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## **Preface**

This book provides a quick introduction to Pyomo, which includes a collection of Python software packages that supports a diverse set of optimization capabilities for formulating and analyzing optimization models. A central component of Pyomo is Pyomo, which supports the formulation and analysis of mathematical models for complex optimization applications. This capability is commonly associated with algebraic modeling languages (AMLs), which support the description and analysis of mathematical models with a high-level language. Although most AMLs are implemented in custom modeling languages, Pyomo's modeling objects are embedded within Python, a full-featured high-level programming language that contains a rich set of supporting libraries.

Pyomo has also proven an effective framework for developing high-level optimization and analysis tools. For example, the PySP package provides generic solvers for stochastic programming. PySP leverages the fact that Pyomo's modeling objects are embedded within a full-featured high-level programming language, which allows for transparent parallelization of subproblems using Python parallel communication libraries.

### Goals of the Book

This book provides a broad overview of different components of the Pyomo software. There are roughly two main goals for this book:

- 1. Help users get started with different Pyomo capabilities. Our goal is not to provide a comprehensive reference, but rather to provide a tutorial with simple and illustrative examples. Also, we aim to provide explanations behind the design and philosophy of Pyomo.
- 2. Provide preliminary documentation of new features and capabilities. We know that a new feature or capability probably will not be used unless it is documented. As Pyomo evolves, we plan to use this book to document these features. This provides users some context concerning the focus of Pyomo development, and it also provides an opportunity to get early feedback on new features before they are documented in other contexts.

## Who Should Read This Book

This book is intended to be a reference for students, academic researchers and practitioners. Pyomo has been effectively used in the classroom with undergraduate and graduate students. However, we assume that the reader is generally familiar with optimization and mathematical modeling. Although this book does not contain a glossary, we recommend the Mathematical Programming Glossary [MPG] as a reference for the reader. We also assume that the reader is generally familiar with the Python programming language. There are a variety of books describing Python, as well as excellent documentation of the Python language and the software packages bundled with Python distributions.

## **Comments and Questions**

The Pyomo home page provides resources for Pyomo users:

http://pyomo.org

Pyomo development is hosted by Sandia National Laboratories and COIN-OR:

- https://software.sandia.gov/pyomo
- https://projects.coin-or.org/Pyomo

See the Pyomo Forum for online discussions of Pyomo:

• http://groups.google.com/group/pyomo-forum/

We welcome feedback on typos and errors in our examples, as well as comments on the presentation of this material. Good Luck!

# **Citing Pyomo**

## 1.1 Pyomo

Hart, William E., Jean-Paul Watson, and David L. Woodruff. "Pyomo: modeling and solving mathematical programs in Python." Mathematical Programming Computation 3, no. 3 (2011): 219-260.

Hart, William E., Carl Laird, Jean-Paul Watson, and David L. Woodruff. Pyomo – Optimization Modeling in Python. Vol. 67. Springer, 2012.

## 1.2 PySP

Watson, Jean-Paul, David L. Woodruff, and William E. Hart. "PySP: modeling and solving stochastic programs in Python." Mathematical Programming Computation 4, no. 2 (2012): 109-149.

## **Release Notes**

## 2.1 Highlights

The following are highlights of the Pyomo 4.1 release:

- Modeling
  - API changes for model transformations
  - Revised API for SOSConstraint, Suffix and Block components
  - Optimization results are now loaded into models
  - Removed explicit specification of model preprocessing
- Solvers
  - Resolved many issues with writing and solving MPECs
  - Changes to MPEC meta-solver names
  - The solution output for runph has been changed to
- Other
  - Pyomo subcommands can now use configuration files (e.g. pyomo solve config.json)
  - New JSON/YAML format for parameter data
  - Added a script to install pyomo.extras

## 2.2 Determining Your Version of Pyomo

To determine your current version of Pyomo, use the command

pyomo --version

# **Pyomo Overview**

## 3.1 Mathematical Modeling

This chapter provides an introduction to Pyomo: Python Optimization Modeling Objects. A more complete description is contained in the Pyomo book. Pyomo supports the formulation and analysis of mathematical models for complex optimization applications. This capability is commonly associated with algebraic modeling languages (AMLs) such as AMPL [AMPL] AIMMS [AIMMS] and GAMS [GAMS]. Pyomo's modeling objects are embedded within Python, a full-featured high-level programming language that contains a rich set of supporting libraries.

Modeling is a fundamental process in many aspects of scientific research, engineering and business. Modeling involves the formulation of a simplified representation of a system or real-world object. Thus, modeling tools like Pyomo can be used in a variety of ways:

- Explain phenomena that arise in a system,
- Make predictions about future states of a system,
- · Assess key factors that influence phenomena in a system,
- · Identify extreme states in a system, that might represent worst-case scenarios or minimal cost plans, and
- Analyze trade-offs to support human decision makers.

Mathematical models represent system knowledge with a formalized mathematical language. The following mathematical concepts are central to modern modeling activities:

#### variables

Variables represent unknown or changing parts of a model (e.g. whether or not to make a decision, or the characteristic of a system outcome). The values taken by the variables are often referred to as a *solution* and are usually an output of the optimization process.

#### parameters

Parameters represents the data that must be supplied to perform the optimization. In fact, in some settings the word *data* is used in place of the word *parameters*.

### relations

These are equations, inequalities or other mathematical relationships that define how different parts of a model are connected to each other.

#### goals

These are functions that reflect goals and objectives for the system being modeled.

The widespread availability of computing resources has made the numerical analysis of mathematical models a commonplace activity. Without a modeling language, the process of setting up input files, executing a solver and extracting the final results from the solver output is tedious and error prone. This difficulty is compounded in complex, large-scale real-world applications which are difficult to debug when errors occur. Additionally, there are many different formats used by optimization software packages, and few formats are recognized by many optimizers. Thus the application of multiple optimization solvers to analyze a model introduces additional complexities.

Pyomo is an AML that extends Python to include objects for mathematical modeling. Hart et al. Pyomo, [PyomoJournal] compare Pyomo with other AMLs. Although many good AMLs have been developed for optimization models, the following are motivating factors for the development of Pyomo:

### **Open Source**

Pyomo is developed within Pyomo's open source project to promote transparency of the modeling framework and encourage community development of Pyomo capabilities.

#### **Customizable Capability**

Pyomo supports a customizable capability through the extensive use of plug-ins to modularize software components.

### **Solver Integration**

Pyomo models can be optimized with solvers that are written either in Python or in compiled, low-level languages.

#### **Programming Language**

Pyomo leverages a high-level programming language, which has several advantages over custom AMLs: a very robust language, extensive documentation, a rich set of standard libraries, support for modern programming features like classes and functions, and portability to many platforms.

### 3.2 Overview of Modeling Components and Processes

Pyomo supports an object-oriented design for the definition of optimization models. The basic steps of a simple modeling process are:

- Create model and declare components
- · Instantiate the model
- · Apply solver
- Interrogate solver results

In practice, these steps may be applied repeatedly with different data or with different constraints applied to the model. However, we focus on this simple modeling process to illustrate different strategies for modeling with Pyomo.

A Pyomo *model* consists of a collection of modeling *components* that define different aspects of the model. Pyomo includes the modeling components that are commonly supported by modern AMLs: index sets, symbolic parameters, decision variables, objectives, and constraints. These modeling components are defined in Pyomo through the following Python classes:

### Set

set data that is used to define a model instance

#### **Param**

parameter data that is used to define a model instance

### Var

decision variables in a model

#### **Objective**

expressions that are minimized or maximized in a model

### Constraint

constraint expressions that impose restrictions on variable values in a model

### 3.3 Abstract Versus Concrete Models

A mathematical model can be defined using symbols that represent data values. For example, the following equations represent a linear program (LP) to find optimal values for the vector x with parameters n and b, and parameter vectors a and c:

$$\begin{array}{ll}
\min & \sum_{j=1}^{n} c_j x_j \\
\text{s.t.} & \sum_{j=1}^{n} a_{ij} x_j \ge b_i \quad \forall i = 1 \dots m \\
& x_j \ge 0 \quad \forall j = 1 \dots n
\end{array}$$

#### Note

As a convenience, we use the symbol  $\forall$  to mean "for all" or "for each."

