is generally 2, if the bus has any measurements associated with it or is attached to a bus or branch that has measurements, otherwise the value is 0. If the bus has measurements and is the reference bus, then the function returns 1. These functions allow the generalized mappers to determine the dimensions of the matrix (for state estimation, the Jacobian is not necessarily square).

Unlike the original matrix-vector interface, the user has to assign the row and column indices to each matrix element. The actual values of these indices are evaluated by the mapper but it is up to the user to take the row index for a particular dependent variable (measurement) and the column index for a particular independent variable (voltage magnitude or phase angle) and pair them with a matrix element (contribution to the Jacobian). The functions that are used for this purpose are

```
virtual void matrixSetRowIndex(int irow, int idx)  
virtual void matrixSetColIndex(int icol, int idx)  
virtual int matrixGetRowIndex(int irow)  
virtual int matrixGetColIndex(int icol)
```

The first two functions are used by the mapper to assign indices for each of the rows and columns contributed by a component. The values of the indices need to be stored in the component so that they can be accessed by other components when evaluating matrix elements. Although these functions are only called by the mapper, they need to be implemented by the user, since multiple matrices may be generated by the application. The variables **irow** and **icol** refer to the list of rows and columns contributed by the component, while the index **idx** is the global index for that row or column in the full matrix. The point of the first two functions is to create a map between the local index of the row or column and the global index of the corresponding row or column in the full matrix. This map is needed because matrix elements constructed on one component may refer to rows or columns on other components. The second pair of functions allow users to recover the global index from the local index.

For example, the state estimation calculation needs to be able to build the Jacobian matrix plus a diagonal matrix that represents the inverse of the uncertainties in all the measurements. The state

estimation components have two modes, **Jacobian\_H** and **R\_inv** for each of these calculations. The **matrixSetRowIndex** method for the buses has the form

```
void SEBus::matrixSetRowIndex(int irow, int idx)
{
    if (p_mode == Jacobian_H) {
        if (irow < p_rowJidx.size()) {
            p_rowJidx[irow] = idx;
        } else {
            p_rowJidx.push_back(idx);
        }
    } else if (p_mode == R_inv) {
        if (irow < p_rowRidx.size()) {
            p_rowRidx[irow] = idx;
        } else {
            p_rowRidx.push_back(idx);
        }
    }
}
```

The row indices for the Jacobian and  $R^{-1}$  are stored in two separate STL arrays **p\_rowJidx** and **p\_rowRidx**. For the state estimation example, the number of rows (for both the Jacobian and  $R^{-1}$ ) is equal to the number of measurements associated with the component. These measurements are held in an internal list in some order. If the number of measurements on the bus is  $M$  then the **irow** index will run from 0,..., $M-1$ , with the **irow** index corresponding to the **irow** element in the list of measurements. The independent variables are also assumed to be ordered in some fashion. Again, for the state estimation example, the phase angle is indexed by 0 and the voltage magnitude is indexed by 1.

The function for accessing the row indices is implemented as

```
int gridpack::state_estimation::SEBus::matrixGetRowIndex(int idx)
{
    if (p_mode == Jacobian_H) {
        return p_rowJidx[idx];
    } else if (p_mode == R_inv) {
        return p_rowRidx[idx];
    }
}
```

Again, depending on the mode, this function will return different values and for this reason, these functions need to be implemented by the user. They cannot be implemented as part of the framework because the number of modes is application-specific and controlled by the developer.

The functions that are used to actually evaluate matrix elements are

```
virtual int matrixNumValues(void) const  
virtual void matrixGetValues(ComplexType *values,  
int *rows, int *cols)
```

The first function returns the total number of matrix elements that will be evaluated by the component. This is used inside the mapper to allocate arrays that hold matrix elements coming from the components. The second function is used to evaluate actual matrix elements, along with their row and column indices. The real-valued version of **matrixGetValues** replaces **ComplexType** with **double**. This function is the one that will make use of the **matrixGetRowIndex** and **matrixGetColIndex** functions. The evaluation of the **matrixNumValues** function can be quite complicated. For the state estimation Jacobian matrix, the number of matrix elements contributed by a component depends on the number of measurements associated with that component and the number of variables that couple to that measurement. A measurement on a bus will usually contribute two values for the independent variables on the bus, plus an additional two values for each bus that is attached to the center bus via a branch. This number will be modified slightly if one of the buses in this group is a reference bus. For branches, the number of matrix elements contributed by each measurement is approximately four, two elements for each bus at either end of the branch. This number may drop if one of the buses is a reference bus.

The **matrixGetValues** function is used to evaluate each of the matrix elements. It also gets the matrix indices for this element from the appropriate network component. The number of matrix elements returned by this function must correspond to the number returned by the **matrixNumValues** function. To see how the assignment of the indices works, we can look at the matrix element of the Jacobian corresponding to the gradient of a real power injection measurement  $P_i$  on bus  $i$  with respect to the phase angle on another bus  $j$  that is connected to  $i$  via a single branch. The contribution to the Jacobian from this measurement is given by the formula

$$\frac{\partial P_i}{\partial \theta_j} = V_i V_j (G_{ij} \sin(\theta_i - \theta_j) - B_{ij} \cos(\theta_i - \theta_j))$$

Suppose  $P_i$  is measurement  $k$  on the bus. Then the row index **im** for this matrix element can be evaluated by calling the function

```
im = matrixGetRowIndex(k);
```



The column index is associated with the phase angle variable on the remote bus  $j$ . Assuming that a pointer (**bus\_j**) to the remote bus is already available, then the column index **jm** for this matrix element could be obtained by calling

```
jm = bus_j->matrixGetColIndex(0);
```

This function is called with the argument 0 since the dependent variables are always ordered as phase angle (0) followed by voltage magnitude (1). The full list of Jacobian matrix elements can be obtained by looping over all measurements. For each bus measurement, there are contributions from the dependent variables on each connected bus plus two contributions from the calling bus. Similarly, for each branch measurement there are approximately four contributions coming from the independent variables associated with the buses at each end of the branch. A simple counter variable can be used to make sure that the matrix element value and the corresponding row and column indices stored in the same location of the **values**, **rows** and **cols** arrays that are returned by the **getMatrixValues** function.

The **GenMatVecInterface** also includes functions for setting up vectors. These work in a very similar way to the generalized matrix functions, so they will only be described briefly. The two functions

```
virtual void vectorSetElementIndex(int ielem, int idx)  
virtual void vectorGetElementIndices(int *idx)
```

can be used to set and retrieve vector indices. The index **ielem** is the local index within the element while **idx** is the global index within the distributed vector. In this case it is usually more convenient to get all indices associate with a component at once, so the **vectorGetElementIndices** returns an array instead of a single value. The function

```
virtual int vectorNumElements() const
```

returns the number of vector elements contributed by a component and the function

```
virtual void vectorGetElementValues(ComplexType *values, int *idx)
```

returns a list of the values along with their global indices. For real vectors, replace the **ComplexType** array with an array of type **double**. Again, the index value can be obtained by first calling the **vectorGetElementIndices** function and using this to obtain the correct index for each element.

The vector interface includes one additional function that does not have a counterpart in the matrix interface. This is the function

```
virtual void vectorSetElementValues(ComplexType *values)
```

This function can be used to push values from a solution vector back into the network components. The values are ordered in the same way as the values in the corresponding **vectorGetElementValues** call, so it is possible to unpack them and assign them to the correct internal variables for each component. This function is analogous to the **setValues** call in the regular **MatVecInterface**.

The functions in the **GenMatVecInterface** are invoked in the generalized mappers. These reside in the **GenMatrixMap** and **GenVectorMap** classes. Like the standard mappers, these classes are relatively simple and contain only a few methods. The **GenMatrixMap** class consists of the constructor

```
GenMatrixMap<MyNetwork>(boost::shared_ptr<MyNetwork> network)
```

and the methods

```
boost::shared_ptr<gridpack::math::Matrix> mapToMatrix(void)  
void mapToMatrix(boost::shared_ptr<gridpack::math::Matrix> matrix)  
void mapToMatrix(gridpack::math::Matrix &matrix)  
void overwriteMatrix(boost::shared_ptr<gridpack::math::Matrix> matrix)  
void overwriteMatrix(gridpack::math::Matrix &matrix)  
void incrementMatrix(boost::shared_ptr<gridpack::math::Matrix> matrix)  
void incrementMatrix(gridpack::math::Matrix &matrix)
```

These functions all have the same behaviors as the analogous functions in the standard **FullMatrixMap**. The **GenVectorMap** class has the constructor

```
GenVectorMap<MyNetwork>(boost::shared_ptr<MyNetwork> network)
```

and supports the methods

```
boost::shared_ptr<gridpack::math::Vector> mapToVector(void)  
void mapToVector(boost::shared_ptr<gridpack::math::Vector> &vector)  
void mapToVector(gridpack::math::Vector &vector)
```

These functions have the same interpretations as the analogous functions in the **BusVectorMap** class. A new function is

```
mapToNetwork(boost::shared_ptr<gridpack::math::Vector> &vector)
```

which can be used to push data from a vector back into the network components (both buses and branches).

## Generalized Slab Mapper

The generalized slab mapper also uses functions in the generalized matrix-vector interface to build dense matrices. These matrices are dense since they are generated by taking a typical vector that corresponds to a set of variables on the buses and branches and replicating the vector for different values of the variables. An example would be a matrix formed from a time series of values for a set of variables on the buses and branches. One set of indices for the matrix corresponds to the set of variables and the other set of indices corresponds to the time series. In a certain sense, these matrices are “fat” vectors since instead of each variable having only one value, they have multiple values. In general, slab matrices are not square. The slab matrices are used in the Kalman filter application, but they may have applicability elsewhere.

The slab mappers use additional functions from the **GenMatVecInterface** in order to construct matrices. These functions are analogous to the functions for setting up vectors using the **GenVectorMap**. The main difference is that instead of describing a list of values, the functions describe a matrix block. The row dimension corresponds to a list of variables and the column dimension describes the number of values taken by each variable. The column dimension must be the same across all variables. The contribution to the matrix from each network component is given by the function

```
void slabSize(int *rows, int *cols) const
```

The index for each row can be stored using the function

```
void slabSetRowIndex(int irow, int idx)
```

This function is called by the mapper and is analogous to the **vectorSetElementIndex** function. For the slab matrices, there is no corresponding call for columns since the matrices are dense and all rows have the same number of non-zero columns. The indices can be retrieved by the function

```
void slabGetRowIndices(int *idx)
```

which is similar to the **vectorGetElementIndices** function.

## Optimization

GridPACK supports optimization via an interface that can be applied to bus and branch components, as well as wrappers to some common optimization libraries. At present, there are no example problems available for the optimization capability in GridPACK.

The optimization interface can be optionally inherited by the bus and branch classes. Unlike the other interfaces described in this document, this is not already included in the **BaseBusComponent** and **BaseBranchComponent** classes. A component that wants to use the optimization interface must explicitly inherit the interface when declaring the class. The

optimization interface is built up around the **Expression** class that is designed to provide some functional programming capability to GridPACK. This functionality matches the interfaces to many optimization packages. Additional classes needed by the optimization interface are the **Variable** class and the **Constraint** class.

The **Variable**, **Expression** and **Constraint** classes are defined in the **gridpack::optimization** namespace. A number of different types of variables inherit from **Variable**. This includes **RealVariable**, **IntegerVariable** and **BinaryVariable**. Variables can take on different values and are designed to be used as parts of expressions. Expressions, in turn, can be composed of variables and can also be composed of other expressions. This allows users to build up quite complicated functions. An example of an expression is the following

```
typedef boost::shared_ptr<Variable> VarPtr;
typedef boost::shared_ptr<Expression> ExpPtr;

VarPtr x;
ExpPtr f;
x.reset(new RealVariable(0.0));
ExpPtr a(new RealConstant(5.0));
ExpPtr b(new RealConstant(-2.0));
f = a*x + b;
```

This code fragment defines the function **f**. Note that the last line is not an assignment in the conventional sense; **f** does not represent the value computed from the current values of **a**, **x**, and **b** but rather the operation of multiplying **x** by the constant value **a** and then adding the constant value **b** to the result.

The functions in the optimization interface are all expressed in terms of the **Expression** and **Variable** classes. Apart from simple constructors and destructors, there are only three functions in this interface. The first of these returns a list of all variables associated with a particular bus or branch.

```
std::vector<boost::shared_ptr<Variable> > getVariables()
```

Note that the variables returned by a network component can be used in expressions generated by other components but all variables in the problem should be returned by one component somewhere in the network and by no more than one component. This function is called by the optimizer to get a complete listing of variables. Ghost buses and branches do not call this function.

The second function is used to determine the constraints in the system. A constraint is a relation between two expressions. Allowed relations are `<=`, `>=` and `==`. The function

```
std::vector<boost::shared_ptr<Constraint> > getLocalConstraints()
```

returns a list of constraints that are associated with a particular network component. This function assumes that all returned constraints can be evaluated by the associated network component. Constraints that are functions of variables on the component plus variables on attached components fall into this category. Simple constraints defined as being numerical bounds on the value of a variable can be incorporated into the variables definitions. The **Constraint** class is designed to handle more complicated constraints that are functions of one or more variables.

The final function in the optimization interface is

```
boost::share_ptr<Expression> getObjectiveFunction()
```

This function assumes that the objective function for the entire system is a sum of terms contributed by each network component. Buses are branches that contribute nothing can return a null pointer. These three functions can be used to define the complete optimization problem. The remainder of this section will expand on the different classes used to implement these functions.

The **Variable** class is used to define different types of variables. Variables can have real values, integer values or they can be binary switches (0 and 1). All types of variables are derived from the **Variable** base class. The **RealVariable** class has two constructors

```
RealVariable(double val)
```

```
RealVariable(double val, double vmin, double vmax)
```

The first constructor initializes the variable with the value **val**, the second constructor initializes the variable with the value **val** and applies the constraints

```
val >= vmin
```

```
val <= vmax
```

to the variable. These constraints could also be applied using the **Constraint** class, but this is obviously simpler, if applicable. The variable can be named using the method

```
void name(const std::string name)
```

The name of the variable can be recovered by the user with the function

```
std::string name() const
```

The name is an important property of the variable. It allows the application to identify two variables as representing the same quantity if they are located on different network components. This can happen with ghost buses and branches, where multiple copies of the same bus or branch may exist in the system. To make sure that the application can identify variables from different copies of the same network components as being the same, the variables defined within the copies must all have the same names. Conversely, variables from physically distinct network components need to have unique names that distinguish them from each other.