We call this an *abstract* or *symbolic* mathematical model since it relies on unspecified parameter values. Data values can be used to specify a *model instance*. The AbstractModel class provides a context for defining and initializing abstract optimization models in Pyomo when the data values will be supplied at the time a solution is to be obtained.

In some contexts a mathematical model can be directly defined with the data values supplied at the time of the model definition and built into the model. We call these *concrete* mathematical models. For example, the following LP model is a concrete instance of the previous abstract model:

min 
$$2x_1 + 3x_2$$
  
s.t.  $3x_1 + 4x_2 \ge 1$   
 $x_1, x_2 \ge 0$ 

The ConcreteModel class is used to define concrete optimization models in Pyomo.

## 3.4 A Simple Abstract Pyomo Model

We repeat the abstract model already given:

min 
$$\sum_{j=1}^{n} c_j x_j$$
  
s.t.  $\sum_{j=1}^{n} a_{ij} x_j \ge b_i$   $\forall i = 1 \dots m$   
 $x_j \ge 0$   $\forall j = 1 \dots n$ 

One way to implement this in Pyomo is as follows:

```
__future_
               _ import division
from pyomo.environ import *
model = AbstractModel()
model.m = Param(within=NonNegativeIntegers)
model.n = Param(within=NonNegativeIntegers)
model.I = RangeSet(1, model.m)
model.J = RangeSet(1, model.n)
model.a = Param(model.I, model.J)
model.b = Param(model.I)
model.c = Param(model.J)
# the next line declares a variable indexed by the set J
model.x = Var(model.J, domain=NonNegativeReals)
def obj_expression(model):
    return summation (model.c, model.x)
model.OBJ = Objective(rule=obj_expression)
```

```
def ax_constraint_rule(model, i):
    # return the expression for the constraint for i
    return sum(model.a[i,j] * model.x[j] for j in model.J) >= model.b[i]

# the next line creates one constraint for each member of the set model.I
model.AxbConstraint = Constraint(model.I, rule=ax_constraint_rule)
```

#### Note

Python is interpreted one line at a time. A line continuation character, backslash, is used for Python statements that need to span multiple lines. In Python, indentation has meaning and must be consistent. For example, lines inside a function definition must be indented and the end of the indentation is used by Python to signal the end of the definition.

We will now examine the lines in this example. The first import line is used to ensure that int or long division arguments are converted to floating point values before division is performed.

```
from __future__ import division
```

In Python versions before 3.0, division returns the floor of the mathematical result of division if arguments are int or long. This import line avoids unexpected behavior when developing mathematical models with integer values.

The next import line that is required in every Pyomo model. Its purpose is to make the symbols used by Pyomo known to Python.

```
from pyomo.environ import *
```

The declaration of a model is also required. The use of the name model is not required. Almost any name could be used, but we will use the name model most of the time in this book. In this example, we are declaring that it will be an abstract model.

```
model = AbstractModel()
```

We declare the parameters m and n using the Pyomo Param function. This function can take a variety of arguments; this example illustrates use of the within option that is used by Pyomo to validate the data value that is assigned to the parameter. If this option were not given, then Pyomo would not object to any type of data being assigned to these parameters. As it is, assignment of a value that is not a non-negative integer will result in an error.

```
model.m = Param(within=NonNegativeIntegers)
model.n = Param(within=NonNegativeIntegers)
```

Although not required, it is convenient to define index sets. In this example we use the RangeSet function to declare that the sets will be a sequence of integers starting at 1 and ending at a value specified by the the parameters model.m and model.n.

```
model.I = RangeSet(1, model.m)
model.J = RangeSet(1, model.n)
```

The coefficient and right-hand-side data are defined as indexed parameters. When sets are given as arguments to the Param function, they indicate that the set will index the parameter.

```
model.a = Param(model.I, model.J)
model.b = Param(model.I)
model.c = Param(model.J)
```

#### Note

In Python, and therefore in Pyomo, any text after pound sign is considered to be a comment.

The next line interpreted by Python as part of the model declares the variable x. The first argument to the Var function is a set, so it is defined as an index set for the variable. In this case the variable has only one index set, but multiple sets could be used as was the case for the declaration of the parameter model.a. The second argument specifies a domain for the variable. This information is part of the model and will passed to the solver when data is provided and the model is solved. Specification of the NonNegativeReals domain implements the requirement that the variables be greater than or equal to zero.

```
# the next line declares a variable indexed by the set J
model.x = Var(model.J, domain=NonNegativeReals)
```

In abstract models, Pyomo expressions are usually provided to objective function and constraint declarations via a function defined with a Python def statement. The def statement establishes a name for a function along with its arguments. When Pyomo uses a function to get objective function or constraint expressions, it always passes in the model (i.e., itself) as the the first argument so the model is always the first formal argument when declaring such functions in Pyomo. Additional arguments, if needed, follow. Since summation is an extremely common part of optimization models, Pyomo provides a flexible function to accommodate it. When given two arguments, the summation function returns an expression for the sum of the product of the two arguments over their indexes. This only works, of course, if the two arguments have the same indexes. If it is given only one argument it returns an expression for the sum over all indexes of that argument. So in this example, when summation is passed the arguments model.x it returns an internal representation of the expression  $\sum_{i=1}^{n} c_i x_i$ .

```
def obj_expression(model):
    return summation(model.c, model.x)
```

To declare an objective function, the Pyomo function called Objective is used. The rule argument gives the name of a function that returns the expression to be used. The default *sense* is minimization. For maximization, the sense=maximize argument must be used. The name that is declared, which is OBJ in this case, appears in some reports and can be almost any name.

```
model.OBJ = Objective(rule=obj_expression)
```

Declaration of constraints is similar. A function is declared to deliver the constraint expression. In this case, there can be multiple constraints of the same form because we index the constraints by i in the expression  $\sum_{j=1}^{n} a_{ij}x_j \ge b_i \ \forall i=1...m$ , which states that we need a constraint for each value of i from one to m. In order to parametrize the expression by i we include it as a formal parameter to the function that declares the constraint expression. Technically, we could have used anything for this argument, but that might be confusing. Using an i for an i seems sensible in this situation.

```
def ax_constraint_rule(model, i):
    # return the expression for the constraint for i
    return sum(model.a[i,j] * model.x[j] for j in model.J) >= model.b[i]
```

#### Note

In Python, indexes are in square brackets and function arguments are in parentheses.

In order to declare constraints that use this expression, we use the Pyomo Constraint function that takes a variety of arguments. In this case, our model specifies that we can have more than one constraint of the same form and we have created a set, model.I, over which these constraints can be indexed so that is the first argument to the constraint declaration function. The next argument gives the rule that will be used to generate expressions for the constraints. Taken as a whole, this constraint declaration says that a list of constraints indexed by the set model.I will be created and for each member of model.I, the function ax\_constraint\_rule will be called and it will be passed the model object as well as the member of model.I.

```
# the next line creates one constraint for each member of the set model.I
model.AxbConstraint = Constraint(model.I, rule=ax_constraint_rule)
```

In the object oriented view of all of this, we would say that model object is a class instance of the AbstractModel class, and model. J is a Set object that is contained by this model. Many modeling components in Pyomo can be optionally specified as *indexed components*: collections of components that are referenced using one or more values. In this example, the parameter model.c is indexed with set model.J.

In order to use this model, data must be given for the values of the parameters. Here is one file that provides data.

```
# one way to input the data in AMPL format
# for indexed parameters, the indexes are given before the value

param m := 1;
param n := 2;

param a :=
1 1 3
1 2 4;

param c:=
1 2
2 3;

param b := 1 1;
```

There are multiple formats that can be used to provide data to a Pyomo model, but the AMPL format works well for our purposes because it contains the names of the data elements together with the data. In AMPL data files, text after a pound sign is treated as a comment. Lines generally do not matter, but statements must be terminated with a semi-colon.

For this particular data file, there is one constraint, so the value of model.m will be one and there are two variables (i.e., the vector model.x is two elements long) so the value of model.n will be two. These two assignments are accomplished with standard assignments. Notice that in AMPL format input, the name of the model is omitted.

```
param m := 1 ;
param n := 2 ;
```

There is only one constraint, so only two values are needed for model.a. When assigning values to arrays and vectors in AMPL format, one way to do it is to give the index(es) and the the value. The line 1 2 4 causes model.a[1,2] to get the value 4. Since model.c has only one index, only one index value is needed so, for example, the line 1 2 causes model.c[1] to get the value 2. Line breaks generally do not matter in AMPL format data files, so the assignment of the value for the single index of model.b is given on one line since that is easy to read.

```
param a :=
1 1 3
1 2 4
;

param c:=
1 2
2 3
;

param b := 1 1;
```

When working with Pyomo (or any other AML), it is convenient to write abstract models in a somewhat more abstract way by using index sets that contain strings rather than index sets that are implied by  $1, \ldots, m$  or the summation from 1 to n. When this is done, the size of the set is implied by the input, rather than specified directly. Furthermore, the index entries may have no real order. Often, a mixture of integers and indexes and strings as indexes is needed in the same model. To start with an illustration of general indexes, consider a slightly different Pyomo implementation of the model we just presented.

```
# abstract2.py

from __future__ import division
from pyomo.environ import *

model = AbstractModel()

model.I = Set()
```

```
model.d = Param(model.I, model.J)
model.b = Param(model.I)
model.c = Param(model.J)

# the next line declares a variable indexed by the set J
model.x = Var(model.J, domain=NonNegativeReals)

def obj_expression(model):
    return summation(model.c, model.x)

model.OBJ = Objective(rule=obj_expression)

def ax_constraint_rule(model, i):
    # return the expression for the constraint for i
    return sum(model.a[i,j] * model.x[j] for j in model.J) >= model.b[i]

# the next line creates one constraint for each member of the set model.I
model.AxbConstraint = Constraint(model.I, rule=ax_constraint_rule)
```

To get the same instantiated model, the following data file can be used.

```
# abstract2a.dat AMPL format

set I := 1;
set J := 1 2;

param a :=
1 1 3
1 2 4;

param c:=
1 2
2 3;

param b := 1 1;
```

However, this model can also be fed different data for problems of the same general form using meaningful indexes.

```
# abstract2.dat AMPL data format
set I := TV Film;
set J := Graham John Carol;

param a :=
TV Graham 3
TV John 4.4
TV Carol 4.9
Film Graham 1
Film John 2.4
Film Carol 1.1;

param c := [*]
Graham 2.2
John 3.1416
Carol 3;
```

```
param b := TV 1 Film 1 ;
```

## 3.5 A Simple Concrete Pyomo Model

It is possible to get nearly the same flexible behavior from models declared to be abstract and models declared to be concrete in Pyomo; however, we will focus on a straightforward concrete example here where the data is hard-wired into the model file. Python programmers will quickly realize that the data could have come from other sources.

We repeat the concrete model already given:

min 
$$2x_1 + 3x_2$$
  
s.t.  $3x_1 + 4x_2 \ge 1$   
 $x_1, x_2 \ge 0$ 

This is implemented as a concrete model as follows:

```
from __future__ import division
from pyomo.environ import *

model = ConcreteModel()

model.x = Var([1,2], domain=NonNegativeReals)

model.OBJ = Objective(expr = 2*model.x[1] + 3*model.x[2])

model.Constraint1 = Constraint(expr = 3*model.x[1] + 4*model.x[2] >= 1)
```

Although rule functions can also be used to specify constraints and objectives, in this example we use the expr option that is available only in concrete models. This option gives a direct specification of the expression.

## 3.6 Solving the Simple Examples

Pyomo supports modeling and scripting but does not install a solver automatically. In order to solve a model, there must be a solver installed on the computer to be used. If there is a solver, then the pyomo command can be used to solve a problem instance.

Suppose that the solver named glpk (also known as glpsol) is installed on the computer. Suppose further that an abstract model is in the file named abstract1.py and a data file for it is in the file named abstract1.dat. From the command prompt, with both files in the current directory, a solution can be obtained with the command:

```
pyomo solve abstract1.py abstract1.dat --solver=glpk
```

Since glpk is the default solver, there really is no need specify it so the --solver option can be dropped.

#### Note

There are two dashes before the command line option names such as solver.

To continue the example, if CPLEX is installed then it can be listed as the solver. The command to solve with CPLEX is

```
pyomo solve abstract1.py abstract1.dat --solver=cplex
```

This yields the following output on the screen:

```
[ 0.00] Setting up Pyomo environment
[ 0.00] Applying Pyomo preprocessing actions
[ 0.07] Creating model
[ 0.15] Applying solver
[ 0.37] Processing results
Number of solutions: 1
Solution Information
    Gap: 0.0
    Status: optimal
    Function Value: 0.66666666667
Solver results file: results.json
[ 0.39] Applying Pyomo postprocessing actions
[ 0.39] Pyomo Finished
```

The numbers is square brackets indicate how much time was required for each step. Results are written to the file named results.json, which has a special structure that makes it useful for post-processing. To see a summary of results written to the screen, use the --summary option:

```
pyomo solve abstract1.py abstract1.dat --solver=cplex --summary
```

To see a list of Pyomo command line options, use:

```
pyomo solve --help
```

#### Note

There are two dashes before help.

For a concrete model, no data file is specified on the Pyomo command line.

## Sets

### 4.1 Declaration

Sets can be declared using the Set and RangeSet functions or by assigning set expressions. The simplest set declaration creates a set and postpones creation of its members:

```
model.A = Set()
```

The Set function takes optional arguments such as:

- doc = String describing the set
- dimen = Dimension of the members of the set
- filter = A boolean function used during construction to indicate if a potential new member should be assigned to the set
- initialize = A function that returns the members to initialize the set. ordered = A boolean indicator that the set is ordered; the default is False
- validate = A boolean function that validates new member data
- virtual = A boolean indicator that the set will never have elements; it is unusual for a modeler to create a virtual set; they are typically used as domains for sets, parameters and variables
- within = Set used for validation; it is a super-set of the set being declared.

One way to create a set whose members will be two dimensional is to use the dimen argument:

```
model.B = Set(dimen=2)
```

To create a set of all the numbers in set model. A doubled, one could use

```
def doubleA_init(model):
    return (i*2 for i in model.A)
model.C = Set(initialize=DoubleA_init)
```

As an aside we note that as always in Python, there are lot of ways to accomplish the same thing. Also, note that this will generate an error if model. A contains elements for which multiplication times two is not defined.

The initialize option can refer to a Python set, which can be returned by a function or given directly as in

```
model.D = Set(initialize=['red', 'green', 'blue'])
```

The initialize option can also specify a function that is applied sequentially to generate set members. Consider the case of a simple set. In this case, the initialization function accepts a set element number and model and returns the set element associated with that number:

```
def Z_init(model, i):
    if i > 10:
        return Set.End
    return 2*i+1
model.Z = Set(initialize=Z_init)
```

The Set. End return value terminates input to the set. Additional information about iterators for set initialization is in the Pyomo book.

#### Note

Data specified in an input file will override the data specified by the initialize options.