All variables have an implicit upper and lower bound, even if this is not explicitly set using the constructor. If only the initial value of the variable is set by the constructor, then the upper and low bounds are set internally to machine limited values. The actual limits of the variable, along with the initial value, can be recovered with the functions

```
double initial() const
double lowerBound() const
double upperBound() const
```

The **IntegerVariable** class is similar to the **RealVariable** class, except that only integer values of the variable are allowed. It has the two constructors

```
IntegerVariable(int val)
IntegerVariable(int val, int vmin, int vmax)
```

The behavior of this class is similar to that of **RealVariable**, except all arguments are integers. Similar to the **RealVariable** class, the **IntegerVariable** class supports the functions

```
void name(const std::string name)
std::string name() const
int initial() const
int lowerBound() const
int upperBound() const
```

Finally, the **BinaryVariable** class has the single constructor

```
BinaryVariable(int val)
```

The **BinaryVariable** class inherits from the **IntegerVariable** class and has lower and upper bounds of 0 and 1. It can only take on these two values. Similar to the real and integer variable classes, it has the methods

```
void name(const std::string name)
std::string name() const
int initial() const
int lowerBound() const
```

```
int upperBound() const
```

In this case the last two methods are redundant, since the return values will always be 0 and 1, respectively.

Variables can be used to create expressions. Expressions can consist of a single variable, a constant, or variables and expressions combined together using unary or binary functions. The expression class is designed to support the relatively simple sums over polynomials that are encountered in power grid applications, but even with this limited scope, quite a few complicated functions can be created with it. Expressions never use variables directly, they are composed using variable pointers, so all variables in an expression should be instantiated using a `boost::share_ptr<Variable>` construct. Expressions themselves should also be instantiated using a `boost::share_ptr<Expression>` construct. To create a simple quadratic polynomial requires the code

```
VarPtr x;  
ExpPtr a(new RealConstant(1.0));  
ExpPtr b(new RealConstant(-2.0));  
ExpPtr c(new RealConstant(3.0));  
x.reset(new RealVariable(0.0, -10.0, 10.0));  
ExpPtr x2 = x^2;  
ExpPtr p = a+b*x+c*x2;
```

The exponentiation operation in the `Expression` class does not have the correct precedence behavior so it is necessary to define the `Expression x2` and then using it to define `p`. Expressions can be combined using the binary operations

```
ExpPtr+ExpPtr  
ExpPtr-ExpPtr  
ExpPtr*ExpPtr  
ExpPtr/ExpPtr
```

Expressions can also be used in the functions

```
ExpPtr^int  
sin(ExpPtr)  
cos(ExpPtr)
```

More complicated expressions can be built up from these simple operations. A constant expression can be created using the constructs

```
ExpPtr var(new RealConstant(double));  
ExpPtr var(new IntegerConstant(int));
```

These can then be used in other expressions as parameters.

Finally, constraints can be created using constructs such as

```
boost::shared_ptr<Constraint> const(ExpPtr1 relation constant);
```

where the character string relation can be

```
">="  
"<="  
"=="
```

and constant is a **double** or **int** value.

## Application Modules

Many of the example applications in GridPACK have been converted to modules that can be called from other programs. These modules make it relatively simple to chain different types of calculations together to form larger applications. An example is using power flow or state estimation to initialize a dynamic simulation. The modules are designed to separate the major phases of the calculation into calls so that users have some fine-grained control that allows them to mix different applications together. In most cases, different options for setting up calculations are provided so that once a network has been read in and partitioned, it is not necessary to repeat this process when a new calculation is started based on the results of a previous simulation.

Currently, three applications are available as modules within GridPACK. They include power flow, state estimation, and dynamic simulation using the full Y-matrix. Each of these modules can be used to create a short, standalone application, but the goal is to enable users to combine modules together in more complicated work flows. These modules can also be used as a starting point for users to create their own applications by modifying the existing code in the modules to create new functionality. Each of the modules is described in more detail below. Example codes that use the modules to implement applications can be found in the **src/application** directory. These include powerflow, state estimation, contingency analysis and dynamic simulation. These directories also contain sample input networks and input files. Options for different solvers can be found in these files.

### Power Flow

The power flow module consists of a collection of function calls that can be used to set up and run power flow calculations. Additional routines are designed to support different types of contingency analysis. The power flow application class is **PFAppModule** and belongs to the **gridpack::powerflow** namespace. The constructor and destructor for this class are simple and only create the basis power flow object. In particular, the power flow network must be created outside the power flow object and then assigned to the object when the network configuration file is read in. This can be done with the call



```
void readNetwork(boost::shared_ptr<PFNetwork> &network,  
                Configuration *config)
```

The **Configuration** object should already be pointing to an open file containing a **Powerflow** block. This block contains a **networkConfiguration** field that has the name of the PSS/E format file containing the network information. The network configuration file is read directly from the input deck by the **readNetwork** method. The **PFNetwork** is defined in the the **gridpack.hpp** header file. The configuration module is usually opened in the main calling program and a pointer to the file can be passed through to power flow module. The **readNetwork** routine also partitions the network.

Once the network has been read in, the internal indices and exchange buffers can be set up by calling

```
void initialize()
```

The power flow application is now ready to be used. To solve the current configuration, the calls

```
void solve()
```

```
void nl_solve()
```

can be used. The first call solves the system uses a hand coded Newton-Raphson iteration loop to solve the system, the second call uses a non-linear solver to solve the power flow equations. Both solvers can be controlled through solver options in the input file. The type of linear solver used in the solve routine is controlled by the parameters in the **LinearSolver** block, the non-linear solver is controlled by the properties in the **NonlinearSolver** block

Output from the power flow solution can be written to an output file or standard out using one of the commands

```
void write()
```

```
void writeBus(const char* signal)
```

```
void writeBranch(const char* signal)
```

The first command writes out the real and imaginary parts of the complex power for the branches and the voltage magnitude and phase angle for the buses. The second command only writes out bus properties. If no argument is given, the command writes out the voltage magnitude and phase angle for every bus. For buses, the argument “**pq**” writes out the real and imaginary parts of the complex voltage and “**record**” writes out the type of bus, the total active and reactive constant power loads, and the total active and reactive generator power outputs. For branches, “**flow**” writes out the real and imaginary parts of the complex power and “**record**” writes out the values of the resistance, reactance, charging and A, B, C ratings for each line element.

Additional information can be written to standard out or a file using the command

```
void print(const char* buf)
```

which writes out the contents of the character array **buf**. This command can be called from all processors, but only one processor actually writes out data.

The location of output can be controlled using the commands

```
void open(const char* filename)
```

```
void close()
```

If the write commands or **print** are used without calling **open**, then all output is directed to standard out. If **open** is called, then the output is directed to the file specified in filename until the **close** command is called, after which all output is again directed towards standard out.

If the results of the power flow calculation are needed by another calculation, then the voltage magnitude and phase angle of the bus and the real and imaginary parts of the complex power for each generator can be stored in the **DataCollection** objects on each bus using the command

```
void saveData()
```

If the network is then copied to a new type of network, this information is carried over to the new network. The voltage magnitude and phase angle is stored in the **DataCollection** variables **BUS\_PF\_VMAG** and **BUS\_PF\_VANG** and the generator parameters are stored in the indexed variables **GENERATOR\_PF\_PGEN[i]** and **GENERATOR\_PF\_QGEN[i]**, where the index **i** runs over all generators on the bus.

The remaining methods in the **PFAppModule** class support different kinds of contingency applications. Contingencies are defined using the data structure

```
struct Contingency
{
    int p_type;
    std::string p_name;
    // Line contingencies
    std::vector<int> p_from;
    std::vector<int> p_to;
    std::vector<std::string> p_ckt;
    // Status of line before contingency
    std::vector<bool> p_saveLineStatus;
    // Generator contingencies
    std::vector<int> p_busid;
}
```

```

std::vector<std::string> p_genid;
// Status of generator before contingency
std::vector<bool> p_saveGenStatus;
};

```

The variable **p\_type** corresponds to an enumerated type that can have the values **Generator** and **Branch**. The variables **p\_saveLinesStatus** and **p\_saveGenStatus** are used internally and do not have to be set by the user. The remaining variables are used to describe the lines and generators that may fail during a contingency event. These variables are all vectors, since a single contingency could theoretically represent the failure of multiple elements. For failures of type **Branch**, the variables **p\_from** and **p\_to** are the original indices of the “from” and “to” bus that identify a branch and the variable **p\_ckt** is the 2 character identifier of the individual transmission element. For failures of type **Generator**, **p\_busid** is the original index of the bus and **p\_genid** is the 2 character identifier of the generator that fails. An example of how to use this functionality is given in the contingency analysis application that can be found under **src/applications/contingency\_analysis**. This is also a good example of how to use modules.

Two calls

```

bool setContingency(Contingency &event)
bool unsetContingency(Contingency &event)

```

can be used to set or unset a contingency. The call **unsetContingency** should only be called after calling **setContingency** and it should use the same **event** argument. After calling the **unsetContingency** method, the network should have the same configuration as before calling the **setContingency** method.

The remaining calls in **PFAppModule** can be used to determine the status of a network after solving a configuration with a contingency. The functions

```

bool checkVoltageViolations(double Vmin, double Vmax)
bool checkVoltageViolations(int area, double Vmin, double Vmax)

```

can be used to check for a voltage violation anywhere in the system where **Vmin** and **Vmax** are the minimum and maximum allowable voltage excursions (per unit). The second function only checks for violations on buses with the specified value of **area**. These functions are true if there are no voltage violations and return false if a violation is found on one or more buses. It frequently turns out that many networks have voltage violations even in the absence of any contingencies and it is often desirable to ignore these violations. This can be accomplished using the function

**void ignoreVoltageViolations(double Vmin, double Vmax)**

If this function is called after solving the power flow system in the absence of any contingencies, then buses that contain violations will be ignored in subsequent checks of violations. These settings can be undone by calling

**void clearVoltageViolations()**

Line overload violations can be checked by calling one of the functions

**bool checkLineOverloadViolations()**

**bool checkLineOverloadViolations(int area)**

The limits on the line are contained in parameters read in from the network configuration file so these functions have no arguments describing the line limits. The second function will only check for violations on lines with the specified value of **area**. Like voltage violations, branches that display line overload violations that are present even without contingencies can be ignored in the checks by calling the function

**void ignoreLineOverloadViolations()**

after running a calculation on the system without contingencies. These settings can be cleared using the function

**void clearLineOverloadViolations()**

Finally, the internal voltage variables that are used as the solution variables in the power flow calculation can be reset to their original values (specified in the network configuration file) by calling the function

**void resetVoltages()**

Again, this may be useful in contingency calculations where multiple calculations are run on the same network and it is desirable that they all start with the same initial condition.

### State Estimation Module

The state estimation module can be used to set up and run a state estimation calculation. It does not have the extra functions that the power flow module contains for supporting contingency analysis, so the interface is a bit smaller. In addition to a standard network configuration file, the state estimation calculation needs a second file consisting of measurements. This file has the format

**<Measurements>**

**<Measurement>**

**<Type>VM</Type>**

```

    <Bus>1</Bus>
    <Value>1.0600</Value>
    <Deviation>0.0050</Deviation>
</Measurement>
<Measurement>
    <Type>PIJ</Type>
    <FromBus>1</FromBus>
    <ToBus>2</ToBus>
    <CKT>BL</CKT>
    <Value>1.5688</Value>
    <Deviation>0.0100</Deviation>
</Measurement>
<Measurement>
    <Type>QIJ</Type>
    <FromBus>1</FromBus>
    <ToBus>2</ToBus>
    <CKT>BL</CKT>
    <Value>-0.2040</Value>
    <Deviation>0.0100</Deviation>
</Measurement>
<Measurement>
    <Type>PI</Type>
    <Bus>1</Bus>
    <Value>2.3240</Value>
    <Deviation>0.0100</Deviation>
</Measurement>
<Measurement>
    <Type>QI</Type>
    <Bus>1</Bus>
    <Value>-0.1690</Value>
    <Deviation>0.0100</Deviation>
</Measurement>
</Measurements>

```

for the five types of measurements **VM**, **PIJ**, **QIJ**, **PI**, and **PJ**. Measurements can appear on any element of the network and multiple measurements are allowed on each element. The state estimation module does not have any error checking ability to determine if there are sufficient measurements to guarantee solvability, if not enough measurements are available then the calculation will simply crash or fail to converge.

The state estimation module is represented by the **SEAppModule** class which is in the **gridpack::state\_estimation** namespace. The **gridpack.hpp** file contains a definition for the state estimation network **SENetwork**. After instantiating an **SEAppModule** object and a shared pointer to an **SENetwork**, the state estimation calculation can read in an external network configuration file using the command

```
void readNetwork(boost::shared_ptr<SENetwork> &network,  
    gridpack::utility::Configuration *config)
```

The **Configuration** object should already be pointing at an open file containing a **State\_estimation** block. Inside the **State\_estimation** block there should be a **networkConfiguration** field containing the name of the network configuration file. The file name is parsed directly inside the **readNetwork** method and does not need to be handled by the user.

Alternatively, the **SENetwork** object may have already been cloned from an existing network and therefore there is no need to read in the configuration from an external file and partition it across processors. In this case, the **SEAppModule** can be assigned the network using the command

```
void setNetwork(boost::shared_ptr<SENetwork> &network,  
    gridpack::utility::Configuration *config)
```

This function just assigns the existing network to an internal pointer, as well as a pointer to the input file. It is much more efficient than reading in the network configuration file, if the network already exists. This can occur when different types of calculations are being chained together.

Once a network is in place and has been properly distributed, the measurements can be read in by calling the function

```
void readMeasurements()
```

The name of the measurement file is in the input deck and a pointer to this file has already been internally cached in the **SEAppModule** when the network was assigned. The measurement file name is stored in the **measurementList** field within the **State\_estimation** block.

The network object can be initialized and the exchange buffers set up by calling the

```
void initialize()
```

method followed by

```
void solve()
```

to obtain the solution to the system. Results can be written out to standard out using the method

**void write()**

This function will write out the voltage magnitude and phase angle for each bus and the real and imaginary parts of the reactive power for each branch. In addition, it will print out a comparison of the calculated value and the original measured value for all measurements.

Finally, the results of the state estimation calculation can be saved to the **DataCollection** object assigned to the buses by calling the

**void saveData()**

method. The voltage magnitude and phase angle are stored as the variables **BUS\_SE\_VMAG** and **BUS\_SE\_VANG** and the generator parameters are stored as the indexed variables **GENERATOR\_SE\_PGEN[i]** and **GENERATOR\_SE\_QGEN[i]**, where **i** runs over the set of generators on the bus.