If sets are given as arguments to Set without keywords, they are interpreted as indexes for an array of sets. For example, to create an array of sets that is indexed by the members of the set model. A, use

```
model.E = Set(model.A)
```

Arguments can be combined. For example, to create an array of sets with three dimensional members indexed by set model. A, use

```
model.F = Set(model.A, dimen=3)
```

The initialize option can be used to create a set that contains a sequence of numbers, but the RangeSet function provides a concise mechanism for simple sequences. This function takes as its arguments a start value, a final value, and a step size. If the RangeSet has only a single argument, then that value defines the final value in the sequence; the first value and step size default to one. If two values given, they are the first and last value in the sequence and the step size defaults to one. For example, the following declaration creates a set with the numbers 1.5, 5 and 8.5:

```
model.G = RangeSet(1.5, 10, 3.5)
```

## 4.2 Operations

Sets may also be created by assigning other Pyomo sets as in these examples that also illustrate the set operators union, intersection, difference, and exclusive-or:

```
model.H = model.A
model.I = model.A | model.D # union
model.J = model.A & model.D # intersection
model.K = model.A - model.D # difference
model.L = model.A ^ model.D # exclusive-or
```

The cross-product operator is the asterisk (\*). For example, to assign a set the cross product of two other sets, one could use

```
model.K = model.B * model.c
```

or to indicate the the members of a set are restricted to be in the cross product of two other sets, one could use

```
model.K = Set(within=model.B * model.C)
```

The cross-product operator is the asterisk (\*). For example, to create a set that contains the cross-product of sets A and B, use

```
model.C = Set(model.A * model.B)
```

to instead create a set that can contain a subset of the members of this cross-product, use

```
model.C = Set(within=model.A * model.B)
```

### 4.3 Predefined Virtual Sets

For use in specifying domains for sets, parameters and variables, Pyomo provides the following pre-defined virtual sets:

· Any: all possible values

• Reals: floating point values

• PositiveReals: strictly positive floating point values

• NonPositiveReals: non-positive floating point values

• NegativeReals: strictly negative floating point values

NonNegativeReals: non-negative floating point values

• PercentFraction: floating point values in the interval [0,1]

• Integers: integer values

• PositiveIntegers: positive integer values

• NonPositiveIntegers: non-positive integer values

• NegativeIntegers: negative integer values

• NonNegativeIntegers: non-negative integer values

• Boolean: boolean values, which can be represented as False/True, 0/1, 'False'/'True' and 'F'/'T'

· Binary: same as boolean

For example, if the set model. M is declared to be within the virtual set NegativeIntegers then an attempt to add anything other than a negative integer will result in an error. Here is the declaration:

```
model.M = Set(within=NegativeIntegers)
```

## 4.4 Sparse Index Sets

Sets provide indexes for parameters, variables and other sets. Index set issues are important for modelers in part because of efficiency considerations, but primarily because the right choice of index sets can result in very natural formulations that are condusive to understanding and maintenance. Pyomo leverages Python to provide a rich collection of options for index set creation and use.

The choice of how to represent indexes often depends on the application and the nature of the instance data that are expected. To illustrate some of the options and issues, we will consider problems involving networks. In many network applications, it is useful to declare a set of nodes, such as

```
model.Nodes = Set()
```

and then a set of arcs can be created with reference to the nodes.

Consider the following simple version of minimum cost flow problem:

$$\begin{array}{ll} \text{minimize} & \sum_{a \in \mathscr{A}} c_a x_a \\ \text{subject to:} & S_n + \sum_{(i,n) \in \mathscr{A}} x_{(i,n)} \\ & -D_n - \sum_{(n,j) \in \mathscr{A}} x_{(n,j)} & n \in \mathscr{N} \\ & x_a \geq 0, & a \in \mathscr{A} \end{array}$$

where

- Set: Nodes  $\equiv \mathcal{N}$
- Set: Arcs  $\equiv \mathscr{A} \subseteq \mathscr{N} \times \mathscr{N}$
- Var: Flow on arc (i, j):  $\equiv x_{i, j}, (i, j) \in \mathcal{A}$
- Param: Flow Cost on arc (i,j):  $\equiv c_{i,j}, \ (i,j) \in \mathscr{A}$
- Param: Demand at node  $i : \equiv D_i, i \in \mathcal{N}$
- Param: Supply at node  $i : \equiv S_i, i \in \mathcal{N}$

In the simplest case, the arcs can just be the cross product of the nodes, which is accomplished by the definition

```
model.Arcs = model.Nodes * model.Nodes
```

that creates a set with two dimensional members. For applications where all nodes are always connected to all other nodes this may suffice. However, issues can arise when the network is not fully dense. For example, the burden of avoiding flow on arcs that do not exist falls on the data file where high-enough costs must be provided for those arcs. Such a scheme is not very elegant or robust.

For many network flow applications, it might be better to declare the arcs using

model.Arcs = Set(within=model.Nodes\*model.Nodes)

```
or
```

```
model.Arcs = Set(dimen=2)
```

where the difference is that the first version will provide error checking as data is assigned to the set elements. This would enable specification of a sparse network in a natural way. But this results in a need to change the FlowBalance constraint because as it was written in the simple example, it sums over the entire set of nodes for each node. One way to remedy this is to sum only over the members of the set model.arcs as in

```
def FlowBalance_rule(model, node):
    return model.Supply[node] \
    + sum(model.Flow[i, node] for i in model.Nodes if (i,node) in model.Arcs) \
    - model.Demand[node] \
    - sum(model.Flow[node, j] for j in model.Nodes if (j,node) in model.Arcs) \
    == 0
```

This will be OK unless the number of nodes becomes very large for a sparse network, then the time to generate this constraint might become an issue (admittely, only for very large networks, but such networks do exist).

Another method, which comes in handy in many network applications, is to have a set for each node that contain the nodes at the other end of arcs going to the node at hand and another set giving the nodes on out-going arcs. If these sets are called model. NodesIn and model.NodesOut respectively, then the flow balance rule can be re-written as

```
def FlowBalance_rule(model, node):
    return model.Supply[node] \
    + sum(model.Flow[i, node] for i in model.NodesIn[node]) \
    - model.Demand[node] \
    - sum(model.Flow[node, j] for j in model.NodesOut[node]) \
    == 0
```

The data for NodesIn and NodesOut could be added to the input file, and this may be the most efficient option.

For all but the largest networks, rather than reading Arcs, NodesIn and NodesOut from a data file, it might be more elegant to read only Arcs from a data file and declare model.NodesIn with an initialize option specifying the creation as follows:

```
def NodesIn_init(model, node):
    retval = []
    for (i,j) in model.Arcs:
        if j == node:
            retval.append(i)
    return retval
model.NodesIn = Set(model.Nodes, initialize=NodesIn_init)
```

with a similar definition for model. NodesOut. This code creates a list of sets for NodesIn, one set of nodes for each node. The full model is:

```
# Isinglecomm.py
# NodesIn and NodesOut are intialized using the Arcs
from pyomo.environ import *
model = AbstractModel()
model.Nodes = Set()
model.Arcs = Set(dimen=2)
def NodesOut_init(model, node):
    retval = []
    for (i, j) in model.Arcs:
       if i == node:
           retval.append(j)
    return retval
model.NodesOut = Set(model.Nodes, initialize=NodesOut_init)
def NodesIn_init(model, node):
    retval = []
    for (i, j) in model.Arcs:
        if j == node:
            retval.append(i)
    return retval
model.NodesIn = Set(model.Nodes, initialize=NodesIn_init)
model.Flow = Var(model.Arcs, domain=NonNegativeReals)
model.FlowCost = Param(model.Arcs)
model.Demand = Param(model.Nodes)
model.Supply = Param(model.Nodes)
def Obj_rule(model):
    return summation (model.FlowCost, model.Flow)
model.Obj = Objective(rule=Obj_rule, sense=minimize)
def FlowBalance_rule(model, node):
    return model.Supply[node] \
     + sum(model.Flow[i, node] for i in model.NodesIn[node]) \
     - model.Demand[node] \
      - sum(model.Flow[node, j] for j in model.NodesOut[node]) \
model.FlowBalance = Constraint(model.Nodes, rule=FlowBalance_rule)
```

### for this model, a toy data file would be:

```
# Isinglecomm.dat: data for Isinglecomm.py
set Nodes := CityA CityB CityC;
set Arcs :=
```

```
CityA CityB
CityA CityC
CityC CityB
;

param : FlowCost :=
CityA CityB 1.4
CityA CityC 2.7
CityC CityB 1.6
;

param Demand :=
CityA 0
CityB 1
CityC 1
;

param Supply :=
CityA 2
CityB 0
CityC 0
;
```

This can be done somewhat more efficiently, and perhaps more clearly, using a build action as shown in Isinglebuild.py.

### 4.4.1 Sparse Index Sets Example

One may want to have a constraint that holds

```
for i in model.I, k in model.K, v in model.V[k]
```

There are many ways to accomplish this, but one good way is to create a set of tuples composed of all of model.k, model.V[k] pairs. This can be done as follows:

```
def kv_init(model):
    return ((k,v) for k in model.K for v in model.V[k])
model.KV=Set(dimen=2, initialize=kv_init)
```

So then if there was a constraint defining rule such as

```
def MyC_rule(model, i, k, v):
    return ...
```

Then a constraint could be declared using

```
model.MyConstraint = Constraint(model.I, model.KV, rule=c1Rule)
```

Here is the first few lines of a model that illustrates this:

```
from pyomo.environ import *

model = AbstractModel()

model.I=Set()
model.K=Set()
model.V=Set(model.K)

def kv_init(model):
    return ((k,v) for k in model.K for v in model.V[k])
model.KV=Set(dimen=2, initialize=kv_init)
```

```
model.a = Param(model.I, model.K)

model.y = Var(model.I)
model.x = Var(model.I, model.KV)

#include a constraint
#x[i,k,v] <= a[i,k]*y[i], for i in model.I, k in model.K, v in model.V[k]

def c1Rule(model,i,k,v):
    return model.x[i,k,v] <= model.a[i,k]*model.y[i]
model.c1 = Constraint(model.I,model.KV,rule=c1Rule)</pre>
```

## **Parameters**

The word "parameters" is used in many settings. When discussing a Pyomo model, we use the word to refer to data that must be provided in order to find an optimal (or good) assignment of values to the decision variables. Parameters are declared with the Param function, which takes arguments that are somewhat similar to the Set function. For example, the following code snippet declares sets model. A, model. B and then a parameter array model. P that is indexed by model. A:

```
model.A = Set()
model.B = Set()
model.P = Param(model.A, model.B)
```

In addition to sets that serve as indexes, the Param function takes the following command options:

- default = The value absent any other specification.
- doc = String describing the parameter
- initialize = A function (or Python object) that returns the members to initialize the parameter values.
- validate = A boolean function with arguments that are the prospective parameter value, the parameter indices and the model.
- within = Set used for validation; it specifies the domain of the parameter values.

These options perform in the same way as they do for Set. For example, suppose that Model.A =RangeSet (1, 3), then there are many ways to create a parameter that is a square matrix with 9, 16, 25 on the main diagonal zeros elsewhere, here are two ways to do it. First using a Python object to initialize:

```
 v=\{\} \\ v[1,1] = 9 \\ v[2,2] = 16 \\ v[3,3] = 25 \\ model.S = Param(model.A, model.A, initialize=v, default=0)
```

And now using an initialization function that is automatically called once for each index tuple (remember that we are assuming that model. A contains 1,2,3)

```
def s_init(model, i, j):
    if i == j:
        return i*i
    else:
        return 0.0
model.S = Param(model.A, model.A, initialize=s_init)
```

In this example, the index set contained integers, but index sets need not be numeric. It is very common to use strings.

### Note

Data specified in an input file will override the data specified by the initialize options.

Parameter values can be checked by a validation function. In the following example, the parameter S indexed by model.A and checked to be greater than 3.14159. If value is provided that is less than that, the model instantation would be terminated and an error message issued. The function used to validate should be written so as to return True if the data is valid and False otherwise.

```
def s_validate(model, v, i):
    return v > 3.14159
model.S = Param(model.A, validate=s_validate)
```

## **Variables**

Variables are intended to ultimately be given values by an optimization package. They are declared and optionally bounded, given initial values, and documented using the Pyomo Var function. If index sets are given as arguments to this function they are used to index the variable, other optional directives include:

- bounds = A function (or Python object) that gives a (lower,upper) bound pair for the variable
- domain = A set that is a super-set of the values the variable can take on.
- initialize = A function (or Python object) that gives a starting value for the variable; this is particularly important for non-linear models
- within = (synonym for domain)

The following code snippet illustrates some aspects of these options by declaring a *singleton* (i.e. unindexed) variable named model.LumberJack that will take on real values between zero and 6 and it initialized to be 1.5:

```
model.LumberJack = Var(within=NonNegativeReals, bounds=(0,6), initialize=1.5)
```

Instead of the initialize option, initialization is sometimes done with a Python assignment statement as in

```
model.LumberJack = 1.5
```

For indexed variables, bounds and initial values are often specified by a rule (a Python function) that itself may make reference to parameters or other data. The formal arguments to these rules begins with the model followed by the indexes. This is illustrated in the following code snippet that makes use of Python dictionaries declared as lb and ub that are used by a function to provide bounds:

```
model.A = Set(initialize=['Scones', 'Tea']
lb = {'Scones':2, 'Tea':4}
ub = {'Scones':5, 'Tea':7}
def fb(model, i):
    return (lb[i], ub[i])
model.PriceToCharge = Var(model.A, domain=PositiveInteger, bounds=fb)
```

#### Note

Many of the pre-defined virtual sets that are used as domains imply bounds. A strong example is the set Boolean that implies bounds of zero and one.

# **Objectives**

An objective is a function of variables that returns a value that an optimization package attempts to maximize or minimize. The Objective function in Pyomo declares an objective. Although other mechanisms are possible, this function is typically passed the name of another function that gives the expression. Here is a very simple version of such a function that assumes model.x has previously been declared as a Var:

```
def ObjRule(model):
    return 2*model.x[1] + 3*model.x[2]
model.g = Objective(rule=ObjRule)
```

It is more common for an objective function to refer to parameters as in this example that assumes that model.p has been declared as a parameters and that model.x has been declared with the same index set, while model.y has been declared as a singleton:

```
def profrul(model):
    return summation(model.p, model.x) + model.y
model.Obj = Objective(rule=ObjRule, sense=maximize)
```

This example uses the sense option to specify maximization. The default sense is minimize.

## **Constraints**

Most constraints are specified using equality or inequality expressions that are created using a rule, which is a Python function. For example, if the variable model.x has the indexes *butter* and *scones*, then this constraint limits the sum for them to be exactly three:

```
def teaOKrule(model):
    return(model.x['butter'] + model.x['scones'] == 3)
model.TeaConst = Constraint, rule=teaOKrule)
```

Instead of expressions involving equality (==) or inequalities (<= or >=), constraints can also be expressed using a 3-tuple if the form (lb, expr, ub) where lb and ub can be None, which is interpreted as lb <= expr <= ub. Variables can appear only in the middle expr. For example, the following two constraint declarations have the same meaning:

```
model.x = Var()

def aRule(model):
    return model.x >= 2
Boundx = Constraint(rule=aRule)

def bRule(model):
    return (2, model.x, None)
Boundx = Constraint(rule=bRule)
```

For this simple example, it would also be possible to declare model.x with a bound option to accomplish the same thing.

Constraints (and objectives) can be indexed by lists or sets. When the declaration contains lists or sets as arguments, the elements are iteratively passed to the rule function. If there is more than one, then the cross product is sent. For example the following constraint could be interpreted as placing a budget of i on the ith item to buy where the cost per item is given by the parameter model.a:

```
model.A = RangeSet(1,10)
model.a = Param(model.A, within=PostiveReals)
model.ToBuy = Var(model.A)
def bud_rule(model, i):
    return model.a[i] *model.ToBuy[i] <= i
aBudget = Constraint(model.A)</pre>
```

#### Note

Python and Pyomo are case sensitive so model.a is not the same as model.A.