### Dynamic Simulation Module using Full Y-Matrix

GridPACK supplies a dynamic simulation module that integrates the equations of motion using an algorithm based on inversion of the full Y-matrix. This module has been designed to enable the addition of generator models that extend beyond the classical generator. It also supports exciters, governors, relays and dynamic loads. Models that are currently available include

Generators:

**GENCLS**  
**GENSAL**  
**GENROU**

Exciters:

**EXDC1**  
**ESST1A**

Governors:

**WSIEG1**  
**WSHYGP**

Relays:

**LVSHBL**  
**FRQTPAT**  
**DISTR1**

Dynamic Loads:

**ACMTBLU1**  
**IEEL**  
**MOTORW**  
**CIM6BL**

The full Y-matrix implementation of dynamic simulation is represented by the **DSFullApp** class and the **DSFullNetwork**, both of which reside in the **gridpack::dynamic\_simulation** namespace.

The dynamic simulation module uses an input deck of the form

```
<?xml version="1.0" encoding="utf-8"?>
<Configuration>
  <Dynamic_simulation>
    <networkConfiguration>IEEE_145.raw</networkConfiguration>
    <generatorParameters>IEEE_145.dyr</generatorParameters>
    <simulationTime>30</simulationTime>
    <timeStep>0.005</timeStep>
    <faultEvents>
      <faultEvent>
        <beginFault> 2.00</beginFault>
        <endFault> 2.05</endFault>
        <faultBranch>6 7</faultBranch>
        <timeStep> 0.005</timeStep>
      </faultEvent>
    </faultEvents>
    <generatorWatch>
      <generator>
        <busID> 60 </busID>
        <generatorID> 1 </generatorID>
      </generator>
      <generator>
        <busID> 112 </busID>
        <generatorID> 1 </generatorID>
      </generator>
    </generatorWatch>
    <generatorWatchFrequency> 1 </generatorWatchFrequency>
    <generatorWatchFileName>gen_watch.csv</generatorWatchFileName>
    <LinearMatrixSolver>
```



```

    <PETScOptions>
      -ksp_atol 1.0e-18
      -ksp_rtol 1.0e-10
      -ksp_monitor
      -ksp_max_it 200
      -ksp_view
    </PETScOptions>
  </LinearMatrixSolver>
</Dynamic_simulation>
</Configuration>

```

The input for dynamic simulation module is contained in the **Dynamic\_simulation** block. Two features are important, the blocks describing faults and the blocks describing monitored generators. Faults are described in the **faultEvents** block. The code currently only handles faults on branches. Inside the **faultEvents** block are individual faults, described by a **faultEvent** block. Multiple **faultEvent** blocks can be contained within the **faultEvents** block. As will be described below, it is possible for the faults to be listed in a separate file. This can be convenient for describing a task-based calculation that may contain a lot of faults. The parameters describing the fault include the time (in seconds) that the fault is initiated, the time that it is terminated, the timestep used while integrating the fault and the indices of the two buses at either end of the fault branch.

When running a dynamic simulation, it is generally desirable to monitor the behavior of a few generators in the system and this can be done by setting generator watch parameters. The **generatorWatch** block specifies which generators are to be monitored. Each generator is described within a **generator** block that contains the index of the bus that the generator is located on and the character string ID of the generator. The results of monitoring the generator are written to the file listed in the **generatorWatchFileName** field and the frequency for storing generator parameters in this file is set in the **generatorWatchFrequency** field. This parameter describes the time step interval for writing results (an integer), not the actual time interval.

Before using the dynamic simulation module, a network needs to be instantiated outside the **DSFullApp** and then passed into the module. If the module itself is going to read and partition a network, then it should use the function

```

void readNetwork(boost::shared_ptr<DSFullNetwork> &network,
  gridpack::utility::Configuration *config,
  const char *otherfile = NULL)

```

The **Configuration** object should already be pointing to an input deck with a **Dynamic\_simulation** block that specifies the network configuration file. The optional **otherfile** argument in **readNetwork** can be used to overwrite the **networkConfiguration** field in the input deck with a different filename. This capability has proven useful in some contingency applications where multiple PSS/E files need to be read.

Alternatively, a distributed network may already exist (it may have been cloned from another calculation). In that case, the function

```
void setNetwork(boost::shared_ptr<DSFullNetwork> &network,  
               gridpack::utility::Configuration *config)
```

can be used to assign an internal pointer to the network. Again, the **Configuration** object should already be pointing to an input file.

Additional generator parameters can be assigned to the generators by calling the function

```
void readGenerators()
```

This function opens the file specified in the **generatorParameters** field in the input file and reads the additional generator parameters. The file is assumed to correspond to the PSS/E .dyr format. The devices listed at the start of this section can be included in this file.

After setting up the network and reading in generator parameters, the module can be initialized by calling

```
void initialize()
```

This sets up internal parameters and initializes the network so that it is ready for calculations.

A list of faults can be generated from the input file by calling

```
std::vector<gridpack::dynamic_simulation::DSFullBranch::Event>  
getFaults(gridpack::utility::Configuration::CursorPtr cursor)
```

If the cursor variable is pointed at a **Dynamic\_simulation** block inside the input file (as in the example input block above) then this function will return a list of faults from the input deck. However, it is also possible that the cursor could be pointed to the contents of another file. As long as it is pointed to a block containing a **faultEvents** block, this function will return a list of faults. This allows users to declare a large list of faults in a separate file and then access the list by including the external file name as a parameter in the input deck of their application.

The monitoring of generators specified in the input deck can be set up by calling

```
void setGeneratorWatch()
```

This will guarantee that all generators specified in the input deck are monitored and that the results are written out to the specified file. If this function is not called, the generator watch parameters in the input file are ignored.

Simulations can be run using the function

```
void solve(gridpack::dynamic_simulation::DSFullBranch::Event fault)
```

Some additional results can be written at the end of the simulation using the function

```
void write(const char *signal)
```

The signal parameter can be used to control which results are written out. This function currently does not support any output. All output results are controlled using the generator watch parameters.

Some additional functions can be used to control where output generated during the course of a simulation is directed. The following two functions can be used to direct output from the **write** function to a file

```
void open(const char* filename)
```

```
void close()
```

The function

```
void print(const char* buf)
```

can be used to print out a string to standard out. If the **open** function has been used to open a file, then the output is directed to the file. This function is equivalent to the **header** convenience function in the serial IO classes.

Additional functions can be used to stored data from the generator watch variables. These can be used to save the time series data from a simulation in a collection of vectors. The application can then use these series in whatever way it wants. There are four functions that enable this capility. The first is

```
void saveTimeSeries(bool flag)
```

This function must be called with the argument set to “true” in order for the time series data to be saved. Otherwise it is only written to output and no data is saved between time steps. The second function can be called after the solve function has been called and the simulation is completed. It returns a vector of time series

```
std::vector<std::vector<double> > getGeneratorTimeSeries()
```

This function returns a vector containing the time series data for all the watched generators located *on this processor* (generators on buses owned by neighboring processors are not included).

To find out which variables are actually in the list returned by **getGeneratorTimeSeries** requires the remaining two functions. The function

```
void getListWatchedGenerators(std::vector<int> &bus_ids,  
    std::vector<std::string> &gen_ids)
```

returns a list of the bus IDs and 2-character generator tags for all monitored generators. In particular, it assigns an ordering to these generators that is used by function

```
std::vector<int> getTimeSeriesMap()
```

This function returns a map between the elements in the list of time series returned by **getGeneratorTimeSeries** and the generators that those time series correspond to. For example suppose the time series list has four elements in it that happen to correspond to two generators on processor. There are a total of six monitored generators in the system. The vectors returned by **getListWatchedGenerators** have length six, the vector returned by **getTimeSeriesMap** has length four. The value in the map vector for the corresponding element in the time series vector points to the location of the bus index and generator tag for that time series variable in the lists returned by **getListWatchedGenerators**. This still leaves it up to the user to identify the actual variable being watched within the generator. In this example there are four variables that are watched but only two generators. Currently, the generator watch capability only watches the rotor speed and rotor angle of each generator. The first time series is the speed and the second time series is the angle.

### Kalman Filter

GridPACK includes a Kalman filter module that can be used for dynamic state estimation. The Kalman filter relies heavily on parallel matrix multiplies that are not currently very high performing, so users will probably find this module too slow for large grids. However, we include it for users interested in exploring the use of Kalman filters in smaller applications. We hope to improve performance in future releases.

The current implementation of the Kalman filter only supports classical generators. These are described in a PSS/E .dyr formatted file. The network itself can be described using a standard PSS/E .raw file. In addition to the .raw and .dyr files, users need to supply times series data for the voltage magnitude and voltage phase angle on all buses. These are stored as .csv files. The format for both the voltage magnitude and phase angle files is

```
t-3001, Bus-1, Bus-2,...  
0.0, -0.001, -0.135,...
```

```

0.1,      -0.001, -0.135,...
0.2,      -0.001, -0.135,...
:

```

All entries on the same lines are separated by commas. The first row contains the name of all columns. The first column is time and has a name of the form **t-xxx**, where **xxx** is an integer representing the number of time steps in the file. The number of rows in the file corresponds to **xxx+1** (the extra row is the first line with the column names). The number of columns is equal to the number of buses in the file plus one (the extra column contains the times). After the first column, the remaining names are all of the form **Bus-xxx**, where **xxx** is an integer representing the bus ID. The remaining rows contain the time of the measurement and the value for the measurement on each of the buses.

The input file for the Kalman filter module used both for a dynamic simulation as well as input that is unique to the Kalman filter module. The dynamic simulation parameters that are used include

```

<Dynamic_simulation>
  <simulationTime>3</simulationTime>
  <timeStep>0.01</timeStep>
  <!-- = 1 Fault Event is known;
        = 0 Fault event is unknown, switch is skipped.
  -->
  <KnownFault> 1 </KnownFault>
  <TimeOffset> 0 </TimeOffset> <!--skip initial measurement data -->
  <faultEvents>
    <faultEvent>
      <beginFault> 1 </beginFault>
      <endFault> 1.1</endFault>
      <faultBranch>6 7</faultBranch>
      <timeStep> 0.01</timeStep>
    </faultEvent>
  </faultEvents>
</Dynamic_simulation>

```

The fault used in the simulation is specified using the same **faultEvents** block as for dynamic simulation. If the Kalman filter simulation is not being initialized from another calculation, the **networkConfiguration** field can also be added. The KnowFault and TimeOffset parameters are unique to the Kalman filter application and control whether the fault is considered to be a know event and whether all the time series data should be used in the analysis.

The Kalman filter block consists of the fields

```
<Kalman_filter>
  <KalmanAngData>IEEE14_Kalman_input_ang.csv</KalmanAngData>
  <KalmanMagData>IEEE14_Kalman_input_mag.csv</KalmanMagData>
  <generatorParameters>IEEE14_classicGen.dyr</generatorParameters>
  <ensembleSize>21</ensembleSize>
  <gaussianWidth>1e-2</gaussianWidth>
  <noiseScale>1e-4</noiseScale>
  <randomSeed>931316785</randomSeed>
  <maxSteps>3000</maxSteps>
  <LinearSolver>
    <PETScOptions>
      -ksp_view
      -ksp_type richardson
      -pc_type lu
      -pc_factor_mat_solver_package superlu_dist
      -ksp_max_it 1
    </PETScOptions>
  </LinearSolver>
</Kalman_filter>
```

The **KalmanAngData** and **KalmanMagData** fields specify the locations of the files containing the time series data for the voltage magnitude and phase angle. The .dyr file containing the generator parameters (classical generators only) is specified in the **generatorParameters** field. Additional Kalman filter parameters include

- **ensembleSize**: The number of random ensembles generated for the Kalman filter calculation.
- **gaussianWidth**:
- **noiseScale**:
- **randomSeed**: This is an arbitrary integer used to seed the GridPACK random number generator.
- **maxSteps**: this parameter can be used to control the number of steps simulated. If the number of steps is smaller than the number of steps in the time series data files, then only the number of steps set by **maxSteps** will be simulated.

The Kalman filter also needs to make use of linear solvers and the type of solver and its parameters can be specified in this block as well.

The Kalman filter module is represented by the **KalmanApp** class and the **KalmanNetwork**, both of which are in the **gridpack::kalman\_filter** namespace. At present there are only a few functions in this class, more will probably be added as we develop this module further. Apart from the constructor and destructor, the **KalmanApp** class has a method for reading in a network from a PSS/E formatted file and partitioning it among processors

```
void readNetwork(boost::shared_ptr<KalmanNetwork> &network,  
                gridpack::utility::Configuration *config)
```

If the network already exists, then it can be applied to an existing KalmanApp object using the function

```
void readNetwork(boost::shared_ptr<KalmanNetwork> &network,  
                gridpack::utility::Configuration *config)
```

The application can be initialized by calling the function

```
void initialize()
```

This function will read in the files containing the time series data for the voltage magnitude and phase angles and will set up and configure the calculation based on the parameters in the input file. The simulation is run and output generated using

```
void solve()
```

The values of the rotor speed and rotor angle for all generators will be written to the files **omega.dat** and **delta.dat** after this simulation is run.

## GridPACK Examples

This section will expand on the discussion of the power flow application and provide additional examples of how GridPACK can be used to develop applications. Two of these are simple applications that have been provided in GridPACK that illustrate how the code works, without necessarily getting involved in the details that would be needed to implement a realistic power grid model. The third example is an in-depth discussion of the contingency analysis application. This is an actual application but it is fairly short and provides a good illustration of how to create multi-task simulations and also an example of how to use modules.

The first of the simple examples consists of a “hello world” program that writes a message from a small 10 x 10 square grid of buses and branches. The second example calculates the electric current flow through a square grid of resistors. Both examples are designed to show how the basic features of the GridPACK framework interact with each other. More complicated examples for realistic models can be found in the modules and components directories under applications. Although these examples represent more complicated bus and branch models, they contain many of the same characteristics that can be found in the hello world and resistor grid programs.

The contingency analysis application is also discussed in depth since this illustrates a great many of the advanced features of GridPACK in a fairly short code. These features include creating your own parser, using subcommunicators and the task manager, using modules and controlling output.

### “Hello World”

The “Hello world” program is a famous example problem from C programming. Many other packages have adopted the spirit of this program, if not the specifics, to describe the simplest non-trivial program that can be written using the package. In this section, a program that prints out a message from each of the buses and branches on a small grid is described. This application requires us to define branch and bus classes, create a network class and implement a top level application.

We start by implementing the load and serialWrite methods in the **BaseComponent** class for the bus and branch classes of our “Hello world” application. The bus and branch classes for this application are called **HWBus** and **HWBranch** and have the header file

```
#ifndef _hw_components_h_
#define _hw_components_h_

#include "boost/smart_ptr/shared_ptr.hpp"
#include "gridpack/include/gridpack.hpp"

namespace gridpack {
namespace hello_world {

class HWBus
: public gridpack::component::BaseBusComponent {
public:
    HWBus(); // Constructor
    ~HWBus() // Destructor
    void load(const boost::shared_ptr
               <gridpack::component::DataCollection> &data);
    bool serialWrite(char *string, const int bufsize,
                     const char *signal = NULL);
private:
    int p_original_idx;
    friend class boost::serialization::access;
    template<class Archive> void serialize(Archive &ar,
        const unsigned int version)
    {
```



```

        ar & boost::serialization::base_object
            <gridpack::component::BaseBusComponent>(*this)
            & p_original_idx;
    }
};

class HWBranch
: public gridpack::component::BaseBranchComponent {
public:
    HWBranch();    //Constructor
    ~HWBranch();   //Destructor
    void load(const boost::shared_ptr
                <gridpack::component::DataCollection> &data);
    bool serialWrite(char *string, const int bufsize,
                    const char *signal = NULL);

private:
    int p_original_idx1;
    int p_original_idx2;

friend class boost::serialization::access;

template<class Archive>
void serialize(Archive & ar, const unsigned int version)
{
    ar & boost::serialization::base_object
        <gridpack::component::BaseBranchComponent>(*this)
        & p_original_idx1
        & p_original_idx2;
}

};

typedef network::BaseNetwork<HWBus, HWBranch > HWNetwork;
}    // hello_world
}    // gridpack
#endif

```

The **HWBus** class has one private member, **p\_original\_idx**, which is the index of the bus in the network configuration file. Similarly, the **HWBranch** class has two private members, **p\_original\_idx1** and **p\_original\_idx2**, representing the buses at the “from” and “to” ends of the branch. The name of the file containing this code is **hw\_components.hpp**. The

first two lines of the file are the standard preprocessor protection flags that guarantee that any declarations in this file only appear in another file a single time. The next two lines include the Boost smart pointer header file and the header files from the GridPACK framework. The next two lines declare that all functions and classes in the file are in the **gridpack::hello\_world** namespace. The use of namespaces is up to the user and other choices are possible. The declaration of the **HWBus** class inherits from the **BaseBusComponent** class so all functions in the **BaseBusComponent** class are available to **HWBus**. **BaseBusComponent** also provides some virtual functions, along with their default implementations, that can be overwritten in **HWBus**. Two of these are the **load** and **serialWrite** functions. Only these functions are used in the “Hello world” application, the remaining functions in the bases classes are represented by the default implementations. Inside **HWBus** are declarations for the constructor, destructor, **load** and **serialWrite** functions. These will be implemented in the **hw\_components.cpp** file.

The final component in **HWBus** is the implementation of the serialize method. This method is used when copying the class from one processor to another and allows the program to move all the data associated with a particular instance of **HWBus** to another processor. The friend declaration means that **HWBus** has access to protected methods and data in **boost::serialization::access** and the templated serialization function is used to declare all internal data members that need to be transferred with the **HWBus** instance if it is moved from on processor to another. These elements include whatever base class **HWBus** may be derived from, which is represented by the element

```
boost::serialization::base_object<gridpack::component  
::BaseBusComponent>(*this)
```

The remaining data element is **p\_original\_idx**. The variable **ar** of type **Archive** is appended to using the operator **&**. In this case the data appended to **ar** is any serialized data coming from the base class and the variable **p\_original\_idx**. The serialization function is recursive, so including the base class is enough to guarantee that any variables beneath that are also included in the serialization.

The declaration for **HWBranch** is very similar. The only major difference is that there are two private variables representing the buses at either end of the branch and these must both be included in the **serialize** function.

The bottom of the file contains a typedef declaration for a network using **HWBus** and **HWBranch** for it bus and branch classes. This is a convenience and makes it easier to define other functions and classes in the application.

The `hw_components.cpp` file contains the actual implementation of the functions declared in `hw_components.hpp`. The declarations for STL vectors and iostreams and the `hw_components.hpp` file are included at the top of the file so that all functions in the class are defined. For `HWBus`, the constructor and destructor are trivial and are given by

```
gridpack::hello_world::HWBus::HWBus()
{
    p_original_idx = 0;
}
gridpack::hello_world::HWBus::~~HWBus()
{
}
```

The `load` function is more interesting and is designed to transfer data that was read in from the network configuration file to the internal parameters of the bus. In this case, there is only one internal parameter, so `load` is fairly simple. The bus ID is stored in the variable `BUS_NUMBER`, so the load implementation is

```
void gridpack::hello_world::HWBus::load(const
    boost::shared_ptr<gridpack::component::DataCollection> &data)
{
    data->getValue(BUS_NUMBER,&p_original_idx);
}
```

All the parameters associated with the bus that came from the network configuration file are stored in the `data DataCollection` object, so the `getValue` statement is used to get the value from `data` and assign it to `p_original_index`. A completely listing of all variables that might be found in a `DataCollection` object can be found in the `dictionary.hpp` file located in the `src/parser` directory.

The `serialWrite` function returns a string with a message from the bus if called by some other program (in this case an instance of `SerialBusIO`). For “Hello world”, the bus reports back the bus index using the function

```
bool gridpack::hello_world::HWBus::serialWrite(char *string,
    const int bufsize, const char *signal)
{
    sprintf(string,"Hello world from bus %d\n",p_original_idx);
    return true;
}
```

For this case, both the incoming variables `bufsize` and `signal` are ignored since “Hello world” only has one type of output and it is guaranteed to fit in the buffer, but both variables

could be used in more complicated implementations. The **bufsize** variable can be used to make sure that the string does not exceed an internal buffer size and **signal** can be used to produce different outputs depending on what the actual contents of signal are. For the **serialWrite** implementations described for this application, guaranteeing that the strings fit inside the buffer is straightforward, since all strings are the same size. For real applications, this may not be the case. For example, when printing out generator properties, the strings from buses can vary in size because the number of generators on a bus can vary.

The implementations of the functions in **HWBranch** are similar. The constructor and destructor are

```
gridpack::hello_world::HWBranch::HWBranch(void)
{
    p_original_idx1 = 0;
    p_original_idx2 = 0;
}
gridpack::hello_world::HWBranch::~~HWBranch(void)
{
}
```

The **load** function is given by

```
void gridpack::hello_world::HWBranch::load(
    const boost::shared_ptr<gridpack::component::DataCollection>
    &data)
{
    data->getValue(BRANCH_FROMBUS,&p_original_idx1);
    data->getValue(BRANCH_TOBUS,&p_original_idx2);
}
```

This is similar to the implementation of the load function for **HWBus**, except that the internal data members are mapped to the values of the **BRANCH\_FROMBUS** and **BRANCH\_TOBUS** elements of the data collection object. The **serialWrite** function is

```
bool gridpack::hello_world::HWBranch::serialWrite(char *string,
    const int bufsize, const char *signal)
{
    sprintf(string,
        "Hello world from the branch connecting bus %d to bus %d\n",
        p_original_idx1, p_original_idx2);
    return true;
}
```

Every branch prints out a string describing the branch in terms of the bus IDs at each end of the branch. Again, the incoming bufsize and signal variables are ignored in this case and it is assumed that the buffer size assigned to the **SerialBranchIO** object when it is instantiated is sufficiently large to guarantee that all strings from every branch will fit.

The implementation of the factory class for the “Hello world” application is straightforward, since the class only needs the functionality in the **BaseFactory** class. The complete class is given by

```
#ifndef _hw_factory_h_
#define _hw_factory_h_

#include "boost/smart_ptr/shared_ptr.hpp"
#include "gridpack/include/gridpack.hpp"
#include "hw_components.hpp"

namespace gridpack {
namespace hello_world {

class HWFactory
: public gridpack::factory::BaseFactory<HWNetwork> {
public:
    HWFactory(boost::shared_ptr<HWNetwork> network)
        : gridpack::factory::BaseFactory<HWNetwork>(network)
    {
    }
    ~HWFactory() {}
};
} // hello_world
} // gridpack
#endif
```

This class is defined in the **hw\_factory.hpp** file. Because the class is so simple, the complete class declaration is given in **hw\_factory.hpp** and there is no corresponding **.cpp** file. In addition to including the **gridpack.hpp** header, this file also includes **hw\_components.hpp**, so it has the definitions of **HWNetwork**. The **HWFactory** constructor is used to initialize the underlying **BaseFactory** object with the network that is passed in through the argument list. That is the only functionality that is defined in this class.

The application class that is built on top of the component and factory classes consists of the class

```

#ifndef _hw_app_h_
#define _hw_app_h_

namespace gridpack {
namespace hello_world {

class HwApp
{
public:
    HwApp(void);
    ~HwApp(void);
    void execute(int argc, char** argv);
};

} // hello_world
} // gridpack
#endif

```

This class is declared in **hw\_app.hpp**. Apart from the constructor and destructor, there is only the function **execute**, which is used to actually run the program. This takes the standard **argc** and **argv** variables as arguments, which could be passed in from the top level calling program.

The implementation of these functions are relatively simple, most of the complexity for this program is in defining the bus and branch classes. The implementations are defined in the file **hw\_app.cpp**

```

#include <iostream>
#include "boost/smart_ptr/shared_ptr.hpp"
#include "gridpack/include/gridpack.hpp"
#include "hw_app.hpp"
#include "hw_factory.hpp"

gridpack::hello_world::HwApp::HwApp(void)
{
}

gridpack::hello_world::HwApp::~~HwApp(void)
{
}

void gridpack::hello_world::HwApp::execute(int argc, char** argv)

```

```

{
    gridpack::parallel::Communicator world;
    boost::shared_ptr<HWNetwork> network(new HWNetwork(world));

    std::string filename = "10x10.raw";
    gridpack::parser::PTI23_parser<HWNetwork> parser(network);
    parser.parse(filename.c_str());
    gridpack::hello_world::HWFactory factory(network);
    factory.load();

    gridpack::serial_io::SerialBusIO<HWNetwork> busIO(128,network);
    busIO.header("\nMessage from buses\n");
    busIO.write();
    gridpack::serial_io::SerialBranchIO<HWNetwork>
        branchIO(128,network);
    branchIO.header("\nMessage from branches\n");
    branchIO.write();
}

```

The top of the file contains the **gridpack.hpp** header as well as the application headers. The constructor and destructors for the **HWApp** class are the standard defaults, so only the **execute** function has any significant behavior. This function starts by defining a communicator on the set of all processors and using that to instantiate and instance of an **HWNetwork**. At this point the network exists, but it contains no buses or branches. The next step is to read in a network configuration file with the name **10x10.raw**. This file is written using the standard PSS/E version 23 format. For this simple application, it is assumed that the file is available in the directory in which the program is being run (this file is included in the **hello\_world** directory as part of the GridPACK distribution). The program creates an instance of a **PTI23\_parser** and uses this to parse the configuration file. The program now has a copy of the full network stored internally, but the buses and nodes are not distributed in a way that is convenient for computation. Calling the partition method on the network redistributes all buses and branches so that each process has a relatively connected chunk of the network.

The next step is to create an **HWFactory** instance and use this to call the base class **load** method. This method in turn calls the **load** method on all the individual buses and branches and transfers data from the data collection objects to the internal parameters of the buses and branches. The data collection objects were initialized with data collected from the **10x10.raw** file when the **parse** function was called. The remaining lines create **SerialBusIO** and **SerialBranchIO** objects that are used to print out the messages from individual bus and branch objects. The **busIO** object is used to print out a header (“Message from buses”) and then

a message from each bus identifying itself by the bus ID defined in the PSS/E file. Similarly, the **branchIO** object writes out a header and then a message from each branch identifying itself by the IDs of the buses at either end.

The final part of the “Hello world” application is the main calling program, located in the file `hw_main.cpp`. This program consists of the lines

```
#include "gridpack/include/gridpack.hpp"
#include "hw_app.hpp"

int main(int argc, char **argv)
{
    gridpack::parallel::Environment env(argc, argv);

    gridpack::hello_world::HWApp app;
    app.execute(argc, argv);
    return 0;
}
```

The program consists of a line creating a parallel environment, a line instantiating an **HWApp**, and a line calling the `execute` method on the application. The constructor for the parallel environment initializes the underlying parallel communication libraries. The destructor is called at the end of `main` and terminates all communication libraries so that the program exits cleanly. The **HWApp** instance runs the application when **`execute`** is called. A portion of the output looks like

#### Message from buses

```
Hello world from bus 1
Hello world from bus 2
Hello world from bus 3
Hello world from bus 4
Hello world from bus 5
Hello world from bus 6
Hello world from bus 7
:
```

#### Message from branches

```
Hello world from the branch connecting bus 1 to bus 2
Hello world from the branch connecting bus 2 to bus 3
Hello world from the branch connecting bus 3 to bus 4
Hello world from the branch connecting bus 4 to bus 5
Hello world from the branch connecting bus 5 to bus 6
```

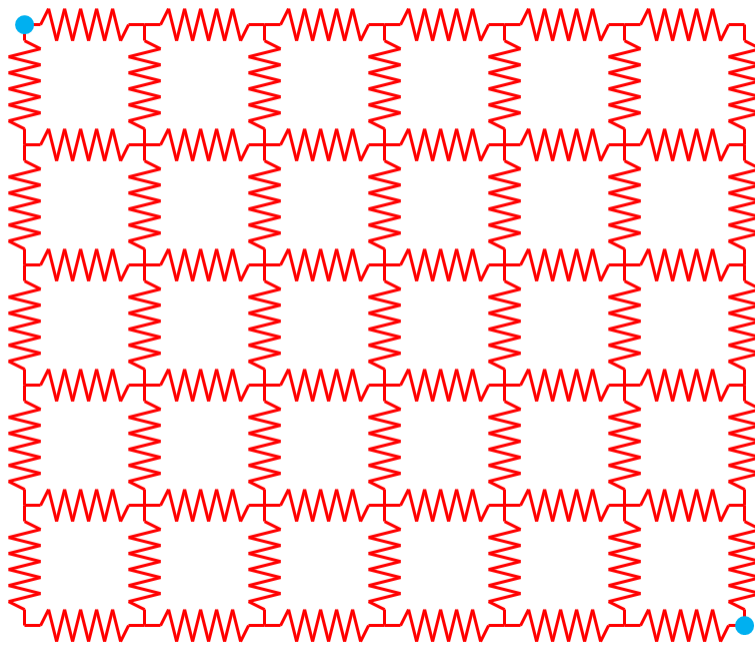


:

Note that this output would be the same, regardless of the number of processors that are used to run the code. This is in spite of the fact that the distribution of buses and branches may be different for different numbers of processors.

### Resistor Grid Application

The resistor grid is a more complicated example that illustrates how GridPACK can be used to set up equations describing a physical system and then solve the system using a linear solver. The physical system is a rectangular grid with resistors connecting all the nodes. Two nodes are chosen to be set at fixed potentials, these then drive currents through the rest of the network resulting in different currents on the individual branches and different voltages on the different buses (nodes). The system is illustrated schematically in Figure 12.