# **Expressions**

In this chapter, we use the word "expression" is two ways: first in the general sense of the word and second to desribe a class of Pyomo objects that have the name expression as described in the subsection on expression objects.

## 9.1 Rules to Generate Expressions

Both objectives and constraints make use of rules to generate expressions. These are Python functions that return the appropriate expression. These are first-class functions that can access global data as well as data passed in, including the model object.

Operations on model elements results in expressions, which seems natural in expression like the constraints we have seen so far. It is also possible to build up expressions. The following example illustrates this along with a reference to global Pyton data in the form of a Python variable called switch:

```
switch = 3

model.A = RangeSet(1, 10)
model.c = Param(model.A)
model.d = Param()
model.x = Var(model.A, domain=Boolean)

def pi_rule(model)
    accexpr = summation(model.c, model.x)
    if switch >= 2:
        accexpr = accexpr - model.d
    return accexpr >= 0.5
PieSlice = Constraint(rule=pi_rule)
```

In this example, the constraint that is generated depends on the value of the Python variable called switch. If the value is 2 or greater, then the constraint is summation (model.c, model.x) -model.d >=0.5; otherwise, the model.d term is not present.



### Caution

Because model elements result in expressions, not values, the following does not work as expected in an abstract model!

```
model.A = RangeSet(1, 10)
model.c = Param(model.A)
model.d = Param()
model.x = Var(model.A, domain=Boolean)
```

```
def pi_rule(model)
    accexpr = summation(model.c, model.x)
    if model.d >= 2: # NOT in an abstract model!!
        accexpr = accexpr - model.d
    return accexpr >= 0.5
PieSlice = Constraint(rule=pi_rule)
```

The trouble is that model.d >=2 results in an expression, not its evaluated value. Instead use if value (model.d) >=2

## 9.2 Piecewise Linear Expressions

Pyomo has facilities to add piecewise constraints of the form y=f(x) for a variety of forms of the function f.

The piecewise types other than SOS2, BIGM\_SOS1, BIGM\_BIN are implement as described in the paper [Vielma\_et\_al].

There are two basic forms for the declaration of the constraint:

```
model.pwconst = Piecewise(indexes, yvar, xvar, **Keywords)
model.pwconst = Piecewise(yvar, xvar, **Keywords)
```

where pwconst can be replaced by a name appropriate for the application. The choice depends on whether the x and y variables are indexed. If so, they must have the same index sets and these sets are give as the first arguments.

#### KEYWORDS:

- pw\_pts={},[],() A dictionary of lists (keys are index set) or a single list (for the non-indexed case or when an identical set of breakpoints is used across all indices) defining the set of domain breakpoints for the piecewise linear function. NOTE: pw\_pts is always required. These give the breakpoints for the piecewise function and are expected to full span the bounds for the independent variable(s).
- pw\_repn=<Option> Indicates the type of piecewise representation to use. This can have a major impact on solver performance. Options: (Default 'SOS2')
  - 'SOS2' Standard representation using sos2 constraints.
  - 'BIGM\_BIN' BigM constraints with binary variables. The theoretically tightest M values are automatically determined.
  - 'BIGM\_SOS1' BigM constraints with sos1 variables. The theoretically tightest M values are automatically determined.
  - 'DCC' Disaggregated convex combination model.
  - 'DLOG' Logarithmic disaggregated convex combination model.
  - 'CC' Convex combination model.
  - 'LOG' Logarithmic branching convex combination.
  - 'MC' Multiple choice model.
  - 'INC' Incremental (delta) method. NOTE: Step functions are supported for all but the two BIGM options. Refer to the force\_pw option.
- pw\_constr\_type= <Option> Indicates the bound type of the piecewise function. Options:
  - 'UB' y variable is bounded above by piecewise function
  - 'LB' y variable is bounded below by piecewise function
  - 'EQ' y variable is equal to the piecewise function
- f\_rule=f(model,i,j,...,x), {}, [], ()

An object that returns a numeric value that is the range value corresponding to each piecewise domain point. For functions, the first argument must be a Pyomo model. The last argument is the domain value at which the function evaluates (Not a Pyomo Var). Intermediate arguments are the corresponding indices of the Piecewise component (if any). Otherwise, the object can be a dictionary of lists/tuples (with keys the same as the indexing set) or a singe list/tuple (when no indexing set is used or when all indices use an identical piecewise function). Examples:

```
# A function that changes with index
def f(model,j,x):
    if (j == 2):
        return x**2 + 1.0
    else:
        return x**2 + 5.0

# A nonlinear function
f = lambda model,x: return exp(x) + value(model.p)
        (model.p is a Pyomo Param)

# A step function
f = [0,0,1,1,2,2]
```

#### force\_pw=True/False

Using the given function rule and pw\_pts, a check for convexity/concavity is implemented. If (1) the function is convex and the piecewise constraints are lower bounds or if (2) the function is concave and the piecewise constraints are upper bounds then the piecewise constraints will be substituted for linear constraints. Setting *force\_pw=True* will force the use of the original piecewise constraints even when one of these two cases applies.

• warning\_tol=<float>

To aid in debugging, a warning is printed when consecutive slopes of piecewise segments are within <warning\_tol> of each other. Default=1e-8

• warn\_domain\_coverage=True/False

Print a warning when the feasible region of the domain variable is not completely covered by the piecewise breakpoints. Default=True

• unbounded\_domain\_var=True/False

Allow an unbounded or partially bounded Pyomo Var to be used as the domain variable. Default=False NOTE: This does not imply unbounded piecewise segments will be constructed. The outermost piecwise breakpoints will bound the domain variable at each index. However, the Var attributes .lb and .ub will not be modified.

Here is an example of an assignment to a Python dictionary variable that has keywords for a picewise constraint:

```
kwds = {'pw_constr_type':'EQ','pw_repn':'SOS2','sense':maximize,'force_pw':True}
```

Here is a simple example based on the abstract2.py example given early. In this new example, the objective function is the sum of c times x to the fourth. In this example, the keywords are passed directly to the Piecewise function without being assigned to a dictionary variable. The upper bound on the x variables was chosen whimsically just to make the example. The important thing to note is that variables that are going to appear as the independent variable in a piecewise constraint must have bounds.

```
# abstract2piece.py
# Similar to abstract2.py, but the objective is now c times x to the fourth power

from pyomo.environ import *

model = AbstractModel()

model.I = Set()
model.J = Set()

Topx = 6.1 # range of x variables

model.a = Param(model.I, model.J)
model.b = Param(model.I)
model.c = Param(model.J)

# the next line declares a variable indexed by the set J
```

```
model.x = Var(model.J, domain=NonNegativeReals, bounds=(0, Topx))
model.y = Var(model.J, domain=NonNegativeReals)
# to avoid warnings, we set breakpoints at or beyond the bounds
PieceCnt = 100
bpts = []
for i in range(PieceCnt+2):
   bpts.append(float((i*Topx)/PieceCnt))
def f4(model, j, xp):
    # we not need j, but it is passed as the index for the constraint
    return xp**4
model.ComputeObj = Piecewise(model.J, model.y, model.x, pw_pts=bpts, pw_constr_type='EQ',
   f_rule=f4)
def obj_expression(model):
    return summation (model.c, model.y)
model.OBJ = Objective(rule=obj_expression)
def ax_constraint_rule(model, i):
    # return the expression for the constraint for i
    return sum(model.a[i,j] * model.x[j] for j in model.J) >= model.b[i]
# the next line creates one constraint for each member of the set model.I
model.AxbConstraint = Constraint(model.I, rule=ax_constraint_rule)
```

A more advanced example is provided as abstract2piecebuild.py.