**Figure 12.** A schematic diagram of a simple resistor grid network. The buses (nodes) in blue are set at fixed external voltages, the remaining bus voltages and branch currents are calculated by the application.

The topology and choice of nodes held at fixed potential is determined by the network configuration file, as are the values of the resistance on each of the branches. The system is described by a set of coupled equations representing the application of Kirkoff's law to each of the nodes that is not held at a fixed potential. Kirkoff's law is expressed by the equations

$$\sum_{\beta \in \{\alpha\}} i_{\alpha\beta} = 0$$

where  $i_{\alpha\beta}$  is the current flowing between nodes  $\alpha$  and  $\beta$  and  $\{\alpha\}$  is the set of nodes connected directly to  $\alpha$ . This current can be found from Ohm's law

$$i_{\alpha\beta} = \frac{V_{\alpha} - V_{\beta}}{R_{\alpha\beta}}$$

Where  $V_{\alpha}$  and  $V_{\beta}$  are the voltage potentials on nodes  $\alpha$  and  $\beta$  and  $R_{\alpha\beta}$  is the resistance on the branch connecting nodes  $\alpha$  and  $\beta$ . Plugging the expression for the current back into Kirkoff's law gives the equation

$$\sum_{\beta \in \{\alpha\}} \frac{V_{\alpha} - V_{\beta}}{R_{\alpha\beta}} = V_{\alpha} \sum_{\beta \in \{\alpha\}} \frac{1}{R_{\alpha\beta}} - \sum_{\beta \in \{\alpha\}} \frac{V_{\beta}}{R_{\alpha\beta}} = 0$$

The unknowns in this system are the potentials  $V_{\alpha}$ . Kirkoff's law applies to any node that does not have an applied value of the potential. The nodes that do have a fixed potential appear as part of the right hand side vector. Assuming that any node with a non-fixed value of the potential is attached to at most one fixed node, then the  $\alpha$ th element of the right hand side vector is

$$\frac{V_{\beta}^0}{R_{\alpha\beta}}$$

where  $V_{\beta}^0$  is the value of the fixed potential on node  $\beta$  and  $\alpha$  is attached to  $\beta$ . If  $\alpha$  is not attached to  $\beta$ , then the element is zero. The voltages can be evaluated by solving the matrix equation

$$\bar{\bar{C}} \cdot \bar{V} = \bar{I}_0$$

The voltage vector and right hand side have already been discussed. The matrix elements have the form

$$C_{\alpha\alpha} = \sum_{\beta \in \{\alpha\}} \frac{1}{R_{\alpha\beta}}$$

$$C_{\alpha\beta} = -\frac{1}{R_{\alpha\beta}} \text{ if } \alpha \neq \beta$$

With this background, we can talk about the implementation of the resistor grid application.

Much of the basic structure of the classes has already been discussed in the previous example of "Hello world", so we will limit ourselves to discussing new features. The **RGBus** class inherits

from the **BaseBusComponent** class and implements the following functions (in addition to the constructor and destructor)

```
void load(const boost::shared_ptr
    <gridpack::component::DataCollection> &data);
bool isLead() const;
double voltage() const;
bool matrixDiagSize(int *isize, int *jsize) const;
bool matrixDiagValues(ComplexType *values);
bool vectorSize(int *isize) const;
bool vectorValues(ComplexType *values);
void setValues(gridpack::ComplexType *values);
int getXCBufSize();
void setXCBuf();
bool serialWrite(char *string, const int bufsize,
    const char *signal = NULL);
```

In addition, the **RGBus** class has three private members

```
bool p_lead;
double *p_voltage;
double p_v;
```

The variable **p\_lead** keeps track of whether a bus has a fixed voltage applied to it. In order to correctly calculate the currents, it is necessary to exchange voltages at the end of the calculation. The voltages at each bus are stored in an exchange buffer that can be accessed by the pointer **p\_voltage**. The voltages in the external PSS/E file are read in before the exchange buffer is allocated, so to make sure there is a variable to store the value, the variable **p\_v** is also included as a private member. In addition to implementing **load** and **serialWrite**, the **RGBus** class implements several functions in the **MatVecInterface**, as well as two functions that are unique to this class.

Similarly, the **RGBranch** class implements the functions

```
void load(const boost::shared_ptr
    <gridpack::component::DataCollection> &data);
double resistance(void) const;
bool matrixForwardSize(int *isize, int *jsize) const;
bool matrixReverseSize(int *isize, int *jsize) const;
bool matrixForwardValues(ComplexType *values);
bool matrixReverseValues(ComplexType *values);
bool serialWrite(char *string, const int bufsize,
```

```
const char *signal = NULL);
```

and has the private member

```
double p_resistance;
```

The **RGBus** load method has the implementation

```
void gridpack::resistor_grid::RGBus::load(const
    boost::shared_ptr<gridpack::component::DataCollection> &data)
{
    int type;
    data->getValue(BUS_TYPE,&type);
    if (type == 2) {
        p_lead = true;
        data->getValue(BUS_BASEKV,&p_v);
    }
}
```

The PSS/E file that is used to run this application has been configured so that the bus type parameter is set to 2 if the bus has a fixed voltage and the value of the voltage is stored in the **BUS\_BASEKV** variable. The private member **p\_lead** is initialized to false in the **RGBus** constructor and **p\_v** is initialized to zero. In the **load** method, the bus type is assigned from the **BUS\_TYPE** variable in the data collection. If it is 2, the bus has a fixed value of the potential and **p\_lead** is set to true. The value of **p\_v** is assigned to whatever is stored in the **BUS\_BASEKV** variable when the bus type is 2. The contents of **p\_v** will eventually be mapped to **p\_voltage**, once the exchange buffers are allocated.

The **load** function for **RGBranch** simply assigns the data collection variable **BRANCH\_R** to the private member **p\_resistance**.

```
void gridpack::resistor_grid::RGBranch::load(
    const boost::shared_ptr
    <gridpack::component::DataCollection> &data)
{
    data->getValue(BRANCH_R,&p_resistance,0);
}
```

Once the bus and branch private members have been set using the load methods, the values can be recovered by other objects using the accessors **isLead**, **voltage**, and **resistance**. These functions are used in the math interface implementations to calculate values of the matrix elements and right hand side vectors and have the relatively simple forms

```

bool gridpack::resistor_grid::RBus::isLead() const
{
    return p_lead;
}
double gridpack::resistor_grid::RBus::voltage() const
{
    return *p_voltage;
}
double gridpack::resistor_grid::RBranch::resistance(void) const
{
    return p_resistance;
}

```

Note that the **voltage** function is returning the contents of **p\_voltage**, which will contain up-to-date values of the voltage once the calculation begins.

The diagonal matrix block routines in the bus class have the implementations

```

bool gridpack::resistor_grid::RBus::matrixDiagSize(int *isize,
    int *jsize) const
{
    if (!p_lead) {
        *isize = 1;
        *jsize = 1;
        return true;
    } else {
        return false;
    }
}
bool gridpack::resistor_grid::RBus::matrixDiagValues(
    ComplexType *values)
{
    if (!p_lead) {
        gridpack::ComplexType ret(0.0,0.0);
        std::vector<boost::shared_ptr<BaseComponent> > branches;
        getNeighborBranches(branches);
        int size = branches.size();
        int i;
        for (i=0; i<size; i++) {
            gridpack::resistor_grid::RBranch *branch
                = dynamic_cast<gridpack::resistor_grid::RBranch*>

```

```

        (branches[i].get());
    ret += 1.0/branch->resistance();
}
values[0] = ret;
return true;
} else {
    return false;
}
}

```

The **matrixDiagSize** routine returns a single element in the **values** array if the bus is not a lead with a fixed voltage, otherwise it returns false and there are no values in the **values** array. The **matrixDiagValues** function sets the first element of the **values** array equal to the sum of the reciprocal of the resistances on all the attached branches, if the bus is not a lead. To calculate this quantity, it starts by calling the **getNeighborBranches** function to get a list of pointers to attached branches. These pointers are all of type **BaseComponent**, so they need to be cast to pointers of type **RGBranch** before functions like **resistance** can be called on them. This is done by first calling the **get** function on the **shared\_ptr** to the **BaseComponent** object to get a bare pointer to the neighboring branch and then doing a dynamic cast to a pointer of type **RGBranch**. The resistance method can now be called on the **RGBranch** pointer to get the resistance of the branch and use it to calculate the contribution to the diagonal matrix element. This value is assigned to **values[0]**. If the bus is a lead, then no values are calculated and the function returns false. It is also worth noting that this function will only be called on buses that are local to the process, so each bus that evaluates a diagonal matrix element will have a complete set of branches attached to it. This is not the case for ghost buses. These have only one branch attached to them, no matter how many branches are attached to it in the original network.

The off-diagonal elements are calculated by the branch components in the functions **matrixForwardSize**, **matrixReverseSize**, **matrixForwardValues**, and **matrixReverseValues**. The matrix  $\bar{\bar{C}}$  for the resistor grid problem is completely symmetric, so in this case, the forward and reverse calculations are identical. For realistic power problems, this is not generally true, and the forward and reverse functions will have different implementations. The forward functions are described below, the implementation of the reverse functions is identical. The branch forward size and value functions are

```

bool gridpack::resistor_grid::RGBranch::matrixForwardSize(
    int *isize, int *jsize) const
{
    gridpack::resistor_grid::RBus *bus1

```

```

        = dynamic_cast<gridpack::resistor_grid::RBus*>(getBus1().get());
gridpack::resistor_grid::RBus *bus2
        = dynamic_cast<gridpack::resistor_grid::RBus*>(getBus2().get());
if (!bus1->isLead() && !bus2->isLead()) {
    *isize = 1;
    *jsize = 1;
    return true;
} else {
    return false;
}
}
bool gridpack::resistor_grid::RGBranch::matrixForwardValues(
    ComplexType *values)
{
    gridpack::resistor_grid::RBus *bus1
        = dynamic_cast<gridpack::resistor_grid::RBus*>(getBus1().get());
    gridpack::resistor_grid::RBus *bus2
        = dynamic_cast<gridpack::resistor_grid::RBus*>(getBus2().get());
    if (!bus1->isLead() && !bus2->isLead()) {
        values[0] = -1.0/p_resistance;
        return true;
    } else {
        return false;
    }
}
}

```

Before these functions can calculate return values, they must first determine if one of the buses at either end of the branch is a lead bus. To do this, the functions need to get pointers to the “from” and “to” buses at either end of the branch. They can do this through the **getBus1** and **getBus2** calls in the **BaseBranchComponent** class which return pointers of type **BaseComponent**. These pointers can then be converted to **RBus** pointers by a dynamic cast. The **isLead** functions can be called to find out if either bus is a lead bus. If neither bus is a lead bus, the size of the off-diagonal block is returned as a 1x1 matrix and the off-diagonal matrix element is calculated and returned in **values[0]**. Otherwise both functions return false to indicate that there is no contribution to the matrix from this branch.

In addition to calculating values of the matrix  $\bar{\bar{C}}$ , it is also necessary to set up the right hand side vector. This is done via the functions **vectorSize** and **vectorValues** defined on the buses. Only buses that are not lead buses contribute to the right hand side vector. On the other hand, the

only non-zero values in the right hand side vector come from lead buses that are attached to non-lead buses. The **vectorSize** function has the implementation

```
bool gridpack::resistor_grid::RBus::vectorSize(int *isize) const
{
    if (!p_lead) {
        *isize = 1;
        return true;
    } else {
        return false;
    }
}
```

If a bus is not a lead bus, it contributes a single value, otherwise it does not and the function returns false. The **vectorValues** function is a bit more complicated. It has the form

```
bool gridpack::resistor_grid::RBus::vectorValues(ComplexType *values)
{
    if (!p_lead) {
        std::vector<boost::shared_ptr<BaseComponent> > branches;
        getNeighborBranches(branches);
        int size = branches.size();
        int i;
        gridpack::ComplexType ret(0.0,0.0);
        for (i=0; i<size; i++) {
            gridpack::resistor_grid::RBranch *branch
                = dynamic_cast<gridpack::resistor_grid::RBranch*>
                    (branches[i].get());
            gridpack::resistor_grid::RBus *bus1
                = dynamic_cast<gridpack::resistor_grid::RBus*>
                    (branch->getBus1().get());
            gridpack::resistor_grid::RBus *bus2
                = dynamic_cast<gridpack::resistor_grid::RBus*>
                    (branch->getBus2().get());
            if (bus1 != this && bus1->isLead()) {
                ret += bus1->voltage()/branch->resistance();
            } else if (bus2 != this && bus2->isLead()) {
                ret += bus2->voltage()/branch->resistance();
            }
        }
        values[0] = ret;
    }
}
```



```

        return true;
    } else {
        return false;
    }
}

```

The **vectorValues** function starts by getting a list of branches that are attached to the calling bus and then looping over the list. Pointers to each of the branches, as well as the buses at each end of the branch are obtained using the **getBus1** and **getBus2** functions. It is still necessary to determine which end of the branch is opposite the calling bus and this can be done by checking the conditions **bus1 != this** and **bus2 != this**. One of these will be true for the bus opposite the calling bus. If this bus is also a lead bus, then a contribution is added to the right hand side vector element. The contribution can be calculated by getting the value of the fixed voltage from the lead bus and dividing it by the resistance of the branch. These values can be obtained by calling the bus **voltage** function and the branch **resistance** function. The **\*p\_voltage** value of the calling bus is not used. If the calling bus is a lead bus, then the function returns false.

The last function related to vectors that is implemented in the **MatVecInterface** is the **setValues** function

```

void gridpack::resistor_grid::RBus::setValues(
    gridpack::ComplexType *values)
{
    if (!p_lead) {
        p_voltage = real(values[0]);
    }
}

```

Once the voltages have been calculated by solving Kirkoff's equations, it is necessary to have some way of pushing these back on the buses so they can be written to output. The results of the linear solver are returned in the **values** array. The number of values in this array corresponds to the number of values contributed to the right hand side vector (in this case 1 if the bus is not a lead). Thus, the value is assigned to the internal **p\_voltage** variable if the bus is not a lead bus. This function will be called by all buses as part of the **mapToBus** function in the **BusVectorMap**.