## 9.3 Expression Objects

Pyomo Expression objects are very similar to the Param component (with mutable=True) except that the underlying values can be numeric constants or Pyomo expressions. Here's an illustration of expression objects in an AbstractModel. An expression object with an index set that is the numbers 1, 2, 3 is created and initialized to be the model variable x times the index. Later in the model file, just to illustrate how to do it, the expression is changed but just for the first index to be x squared.

```
model = AbstractModel()
model.x = Var(initialize=1.0)
def _e(m,i):
   return m.x*i
model.e = Expression([1,2,3],initialize=_e)
instance = model.create_instance()
print value(instance.e[1]) # -> 1.0
print instance.e[1]()
                               # -> 1.0
print instance.e[1].value # -> a pyomo expression object
# Change the underlying expression
instance.e[1].value = instance.x**2
... solve
... load results
# print the value of the expression given the loaded optimal solution
print value(instance.e[1])
```

An alternative is to create Python functions that, potentially, manipulate model objects. E.g., if you define a function

```
def f(x, p):
    return x + p
```

You can call this function with or without Pyomo modeling components as the arguments. E.g., f(2,3) will return a number, whereas f(model.x, 3) will return a Pyomo expression due to operator overloading.

If you take this approach you should note that anywhere a Pyomo expression is used to generate another expression (e.g., f(model.x, 3) + 5), the initial expression is always cloned so that the new generated expression is independent of the old. For example:

```
model = ConcreteModel()
model.x = Var()

# create a Pyomo expression
e1 = model.x + 5

# create another Pyomo expression
# e1 is copied when generating e2
e2 = e1 + model.x
```

If you want to create an expression that is shared between other expressions, you can use the Expression component.

## **Chapter 10**

# **Data Input**

Pyomo can initialize models in two general ways. When executing the pyomo command, one or more data command files can be specified to declare data and load data from other data sources (e.g. spreadsheets and CSV files). When initializing a model within a Python script, a DataPortal object can be used to load data from one or more data sources.

### 10.1 Data Command Files

The following commands can be used in data command files:

- set declares set data,
- param declares a table of parameter data, which can also include the declaration of the set data used to index parameter data,
- load loads set and parameter data from an external data source such as ASCII table files, CSV files, ranges in spreadsheets, and database tables,
- table loads set and parameter data from a table,
- include specifies a data command file that is to be processed immediately,
- the data and end commands do not perform any actions, but they provide compatibility with AMPL scripts that define data commands, and
- namespace defines groupings of data commands.

The syntax of the set and param data commands are adapted from AMPL's data commands. However, other Pyomo data commands do not directly correspond to AMPL data commands. In particular, Pyomo's table command was introduced to work around semantic ambiguities in the param command. Pyomo's table command does not correspond to AMPL's table command. Instead, the load command mimics AMPL's table command with a simplified syntax.



#### Warning

The data command file was initially developed to provide compatability in data formats between Pyomo and AMPL. However, these data formats continue to diverge in their syntax and semantics. Simple examples using set and param data commands are likely to work for both AMPL and Pyomo, particularly with abstract Pyomo models. But in general a user should expect to need to adapt their AMPL data command files for use with Pyomo.

See the Pyomo book for detailed descriptions of these commands. The following sections provide additional details, particularly for new data commands that are not described in the Pyomo book: table.

#### 10.1.1 table

The table data command was developed to provide a more flexible and complete data declaration than is possible with the param declaration. This command has a similar syntax to the load command, but it includes a complete specification of the table data.

The following example illustrates a simple table command that declares data for a single parameter:

```
table M(A):
A B M N:=
A1 B1 4.3 5.3
A2 B2 4.4 5.4
A3 B3 4.5 5.5;
```

The parameter M is indexed by column A. The column labels are provided after the colon and before the :=. Subsequently, the table data is provided. Note that the syntax is not sensitive to whitespace. Thus, the following is an equivalent table command:

```
table M(A) :
A B M N :=
A1 B1 4.3 5.3 A2 B2 4.4 5.4 A3 B3 4.5 5.5 ;
```

Multiple parameters can be declared by simply including additional parameter names. For example:

```
table M(A) N(A,B) :
A B M N :=
A1 B1 4.3 5.3
A2 B2 4.4 5.4
A3 B3 4.5 5.5;
```

This example declares data for the M and N parameters. As this example illustrates, these parameters may have different indexing

The indexing columns represent set data, which is specified separately. For example:

```
table A={A} Z={A,B} M(A) N(A,B) :
A B M N :=
A1 B1 4.3 5.3
A2 B2 4.4 5.4
A3 B3 4.5 5.5;
```

This examples declares data for the M and N parameters, along with the A and Z indexing sets. The correspondence between the index set Z and the indices of parameter N can be made more explicit by indexing N by Z:

```
table A={A} Z={A,B} M(A) N(Z) :
A B M N :=
A1 B1 4.3 5.3
A2 B2 4.4 5.4
A3 B3 4.5 5.5
;
```

Set data can also be specified independent of parameter data:

```
table Z={A,B} Y={M,N} :
A B M N :=
A1 B1 4.3 5.3
A2 B2 4.4 5.4
A3 B3 4.5 5.5
;
```

Finally, singleton parameter values can be specified with a simple table command:

```
table pi := 3.1416 ;
```

The previous examples considered examples of the table command where column labels are provided. The table command can also be used without column labels. For example, the file table 0.dat can be revised to omit column labels as follows:

```
table columns=4 M(1)={3} :=
A1 B1 4.3 5.3
A2 B2 4.4 5.4
A3 B3 4.5 5.5
;
```

The columns=4 is a keyword-value pair that defines the number of columns in this table; this must be explicitly specified in unlabeled tables. The default column labels are integers starting from 1; the labels are columns 1, 2, 3, and 4 in this example. The M parameter is indexed by column 1. The braces syntax declares the column where the M data is provided.

Similarly, set data can be declared referencing the integer column labels:

```
table A={1} Z={1,2} M(1) N(1,2) :=
A1 B1 4.3 5.3
A2 B2 4.4 5.4
A3 B3 4.5 5.5
;
```

Declared set names can also be used to index parameters:

```
table A={1} Z={1,2} M(A) N(Z) :=
A1 B1 4.3 5.3
A2 B2 4.4 5.4
A3 B3 4.5 5.5;
```

Finally, we compare and contrast the table and param commands:

- Both commands can be used to declare parameter and set data.
- The param command can declare a single set that is used to index one or more parameters. The table command can declare data for any number of sets, independent of whether they are used to index parameter data.
- The param command can declare data for multiple parameters only if they share the same index set. The table command can declare data for any number of parameters that are may be indexed separately.
- Both commands can be used to declare a singleton parameter.
- The table syntax unambiguously describes the dimensionality of indexing sets. The param command must be interpreted with a model that provides the dimension of the indexing set.

This last point provides a key motivation for the table command. Specifically, the table command can be used to reliably initialize concrete models using a DataPortal object. By contrast, the param command can only be used to initialize concrete models with parameters that are indexed by a single column (i.e. a simple set). See the discussion of DataPortal objects below for an example.

#### 10.1.2 namespace

The namespace command allows data commands to be organized into named groups that can be enabled from the pyomo command line. For example, consider again the abstract2.py example. Suppose that the cost data shown in abstract2.dat were valid only under certain circumstances that we will label as "TerryG" and that there would be different cost data under circumstances that we will label "JohnD." This could be represented using the following data file:

```
# abs2nspace.dat AMPL format with namespaces
set I := TV Film ;
set J := Graham John Carol ;
param a :=
TV Graham 3
TV John 4.4
TV Carol 4.9
Film Graham 1
Film John 2.4
Film Carol 1.1
namespace TerryG {
  param c := [*]
     Graham 2.2
     John 3.1416
     Carol 3
   ;
}
namespace JohnD {
  param c := [*]
     Graham 2.7
     John 3
     Carol 2.1
param b := TV 1 Film 1 ;
```

To use this data file with abstract2.py, a namespace must be indicated on the command line. To select the "TerryG" data specification, --namespace TerryG would be added to the command line. For example:

```
pyomo solve abstract2.py abs2nspace.dat --namespace TerryG --solver=cplex
```

If the --namespace option is omitted, then no data will be given for model.c (and no default was given for model.c). In other words, there is no default namespace selection.