In order to correctly calculate the current on all branches for export to standard out, it is necessary to have up-to-date values of the voltage on all buses, including ghost buses. This requires a data exchange at the end of the calculation. To enable this exchange, the

**getXCBufSize** and **setXCBuf** functions must be implemented in the **RGBus** class. These functions have the form

```
int gridpack::resistor_grid::RGBus::getXCBufSize()
{
    return sizeof(double);
}
void gridpack::resistor_grid::RGBus::setXCBuf(void *buf)
{
    p_voltage = static_cast<double*>(buf);
    *p_voltage = p_v;
}
```

The only variable that needs to be exchange is the value of the potential, so **getXCBufSize** returns the number of bytes in a single double precision variable. The **setXCBuf** function assigns the buffer pointed to by the variable **buf** to the internal data member **p\_voltage**. At the same time, it initializes the contents of **p\_voltage** to the variable **p\_v**, which contains the voltage read in from the external PSS/E file.

The **serialWrite** functions on the buses and branches are used to write the voltages and currents on all buses and branches to standard output. The **serialWrite** function on the buses has the form

```
bool gridpack::resistor_grid::RGBus::serialWrite(char *string,
    const int bufsize, const char *signal)
{
    if (p_lead) {
        sprintf(string,"Voltage on bus %d: %12.6f (lead)\n",
            getOriginalIndex(),*p_voltage);
    } else {
        sprintf(string,"Voltage on bus %d: %12.6f\n",
            getOriginalIndex(),*p_voltage);
    }
    return true;
}
```

All buses return a string so the function always returns true. The printout consists of the bus index, obtained with the **getOriginalIndex** function, and the value of the voltage on the bus. Lead buses are marked in the output, indicating that the voltage is the same as that specified in the input file, the remaining voltages are calculated by solving Kirkoff's equations. For

branches, the `serialWrite` function is used to calculate and print the current flowing across each branch

```
bool gridpack::resistor_grid::RGBBranch::serialWrite(char *string,
const int
    bufsize,  const char *signal)
{

    gridpack::resistor_grid::RBus *bus1
        = dynamic_cast<gridpack::resistor_grid::RBus*>(getBus1().get());
    gridpack::resistor_grid::RBus *bus2
        = dynamic_cast<gridpack::resistor_grid::RBus*>(getBus2().get());
    double v1 = bus1->voltage();
    double v2 = bus2->voltage();
    double icur = (v1 - v2)/p_resistance;
    sprintf(string,"Current on line from bus %d to %d is: %12.6f\n",
        bus1->getOriginalIndex(),bus2->getOriginalIndex(),icur);
    return true;
}
```

All branches report the current flowing through them, so this function also returns true for all branches. To calculate the current, it is necessary to get the value of the voltages at both ends of the branch using methods already described and then calculate the current by dividing the difference in voltages by the resistance of the branch. The print line prints the current and uniquely identifies each branch by including the IDs of the buses at either end.

The factory class for resistor grid application only uses functionality in the `BaseFactory` class and has the simple form

```
class RGFactory
: public gridpack::factory::BaseFactory<RGNetwork> {
public:
    RGFactory(boost::shared_ptr<RGNetwork> network)
        : gridpack::factory::BaseFactory<RGNetwork>(network)
    {
    }
    ~RGFactory() {}
};
```

Again, the `BaseFactory` class from which `RGFactory` inherits is initialized by passing the network argument through the constructor. The declaration for this class is in the file `rg_factory.hpp`. There is no corresponding `.cpp` file.

The **RGApp** class declaration is also simple and consists of the functions

```
class RGApp
{
public:
    RGApp(void);
    ~RGApp(void);
    void execute(int argc, char** argv);
};
```

Again, arguments from the top level main program can be passed through to the execute function, which is responsible for implementing the actual resistor grid calculation. The **RGApp** class declaration is contained in the **rg\_app.hpp** file. The implementation is contained in the **rg\_app.cpp** file. The only complicated function in the implementation is execute, which consists of

```
void gridpack::resistor_grid::RGApp::execute(int argc, char** argv)
{
    // read configuration file
    gridpack::parallel::Communicator world;
    gridpack::utility::Configuration *config =
        gridpack::utility::Configuration::configuration();
    config->open("input.xml",world);
    gridpack::utility::Configuration::CursorPtr cursor;
    cursor = config->getCursor("Configuration.ResistorGrid");

    // create network and read in external PTI file
    // with network configuration
    boost::shared_ptr<RGNetwork> network(new RGNetwork(world));
    gridpack::parser::PTI23_parser<RGNetwork> parser(network);
    std::string filename;
    if (!cursor->get("networkConfiguration",&filename)) {
        filename = "small.raw";
    }
    parser.parse(filename.c_str());

    // partition network
    network->partition();

    // create factory and load parameters from input
    // file to network components
```

```

gridpack::resistor_grid::RGFactory factory(network);
factory.load();

// set network components using factory and set up exchange
// of voltages between buses
factory.setComponents();
factory.setExchange();
network->initBusUpdate();

// create mapper to generate voltage matrix
gridpack::mapper::FullMatrixMap<RGNetwork> vMap(network);
boost::shared_ptr<gridpack::math::Matrix> V = vMap.mapToMatrix();

// create mapper to generate RHS vector
gridpack::mapper::BusVectorMap<RGNetwork> rMap(network);
boost::shared_ptr<gridpack::math::Vector> R = rMap.mapToVector();

// create solution vector by cloning R
boost::shared_ptr<gridpack::math::Vector> X(R->clone());

// create linear solver and solve equations
gridpack::math::LinearSolver solver(*V);
solver.configure(cursor);
solver.solve(*R, *X);

// push solution back on to buses
rMap.mapToBus(X);

// exchange voltages so that all buses have correct values. This
// guarantees that current calculations on each branch are correct
network->updateBuses();

// create serial IO objects to export data
gridpack::serial_io::SerialBusIO<RGNetwork> busIO(128, network);
char ioBuf[128];
busIO.header("\nVoltages on buses\n\n");
busIO.write();

gridpack::serial_io::SerialBranchIO<RGNetwork>
    branchIO(128, network);

```

```

    branchIO.header("\nCurrent on branches\n\n");
    branchIO.write();
}

```

The beginning of the resistor grid application is more complicated than “Hello world” in that it uses an input file to control the properties of the linear solver that is used to solve current equations. To read in the input file, the application starts by creating a communicator on the set of all processors. Only one configuration object is available to the application and the **execute** function gets a pointer to this instance by calling the static function

**Configuration::configuration()**. This pointer can then be used to read in the input file, “**input.xml**”, across all processes in the communicator **world** using the **open** method. All processors now have access to the contents of **input.xml**. The input file contains two pieces of information, the name of the PSS/E formatted resistor grid configuration file and the parameters for the linear solver. The input file has the form

```

<?xml version="1.0" encoding="utf-8"?>
<Configuration>
  <ResistorGrid>
    <networkConfiguration> small.raw </networkConfiguration>
    <LinearSolver>
      <PETScOptions>
        -ksp_view
        -ksp_type richardson
        -pc_type lu
        -pc_factor_mat_solver_package superlu_dist
        -ksp_max_it 1
      </PETScOptions>
    </LinearSolver>
  </ResistorGrid>
</Configuration>

```

The resistor grid file name can be obtained by getting a cursor pointer that is pointed at the **ResistorGrid** block in the input file by using the **getCursor** function and then using the **get** function to retrieve the actual file name located in the **networkConfiguration** field. If no file is specified in the input deck, the file name defaults to “**small.raw**”. At the same time, an **RGNetwork** object is instantiated and used to initialize on instance of **PTI23\_parser**. This can then read in the resistor grid configuration file using the parse function.

At this point, all buses and branches have been created, but they may not be distributed in a way that supports computation. The network **partition** function is called to redistribute the

network so that each process has maximal connections between components located on the process and minimal connections to components located on other processes. The ghost buses and branches are also added by the **partition** function.

After partitioning, an **RGFactory** object is created and the base class **load** method is called to initialize the internal data elements on each bus and branch in the network. This function initializes both locally held components as well as ghost components, so there is no need for a data exchange to guarantee that all components are up to date. The factory also calls the base class **setComponents** method, which determines several types of internal indices that are used to set up calculations. The buffers needed to exchange data at the end of the calculation are set up by a call to the factory **setExchange** method. Additional internal data structures needed for the data exchange between buses are created by calling the network **initBusUpdate** method. No data exchanges are needed between branch components.

The next step in the algorithm is to create the matrix  $\bar{\bar{C}}$ , the right hand side vector and a vector to contain the solution. Two separate mappers are needed, one for the matrix  $\bar{\bar{C}}$  and the other for the right hand side vector. For the matrix, the code creates an instance of a **FullMatrixMap** that is initialized with the resistor grid network. The **mapToMatrix** function is called to create the matrix **V**. The right hand side vector is created by creating instance of a **BusVectorMap** and using the **mapToVector** function to create the vector **R**. The solution vector **X** does not need to be initialized to any particular value, it just needs to be the same size as **R** so it is created by having **R** call the **clone** method in the **Vector** class and using the result to initialize **X** in the **Vector** class constructor.

Once **V**, **R**, and **X** are available, the equations can be solved using a linear solver. The linear solver is created by initializing an instance of **LinearSolver** with the matrix **V**. The solver class **configure** method can be used to transfer solver parameters in the **LinearSolver** block in **input.xml** to the solver. The **cursor** pointer that is taken as an argument to **configure** is already pointing to the **ResistorGrid** block in the input file, so **configure** will pick up any parameters in a **LinearSolver** block within the **ResistorGrid** block. After configuring the solver, the solution vector can be obtained by calling the **solve** method and the resulting voltages are pushed back to buses using the **mapToBus** method in the **BusVectorMap** class.

After calling **mapToBus**, all locally held buses have correct values of the voltage, but ghost buses still have their initial values. To correct the voltages on ghost buses, it is necessary to call the network **updateBuses** function. The buffers **p\_voltage** now contain correct values of the voltage on all buses.

The only remaining step is to write the results to standard output. The voltages are written by creating an instance of **SerialBusIO**. The maximum buffer size is set to 128 characters,

which is enough to hold any lines of output coming from the buses. A header labeling the bus output is written to standard out using the **header** method and then bus voltages are written by calling **write**. Similarly, output from the branches can be written by creating an instance of **SerialBranchIO**, writing a header using the **header** method and then calling **write**. Since only one type of output comes from the branches and buses, no character string is passed in as arguments to the **write** functions. The **execute** function has now completed all tasks associated with solving the resistor grid problem and passes control back to the main calling program.

The main calling program is relatively simple and consists of the code

```
int main(int argc, char **argv)
{
    gridpack::parallel::Environment env(argc, argv);
    gridpack::math::Initialize();
    gridpack::resistor_grid::RGApp app;
    app.execute(argc, argv);
    gridpack::math::Finalize();
    return 0;
}
```

The parallel computing environment is set up by creating an instance of **Environment**. The computing environment is also cleaned up at the end of the calculation when the destructor for this object is called. The math libraries are initialized by a call to the static **Initialize** method and cleaned up at the end of the calculation by a call to **Finalize**. The only remaining calls are to create an instance of an **RGApp** and call its **execute** method.

A portion of the output from the resistor grid calculation is the following

**GridPACK math module configured on 8 processors**

**:**  
**Voltages on buses**

<b>Voltage on bus 1:</b>	<b>1.000000 (lead)</b>
<b>Voltage on bus 2:</b>	<b>0.667958</b>
<b>Voltage on bus 3:</b>	<b>0.467469</b>
<b>Voltage on bus 4:</b>	<b>0.329598</b>
<b>Voltage on bus 5:</b>	<b>0.227289</b>
<b>Voltage on bus 6:</b>	<b>0.148733</b>
<b>Voltage on bus 7:</b>	<b>0.088491</b>
<b>:</b>	



## Current on branches

```
Current on line from bus 1 to 2 is:    20.000000
Current on line from bus 2 to 3 is:     4.009776
Current on line from bus 3 to 4 is:     2.757436
Current on line from bus 4 to 5 is:     2.046167
Current on line from bus 5 to 6 is:     4.545785
:
```

The first line is written by the call to the math library **Initialize** function and reports on the number of processors being used in the calculation. This information is useful in keeping track of the performance characteristics of different calculations. Some information from the solvers is usually printed after this. At the end of the calculation, the values of the voltages on the buses are printed out and then the current on each of the branches. The buses with externally applied voltages are also identified in the output.

## Contingency Analysis

The contingency analysis application will be described in detail since it provides a relatively compact demonstration of some of the advanced features of GridPACK. This application is built entirely around the power flow module, so it has no network component classes of its own. The main functionality is located in the **CADriver** class that consists of two methods (other than the constructor and destructor). One function is used to read in a list of contingencies and convert them to a corresponding **Contingency** data structure and the other function executes the contingency analysis calculation. These two functions will be discussed in detail.

The function for reading in the contingencies and converting them to a list of Contingency data structures has the form

```
std::vector<gridpack::powerflow::Contingency> getContingencies(
    gridpack::utility::Configuration::ChildCursors contingencies)
```

The **Contingency** data structures are defined in as part of the power flow module and exist in the **gridpack::powerflow** namespace. The list of cursors represented by the contingencies variable is obtained by the calling program before calling this function. The function itself is

```
std::vector<gridpack::powerflow::Contingency> ret;
int size = contingencies.size();
int i, idx;
gridpack::utility::StringUtils utils;
for (idx = 0; idx < size; idx++) {
    std::string ca_type;
    contingencies[idx]->get("contingencyType",&ca_type);
```

```

std::string ca_name;
contingencies[idx]->get("contingencyName",&ca_name);
if (ca_type == "Line") {
    std::string buses;
    contingencies[idx]->get("contingencyLineBuses",&buses);
    std::string names;
    contingencies[idx]->get("contingencyLineNames",&names);
    std::vector<std::string> string_vec =
        utils.blankTokenizer(buses);
    std::vector<int> bus_ids;
    for (i=0; i<string_vec.size(); i++) {
        bus_ids.push_back(atoi(string_vec[i].c_str()));
    }
    string_vec.clear();
    string_vec = utils.blankTokenizer(names);
    std::vector<std::string> line_names;
    for (i=0; i<string_vec.size(); i++) {
        line_names.push_back(utils.clean2Char(string_vec[i]));
    }
    if (bus_ids.size() == 2*line_names.size()) {
        gridpack::powerflow::Contingency contingency;
        contingency.p_name = ca_name;
        contingency.p_type = Branch;
        int i;
        for (i = 0; i < line_names.size(); i++) {
            contingency.p_from.push_back(bus_ids[2*i]);
            contingency.p_to.push_back(bus_ids[2*i+1]);
            contingency.p_ckt.push_back(line_names[i]);
            contingency.p_saveLineStatus.push_back(true);
        }
        ret.push_back(contingency);
    }
} else if (ca_type == "Generator") {
    std::string buses;
    contingencies[idx]->get("contingencyBuses",&buses);
    std::string gens;
    contingencies[idx]->get("contingencyGenerators",&gens);
    std::vector<std::string> string_vec =
        utils.blankTokenizer(buses);
    std::vector<int> bus_ids;

```

```

for (i=0; i<string_vec.size(); i++) {
    bus_ids.push_back(atoi(string_vec[i].c_str()));
}
string_vec.clear();
string_vec = utils.blankTokenizer(gens);
std::vector<std::string> gen_ids;
for (i=0; i<string_vec.size(); i++) {
    gen_ids.push_back(utils.clean2Char(string_vec[i]));
}
if (bus_ids.size() == gen_ids.size()) {
    gridpack::powerflow::Contingency contingency;
    contingency.p_name = ca_name;
    contingency.p_type = Generator;
    int i;
    for (i = 0; i < bus_ids.size(); i++) {
        contingency.p_busid.push_back(bus_ids[i]);
        contingency.p_genid.push_back(gen_ids[i]);
        contingency.p_saveGenStatus.push_back(true);
    }
    ret.push_back(contingency);
}
}
return ret;

```

This function is designed to parse input of the form

```

<?xml version="1.0" encoding="utf-8"?>
<ContingencyList>
  <Contingency_analysis>
    <Contingencies>
      <Contingency>
        <contingencyType>Line</contingencyType>
        <contingencyName>CTG1</contingencyName>
        <contingencyLineBuses> 13 14</contingencyLineBuses>
        <contingencyLineNames> B1 </contingencyLineNames>
      </Contingency>
      <Contingency>
        <contingencyType>Generator</contingencyType>
        <contingencyName>CTG2</contingencyName>
        <contingencyBuses> 2 </contingencyBuses>

```

```

        <contingencyGenerators>1 </contingencyGenerators>
    </Contingency>
</Contingencies>
</Contingency_analysis>
</ContingencyList>

```

The **contingencies** list in the argument consists of a vector of **Configuration** module cursors, each of which is pointing to one of the **Contingency** blocks in this input.

The first few lines are used to create the return list, determine the number of contingencies in the **ChildCursors** list and create a **StringUtils** object that can be used to parse the input. The function then loops over all cursors in the **contingencies** list. All contingencies should contain the **contingencyType** and **contingencyName** field, so these values are obtained using the get function from the **Configuration** module. The type can be either “**Line**” or “**Generator**”. Based on the type, the function bifurcates into two branches. The “**Line**” branch looks for the strings corresponding to **contingencyLineBuses** and **contingencyLineNames** and assigns these to the string variables **buses** and **names**. More than one transmission element may be involved in the contingency. The **StringUtils blankTokenizer** function is used to parse the **buses** string into a list of strings that can then be converted to a list of integers. These represent the original indices of the buses at each end of the branch. The **names** string is also converted to a list representing the character tag identifying the individual transmission element between the two buses. This is then reformatted to a consistent 2-character format using the **StringUtils clean2Char** function. The string vector **string\_vec** is used to hold the results from **blankTokenizer**, and the final list of integers and character tags are stored in the variables **bus\_ids** and **line\_names**. Each transmission element is characterized by two buses and a character tag, so the number of bus IDs should be twice the number of tags. If this condition is met, then the contingency is assumed to be well formed and a **Contingency** struct is created for it. After copying the data stored in the variables **ca\_type**, **ca\_name**, **bus\_ids** and **line\_names**, this contingency is added to the return variable **ret**.

The “**Generator**” branch is similar to the “**Line**” branch. The strings in the **contingencyBuses** and **contingencyGenerators** fields are copied into the string variables **buses** and **gens**. These are then converted into a list of bus IDs and generator tags using the **blankTokenizer** function and stored in the list **bus\_ids** and **gen\_ids**. A generator is characterized by the original index of the bus that it is associated with and the 2-character generator tag so the size of the **bus\_ids** and **gen\_ids** vectors must be equal. If this condition is met, then a **Contingency** struct is created, the contingency data is copied to it and the struct is added to the return variable **ret**.

After all cursor in contingencies have been processed, the **getContingencies** function returns a list of **Contingency** structs representing all the contingencies in the original XML input file.

The execute function starts with the code block

```
void gridpack::contingency_analysis::CADriver::execute(int argc,
char** argv)
{
    gridpack::parallel::Communicator world;
    gridpack::utility::CoarseTimer *timer =
        gridpack::utility::CoarseTimer::instance();
    int t_total = timer->createCategory("Total Application");
    timer->start(t_total);

    gridpack::utility::Configuration *config
        = gridpack::utility::Configuration::configuration();
    if (argc >= 2 && argv[1] != NULL) {
        char inputfile[256];
        sprintf(inputfile,"%s",argv[1]);
        config->open(inputfile,world);
    } else {
        config->open("input.xml",world);
    }
}
```

The user can pass in the name of the input file when they invoke the contingency analysis application, and this is transmitted via the variables **argc** and **argv** in the argument list. If an argument is detected, then the code will try and open a file using the argument as the filename, otherwise it will assume the input file is called “**input.xml**”. Once the input file is open, all processors have access to its contents. This section also creates a timing category for the calculation and starts the timer. The call to **CoarseTime::instance** returns the timer object and the **createCategory** call creates a timer category with the name “**Total Application**”. It also returns a handle to this category. The **start** call begins the timer. The timer can be started and stopped multiple times for the same category.

The next few lines are used to parse the input file and determine the size of the communicators that should be used to run individual tasks.

```
gridpack::utility::Configuration::CursorPtr cursor;
cursor = config->getCursor("Configuration.Contingency_analysis");
int grp_size;
```

```

double Vmin, Vmax;
if (!cursor->get("groupSize",&grp_size)) {
    grp_size = 1;
}
if (!cursor->get("minVoltage",&Vmin)) {
    Vmin = 0.9;
}
if (!cursor->get("maxVoltage",&Vmax)) {
    Vmax = 1.1;
}
gridpack::parallel::Communicator task_comm = world.divide(grp_size);

```

A **CursorPtr** is defined and set to point to the contents of the **Contingency\_analysis** block in the input file using the **getCursor** function. This block contains parameters defining some of the properties of the simulation. The **groupSize** parameter sets the size of the communicator on which individual power flow calculations are run. The power flow is not very scalable in GridPACK and it is usually fastest to run it on one processor so the default value is 1. The **minVoltage** and **maxVoltage** parameters are the limits, in p.u., for acceptable voltage variations on individual buses. Once the group size has been set, the world communicator is divided into sub communicators using the divide function. This guarantees that each subcommunicator contains at most the number of processes specified using **groupSize** (one subcommunicator may contain less than this number). Each process is now part of the world communicator and one subcommunicator.

The next block of code creates a power flow application on each task communicator and initializes it.

```

boost::shared_ptr<gridpack::powerflow::PFNetwork>
    pf_network(new gridpack::powerflow::PFNetwork(task_comm));
gridpack::powerflow::PFAppModule pf_app;
pf_app.readNetwork(pf_network, config);
pf_app.initialize();
pf_app.solve();
pf_app.ignoreVoltageViolations(Vmin, Vmax);

```

The first line creates a power flow network on the task communicator. The second line creates a power flow application. The **readNetwork** function assigns the powerflow network (which currently has nothing in it) to the power flow application, along with the pointer to the configuration module. The input file is expected to have a **Powerflow** block that contains parameters for the power flow application. These include the location of the network configuration file and the type of solver that is to be used. An example of a complete input file is

```

<?xml version="1.0" encoding="utf-8"?>
<Configuration>
  <Contingency_analysis>
    <contingencyList>contingencies.xml</contingencyList>
    <groupSize>2</groupSize>
    <maxVoltage>1.1</maxVoltage>
    <minVoltage>0.9</minVoltage>
  </Contingency_analysis>
  <Powerflow>
    <networkConfiguration> IEEE14_ca.raw </networkConfiguration>
    <maxIteration>50</maxIteration>
    <tolerance>1.0e-6</tolerance>
    <LinearSolver>
      <PETScOptions>
        -ksp_type richardson
        -pc_type lu
        -pc_factor_mat_solver_package superlu_dist
        -ksp_max_it 1
      </PETScOptions>
    </LinearSolver>
  </Powerflow>
</Configuration>

```

Note that it has two blocks, **Contingency\_analysis** and **Powerflow**. The parameters describing the contingency calculation and the location of the contingencies are located in the first block and the power flow parameters are located in the second block. The **readNetwork** function will read in the network configuration file and partition the network. The **initialize** function is used to initialize the network components from the **DataCollection** objects and assign exchange buffers. The call to **solve** is used to obtain a power solution to the base problem with no contingencies. Since all tasks have the same data at this point, the network solution is duplicated across all subcommunicators. The final call to **ignoreVoltageViolations** sets a parameter in each network component that violates the voltage bounds for base case. These components will be ignored in any subsequent checks for voltage violations.

The next step is to read in the contingencies and convert these to a list of contingency data structs.

```

std::string contingencyfile;
if (!cursor->get("contingencyList",&contingencyfile)) {
  contingencyfile = "contingencies.xml";
}

```

```

}
bool ok = config->open(contingencyfile,world);
cursor = config->getCursor(
    "ContingencyList.Contingency_analysis.Contingencies");
gridpack::utility::Configuration::ChildCursors contingencies;
if (cursor) cursor->children(contingencies);
std::vector<gridpack::powerflow::Contingency>
    events = getContingencies(contingencies);
if (world.rank() == 0) {
    int idx;
    for (idx = 0; idx < events.size(); idx++) {
        printf("Name: %s\n",events[idx].p_name.c_str());
        if (events[idx].p_type == Branch) {
            int nlines = events[idx].p_from.size();
            int j;
            for (j=0; j<nlines; j++) {
                printf(" Line: (from) %d (to) %d (line) \'%s\'\n",
                    events[idx].p_from[j],events[idx].p_to[j],
                    events[idx].p_ckt[j].c_str());
            }
        } else if (events[idx].p_type == Generator) {
            int nbus = events[idx].p_busid.size();
            int j;
            for (j=0; j<nbus; j++) {
                printf(" Generator: (bus) %d (generator ID) \'%s\'\n",
                    events[idx].p_busid[j],events[idx].p_genid[j].c_str());
            }
        }
    }
}
}
}

```

The location of the contingency file is contained in the **contingencyList** field in the input file. If this field is not present, the code defaults to the file name **contingencies.xml**. The contingency file is then opened using the **open** function in the **Configuration** module and a cursor is set to the **Contingencies** block within this file. The **Configuration** **children** function returns a list of cursor pointers that point to each of the individual **Contingency** blocks. The **getContingencies** function described above parses each of these blocks and returns a vector of contingency data structs. The contingency list is replicated on all processors. Process 0 is used to provide a listing of the contingencies to standard output by looping over the **events** vector returned by the **getContingencies** function.



Once the contingencies have been determined, the code next sets up a task manager on the world communicator and sets the number of tasks equal to the number of contingencies.

```
gridpack::parallel::TaskManager taskmgr(world);
int ntasks = events.size();
taskmgr.set(ntasks);
```

The task loop is created by defining a `task_id` variable and a character string buffer that is used inside the loop to create messages. The task manager then begins iterating over different tasks.

```
int task_id;
char sbuf[128];
while (taskmgr.nextTask(task_comm, &task_id)) {
    printf("Executing task %d on process %d\n", task_id, world.rank());
```

The call to `nextTask` takes the task communicator as one of its arguments so the value of `task_id` that is returned is the same for all processors on the communicator. This guarantees that each of the processors in this copy of the power flow application is working on the same contingency. If the `nextTask` function returns false, the tasks have been completed and the code exits from the `while` loop. At the start of the task, the code prints out a statement to standard out describing which tasks are being executed by each processor.

The next few lines in the task loop are used to open a file so that the output from each task is directed to a separate file. This can be used later to examine individual tasks.

```
    sprintf(sbuf, "%s.out", events[task_id].p_name.c_str());
    pf_app.open(sbuf);
    sprintf(sbuf, "\nRunning task on %d processes\n", task_comm.size());
    pf_app.writeHeader(sbuf);
    if (events[task_id].p_type == Branch) {
        int nlines = events[task_id].p_from.size();
        int j;
        for (j=0; j<nlines; j++) {
            sprintf(sbuf, " Line: (from) %d (to) %d (line) \'%s\'\n",
                events[task_id].p_from[j], events[task_id].p_to[j],
                events[task_id].p_ckt[j].c_str());
        }
    } else if (events[task_id].p_type == Generator) {
        int nbus = events[task_id].p_busid.size();
        int j;
        for (j=0; j<nbus; j++) {
            sprintf(sbuf, " Generator: (bus) %d (generator ID) \'%s\'\n",
```

```

        events[task_id].p_busid[j],
        events[task_id].p_genid[j].c_str());
    }
}
pf_app.writeHeader(sbuf);

```

The first line is used to create a name for the output file using the contingency name. The output from the power flow calculation is then redirected to this file using the power flow **open** function. Next, some information about this particular contingency is written to the file using some calls to the `writeHeader` method. This includes the number of processors used to calculate the contingency and the details of the contingency itself.

The remaining lines in the while loop are used to solve the power flow equations.

```

pf_app.resetVoltages();
pf_app.setContingency(events[task_id]);
if (pf_app.solve()) {
    pf_app.write();
    bool ok = pf_app.checkVoltageViolations(Vmin,Vmax);
    ok = ok & pf_app.checkLineOverloadViolations();
    if (ok) {
        sprintf(sbuf,"\nNo violation for contingency %s\n",
            events[task_id].p_name.c_str());
    } else {
        sprintf(sbuf,"\nViolation for contingency %s\n",
            events[task_id].p_name.c_str());
    }
    pf_app.print(sbuf);
}
pf_app.unSetContingency(events[task_id]);
pf_app.close();
}

```

Before doing the calculation, all voltages are returned to the original values defined in the network configuration file using **resetVoltages**. The contingency parameters are set to the values specified by the **task\_id** element in the **events** list using the **setContingency** method.

The system is then solved using the power flow **solve** function. If the solution succeeds, the calculation writes out the voltages and branch power flow values to the output file. The calculation also checks for voltage violations and line overload violations. The results of these checks are written to the output file for each power flow calculation. After this is complete, the

powerflow calculation returns all contingency related parameters to their original values using **unSetContingency** and closes the output file. This is repeated until all contingencies in the event list have been evaluated.

At this point, the contingency application is essentially complete. The remaining lines of code

```
taskmgrprintStats();
timer->stop(t_total);
if (events.size()*grp_size >= world.size()) {
    timer->dump();
}
```

are used to print out a list of how many tasks were evaluated on each processor and to stop the timing of the “**Total Application**” category. The timer **dump** method will print statistics on the amount of time spent in the total application as well as reporting timings inside the power flow application. The check on the **dump** call is to verify that all processors have participated in at least one power flow calculation.

## Fortran 2003 Interface

GridPACK has developed a Fortran interface that can be used to access most of the functionality in the framework modules. The Fortran interface makes extensive use of the object-oriented features in Fortran, so a compiler that supports the Fortran 2003 standard must be used if creating Fortran applications. The Fortran compiler must also support the `iso_c_binding` module, but this will usually be available if the compiler supports Fortran 2003. Most recent compilers support Fortran 2003. A working power flow application written entirely in Fortran has been included in the current release and demonstrates how to use the Fortran interface. The Fortran implementation is very similar to the C++ interface and most of the C++ documentation applies to the corresponding Fortran functionality. The remainder of this section will highlight the important differences between the C++ and Fortran interfaces.

Because Fortran does not have any support for templates (that we know of), the Fortran interface cannot support multiple different kinds of networks within a single application. This means that only one bus and one branch class can be present in an application, so the bus and branch classes must support all possible types of behavior. It is still possible to have more than one network in an application, but all networks must be of the same type.

The bus and branch classes in the Fortran interface are represented by the Fortran derived types **application\_bus** and **application\_branch**. These types have procedures bound to them, as well as internal data elements. These types are defined in the Fortran file **component\_template.F90** file that is located in the **fortran/component** directory. The application bus and branch classes can be created by modifying a copy of **component\_template.F90**. The functions in the math-vector interface and the component

base classes are all defined in this file, along with default implementations for these functions. Additional data elements and procedures can be added to the bus and branch data types to create appropriate functionality for specific problems.

A brief overview of the **application\_bus** type in the **component\_template.F90** file is provided here. Similar considerations apply to the **application\_branch** type. To use the **component\_template.F90** file it should first be copied to the directory where the application source code resides and renamed to something appropriate. We will use the name **app\_component.F90**. Inside the component file, the Fortran types **bus\_xc\_data**, **branch\_xc\_data**, **application\_bus**, **application\_branch** are defined as part of the **application\_components** module. These are the only types that need concern the application developer. There are also two types defined in this file called **application\_bus\_wrapper** and **application\_branch\_wrapper**. These are only used internally but must be defined in this file. They should not be modified. There is a line at the bottom of the **app\_component.F90** file that includes an external file **component\_inc.F90**. This file contains many functions that are required by the interface and must be included in the **application\_components** module. However, these functions should not be modified by the user so to avoid possible errors and to simplify the file somewhat, these functions are put in an include file.

The **application\_bus** type has four parts. These consist of 1) application-specific data elements, 2) data elements that must be defined in order for the component to interact with rest of the framework, 3) application-specific functions that are defined by the user and 4) framework functions that must be included in the component. The framework functions all have base implementations can be modified to suit the application. The only data elements that must be included in the **application\_bus** type is a variable of type **bus\_xc\_data** and a pointer to this variable. The **bus\_xc\_data** type will be discussed further below and represents all data that might need to be exchanged in a bus update.

The framework functions are directly analogous to the functions defined for the C++ implementation and users should refer to the documentation above to find out how these functions work. This section will primarily discuss differences between the Fortran and C++ interfaces. The Fortran compilers do not have the same name-mangling capabilities as C++ so all function names are preceded by either a **bus\_** or **branch\_** to distinguish between bus and branch versions of the functions. A few functions only appear in the bus class or the branch class and do not necessarily need this prefix, but to be consistent, this convention is used for all functions.

Functions that are bound to the **application\_bus** type are already listed in the **component\_template.F90**. These functions consist of both a declaration within the **application\_bus** type and a function or subroutine implementation within the

**application\_components** module. The declarations within the **application\_bus** type (after the **contains** keyword) have the form

```
procedure::bus_matrix_diag_size
procedure::bus_matrix_diag_values
procedure::bus_matrix_forward_size
procedure::bus_matrix_reverse_size
:
```

The **procedure** keyword distinguishes a function or subroutine bound to the Fortran type from a piece of data (which is declared as a data type using one of the intrinsic Fortran data types or a Fortran type declaration).

After the type declarations within the **applications\_components** module, there is a **contains** keyword followed by the subroutine and function implementations for all the declared procedures. The original implementations in the **component\_template.F90** file are just stubs for these functions and typically don't do much. An example is the **bus\_matrix\_diag\_size** function which originally has the implementation

```
logical function bus_matrix_diag_size(bus, isize, jsize)
  implicit none
  class(application_bus), intent(in) :: bus
  integer, intent(out) :: isize, jsize
  bus_matrix_diag_size = .false.
  return
end function bus_matrix_diag_size
```

The initial implementation just returns false if this function is invoked and doesn't set the variables **isize** or **jsize**. Note the first item in the argument list. This is declared as being of type **class(application\_bus)** with intent in. All functions and subroutines that are bound to the **application\_bus** type must have this argument, even if they do not have any other arguments. This argument provides a mechanism for accessing data items or functions that are related to a particular **application\_bus** instance.

To see how the bus argument is used in actual practice, an implementation of this function in a power flow application is shown below

```
logical function bus_matrix_diag_size(bus, isize, jsize)
  implicit none
  class(application_bus), intent(in) :: bus
  integer, intent(out) :: isize, jsize
  isize = 1
```

```

jsize = 1
bus_matrix_diag_size = .true.
if (bus%p_mode.eq.JACOBIAN) then
  if (.not.bus%bus_is_isolated()) then
    isize = 2
    jsize = 2
    bus_matrix_diag_size = .true.
  else
    bus_matrix_diag_size = .false.
  endif
else if (bus%p_mode.eq.YBUS) then
  if (.not.bus%bus_is_isolated()) then
    bus_matrix_diag_size = .true.
    isize = 1
    jsize = 1
  else
    bus_matrix_diag_size = .false.
  endif
  return
endif
return
end function bus_matrix_diag_size

```

The **application\_bus** implementation for power flow contains the variable **p\_mode** and a user-specified function **bus\_is\_isolated** (this is declared as a type-bound procedure). To access this data and this function inside a type-bound procedure, use the Fortran “%” symbol. The **bus** variable in the argument list is acting in a similar way to the “**this**” pointer in C++ and refers back to the **application\_bus** instance that made the original call to **bus\_matrix\_diag\_size**. Although the **bus\_is\_isolated** function implementation has the variable **bus** in its argument list, it doesn’t need to explicitly pass this argument when making a call from an **application\_bus** instance. The **bus** argument is assumed in this case. Similarly, a call to the **bus\_matrix\_diag\_size** function, which has additional arguments, would have the form

```
ok = bus%bus_matrix_diag_size(isize,jsize)
```

Following this syntax, it is possible to construct a complete set of functions for an arbitrary application. Additional application-specific functions can be added to the component types by declaring them as procedures within the type and adding their implementations to the **application\_components** module.

There are a few procedures in both the bus and branch types that should not be modified. No stubs for these appear in the component\_template.F90 file. For the application\_bus type, these procedures are

```
procedure::bus_get_neighbor_branch
procedure::bus_get_neighbor_bus
procedure::bus_get_xc_buf_size
procedure::bus_get_xc_buf
```

For the application\_branch type, the procedures are

```
procedure::branch_get_bus1
procedure::branch_get_bus2
procedure::branch_get_xc_buf_size
procedure::branch_get_xc_buf
```

These procedures are required by other parts of the framework, but should not be modified by the user. Some other procedures are defined in the base class and do not appear as procedure declarations in application\_bus and application\_branch types. These procedures include

```
procedure::bus_get_num_neighbors
procedure::bus_set_reference_bus
procedure::bus_get_reference_bus
procedure::bus_get_original_index
procedure::bus_compare
```

for buses and

```
procedure::branch_get_bus1_original_index
procedure::branch_get_bus2_original_index
procedure::branch_compare
```

for branches. The bus and branch compare functions are used to determine if a bus or branch is equal to another bus or branch. An example of how to use this function can be found in the function that evaluates transformer contributions on branches for the power flow application. The syntax for calling this function is

```
double complex function branch_get_transformer(branch, bus)
:
class(application_branch), intent(in) :: branch
class(application_bus), intent(in) :: bus
class(application_bus), pointer :: bus1, bus2
:
```

```

if (bus%bus_compare(bus1)) then
:

```

In this fragment, the **bus\_compare** function is being used to check if bus1 is equivalent to bus. The **branch\_compare** function is used in a similar way.

The final issue in implementing the Fortran application bus and branch classes is understanding the exchange buffers. These buffers are declared at the top of the **component\_template.F90** file as the **bus\_xc\_data** and **branch\_xc\_data** data types. Although the underlying Fortran interface implementation makes extensive use of the **iso\_c\_binding** module, we have worked very hard to keep the **iso\_c\_binding** data types out of the Fortran interface itself. However, the one place where this is not possible is in the exchange buffers, so it is important to use these data type declarations for any variables that are included in the exchange buffers. The exchange buffers are declared as follows in the top of the **component\_template.F90** file

```

type, bind(c), public :: bus_xc_data
!
! Example data types. Replace with application-specific values
!
integer(C_INT) int_reg
integer(C_LONG) int_long
real(C_FLOAT) real_s
real(C_DOUBLE) real_d
complex(C_FLOAT_COMPLEX) complex_s
complex(C_DOUBLE_COMPLEX) complex_d
logical(C_BOOL) log_reg
end type

```

The variables **int\_reg**, **int\_long**, **real\_s**, **real\_d**, **complex\_s**, **complex\_d** and **log\_reg** are just examples and should be replaced with the variables used in the actual application. Not all data types will be used in an application. Any buffer variables used in an application should use the **iso\_c\_binding** type declarations (**C\_INT**, **C\_LONG**, **C\_FLOAT**, **C\_FLOAT\_COMPLEX**, **C\_DOUBLE\_COMPLEX**, **C\_BOOL**). Variables declared with the **iso\_c\_binding** types can be cast to regular Fortran variables by relying on the compiler to automatically cast an assignment to the right sized variable. For example

```

integer f_var
integer(C_INT) c_var
:
f_var = c_var

```



If **f\_var** is an 8 byte integer and **c\_var** is a 4 byte integer, the compiler can be relied on to do the cast. This also works in the opposite direction, assuming that **f\_var** does not exceed the capacity of a 4 byte variable.

The functions that access neighboring branches or buses also work differently than the corresponding C++ functions. Fortran does not support anything that looks like an STL vector so neighbors are accessed from buses using a two step process. The first step is to get the total number of neighbors attached to the bus using the **bus\_get\_num\_neighbors** procedure. This allows users to set up a loop that can be used to run over either the neighboring branches or the neighboring buses that are attached to the calling bus via a single branch. The neighboring branches can then be accessed by using the **bus\_get\_neighbor\_branch** function which returns a Fortran pointer to the neighboring branch. The syntax for using this function is

```
integer i, nbranch
type(application_branch), pointer :: branch
nbranch = bus%bus_get_num_neighbors()
do i = 1, nbranch
    branch => bus%bus_get_neighbor_branch(i)
end do
```

The **bus\_get\_neighbor\_bus** function works in a similar way and returns a pointer to the bus at the other end of branch **i**. To get pointers to the buses at either end of a branch, use the functions **branch\_get\_bus1** and **branch\_get\_bus2** procedures. Because the Fortran interface only supports one type of bus or branch per application, these functions return pointers of the correct type and there is no need to cast them to something else.

Most of the remaining differences between the Fortran and C++ interfaces are associated with the GridPACK factory module. As with the component classes, the Fortran interface only supports one kind of factory. This is the **app\_factory** type and it can be created by copying the **factory\_template.F90** file in the **fortran/factory** directory and making application-specific changes to it. The factory base class contains the functions

```
procedure::set_components
procedure::load
procedure::set_exchange
procedure::set_mode
procedure::check_true
```

These functions behave the same way as the equivalent C++ functions. In addition, the **app\_factory** type contains the two functions

```
procedure::create
```

## **procedure::destroy**

Because Fortran does not support constructors and destructors in the same way as C++, it is necessary to create explicit functions that implement whatever behaviors are imbedded in the C++ constructors and destructors. This is accomplished in the Fortran interface by adding **create** and **destroy** functions (or **initialize** and **finalize** functions) to most of the Fortran implementations of the GridPACK modules.

Additional methods can be added to the **app\_factory** type to support application-specific functionality. An example of how to do this is the **set\_y\_bus** procedure for the power flow application. This subroutine is declared as a procedure in the **app\_factory** type. The implementation is written as

```
subroutine set_y_bus(factory)
  class(app_factory), intent(in) :: factory
  class(application_bus), pointer :: bus
  class(application_branch), pointer :: branch
  class(network), pointer :: grid
  integer nbus, nbranch, i
  grid => factory%p_network_int
  nbus = grid%num_buses()
  nbranch = grid%num_branches()
  do i = 1, nbus
    bus => bus_cast(grid%get_bus(i))
    call bus%bus_set_y_matrix()
  end do
  do i = 1, nbranch
    branch => branch_cast(grid%get_branch(i))
    call branch%branch_set_y_matrix()
  end do
  return
end subroutine set_y_bus
```

The functions for accessing the bus and branch objects work differently from the functions that get neighboring branches or buses in the component classes. The neighbor bus and branch functions return a pointer to the appropriate bus or branch directly to the calling application. The **get\_bus** and **get\_branch** functions in the Fortran network class return an opaque object that cannot be directly used in a Fortran code. To convert this to a bus or branch pointer it is necessary to call the **bus\_cast** or **branch\_cast** functions which return a pointer that can be called in Fortran.

The last remaining point is to provide a list of the existing Fortran modules that need to be used in a GridPACK application using the Fortran interface. These modules need to be included in any subroutine or function that is using the associated Fortran types. The existing modules are

```
gridpack_network ! type or class network
application_factory ! type or class app_factory
application_components ! type or class application_bus and
                        ! application_branch
gridpack_configuration ! type or class cursor
gridpack_full_matrix_map ! type or class full_matrix_map
gridpack_bus_vector_map ! type or class bus_vector_map
gridpack_gen_matrix_map ! type or class gen_matrix_map
gridpack_gen_vector_map ! type or class gen_vector_map
gridpack_math ! access to math initialization and
               ! finalization routines
gridpack_matrix ! type or class matrix
gridpack_vector ! type or class vector
gridpack_linear_solver ! type or class linear_solver
gridpack_nonlinear_solver ! type or class funcbuilder
                        ! and nonlinear_solver
gridpack_communicator ! type or class communicator
gridpack_parallel ! access to parallel initialization
                  ! and finalization routines
gridpack_parser ! class or type pti23_parser
gridpack_serial_io ! class or type bus_serial_io
                  ! and branch_serial_io
```

The appropriate module should be included in any function or subroutine that uses objects defined in the module. Modules can be included using the standard Fortran “**use**” statement.