The option -ns (with one dash) is an alias for --namespace (which needs two dashes) Multiple namespaces can be selected by giving multiple --namespace or -ns arguments on the Pyomo command line.

## 10.2 DataPortal Objects

The load and store Pyomo data commands can be used to load set and table data from a variety of data sources. Pyomo's DataPortal object provides this same functionality for users who work with Python scripts. A DataPortal object manages the process of loading data from different data sources, and it is used to construct model instances in a standard manner. Similarly, a DataPortal object can be used to store model data externally in a standard manner.

#### 10.2.1 Loading Data

The load method can be used to load data into Pyomo models from a variety of sources and formats. The most common format is a table representation of set and parameter data. For example, consider the file A.tab, which defines a simple set:

```
A A1 A2
```

АЗ

The following example illustrates how a DataPortal object can be used to load this data into a model:

```
model = AbstractModel()
model.A = Set()

data = DataPortal()
data.load(filename='tab/A.tab', set=model.A)
instance = model.create(data)
```

The load method opens the data file, processes it, and loads the data in a format that is then used to construct a model instance. The load method can be called multiple times to load data for different sets or parameters, or to override data processed earlier.

#### Note

1.1

Subsequent examples omit the model declaration and instance creation.

In the previous example, the set option is used to define the model component that is loaded with the set data. If the data source defines a table of data, then this option is used to specify data for a multi-dimensional set. For example, consider the file D.tab:

```
A B
A1 1
A1 2
A1 3
A2 1
A2 2
A3 3
A3 1
A3 2
A3 3
```

If a two-dimensional set is declared, then it can be loaded with the same syntax:

```
model.A = Set(dimen=2)

data.load(filename='tab/C.tab', set=model.A)
```

This example also illustrates that the column titles do not directly impact the process of loading data. Column titles are only used to select columns that are included in the table that is loaded (see below).

The param option is used to define the a parameter component that is loaded with data. The simplest parameter is a singleton. For example, consider the file Z.tab:

```
This data is loaded with the following syntax:
```

```
model.z = Param()
data.load(filename='tab/Z.tab', param=model.z)
```

Indexed parameters can be defined from table data. For example, consider the file Y.tab:

```
A Y
A1 3.3
A2 3.4
A3 3.5
```

The parameter y is loaded with the following syntax:

```
model.A = Set(initialize=['A1','A2','A3','A4'])
model.y = Param(model.A)

data.load(filename='tab/Y.tab', param=model.y)
```

Pyomo assumes that the parameter values are defined on the rightmost column; the column names are not used to specify the index and parameter data (see below). In this file, the A column contains the index values, and the Y column contains the parameter values.

Similarly, multiple parameters can be initialized at once by specifying a list (or tuple) of component parameters. For example, consider the file XW.tab:

```
A X W
A1 3.3 4.3
A2 3.4 4.4
A3 3.5 4.5
```

The parameters x and w are loaded with the following syntax:

```
model.A = Set(initialize=['A1','A2','A3','A4'])
model.x = Param(model.A)
model.w = Param(model.A)

data.load(filename='tab/XW.tab', param=(model.x,model.w))
```

Note that the data for set A is predefined in this example. The index set can be loaded with the parameter data using the index option:

```
model.A = Set()
model.x = Param(model.A)
model.w = Param(model.A)

data.load(filename='tab/XW.tab', param=(model.x, model.w), index=model.A)
```

We have previously noted that the column names are not used to define the set and parameter data. The select option is used to define the columns in the table that are used to load data. This option specifies a list (or tuple) of column names that are used, in that order, to form the table that defines the component data.

For example, consider the following load declaration:

```
model.A = Set()
model.w = Param(model.A)

data.load(filename='tab/XW.tab', select=('A','W'), param=model.w, index=model.A)
```

The columns A and W are selected from the file XW.tab, and a single parameter is defined.

#### Note

The load method allows for a variety of other options that are supported by the add method for ModelData objects. See the Pyomo book for a detailed description of these options.

## **Chapter 11**

## BuildAction and BuildCheck

This is a somewhat advanced topic. In some cases, it is desirable to trigger actions to be done as part of the model building process. The BuildAction function provides this capability in a Pyomo model. It takes as arguments optional index sets and a function to perform the action. For example,

```
model.BuildBpts = BuildAction(model.J, rule=bpts_build)
```

calls the function <code>bpts\_build</code> for each member of <code>model.J</code>. The function <code>bpts\_build</code> should have the model and a variable for the members of <code>model.J</code> as formal arguments. In this example, the following would be a valid declaration for the function:

```
def bpts_build(model, j):
```

A full example, which extends the abstract2.py and abstract2piece.py examples, is

```
# abstract2piecebuild.py
# Similar to abstract2piece.py, but the breakpoints are created using a build action
from pyomo.environ import *
model = AbstractModel()
model.I = Set()
model.J = Set()
model.a = Param(model.I, model.J)
model.b = Param(model.I)
model.c = Param(model.J)
model.Topx = Param(default=6.1) # range of x variables
model.PieceCnt = Param(default=100)
# the next line declares a variable indexed by the set J
model.x = Var(model.J, domain=NonNegativeReals, bounds=(0, model.Topx))
model.y = Var(model.J, domain=NonNegativeReals)
# to avoid warnings, we set breakpoints beyond the bounds
# we are using a dictionary so that we can have different
# breakpoints for each index. But we won't.
model.bpts = {}
def bpts_build(model, j):
   model.bpts[j] = []
    for i in range(model.PieceCnt+2):
        model.bpts[j].append(float((i*model.Topx)/model.PieceCnt))
# The object model.BuildBpts is not referred to again;
```

```
# the only goal is to trigger the action at build time
model.BuildBpts = BuildAction(model.J, rule=bpts_build)

def f4(model, j, xp):
    # we not need j in this example, but it is passed as the index for the constraint
    return xp**4

model.ComputePieces = Piecewise(model.J, model.y, model.x, pw_pts=model.bpts, \leftarrow
    pw_constr_type='EQ', f_rule=f4)

def obj_expression(model):
    return summation(model.c, model.y)

model.OBJ = Objective(rule=obj_expression)

def ax_constraint_rule(model, i):
    # return the expression for the constraint for i
    return sum(model.a[i,j] * model.x[j] for j in model.J) >= model.b[i]

# the next line creates one constraint for each member of the set model.I
model.AxbConstraint = Constraint(model.I, rule=ax_constraint_rule)
```

This example uses the build action to create a model component with breakpoints for a piecewise function. The BuildAction is triggered by the assignment to model.BuildBpts. This object is not referenced again, the only goal is to cause the execution of bpts\_build, which places data in the model.bpts dictionary. Note that if model.bpts had been a Set, then it could have been created with an initialize argument to the Set declaration. Since it is a special-purpose dictionary to support the piecewise functionality in Pyomo, we use a BuildAction.

Another application of BuildAction can be intialization of Pyomo model data from Python data structures, or efficient initialization of Pyomo model data from other Pyomo model data. Consider the Isinglecomm.py example. Rather than using an initialization for each list of sets NodesIn and NodesOut separately using initialize, it is a little more efficient and probably a little clearer, to use a build action.

#### The full model is:

```
# Isinglecomm.py
# NodesIn and NodesOut are created by a build action using the Arcs
from pyomo.environ import *
model = AbstractModel()
model.Nodes = Set()
model.Arcs = Set(dimen=2)
model.NodesOut = Set(model.Nodes, within=model.Nodes, initialize=[])
model.NodesIn = Set(model.Nodes, within=model.Nodes, initialize=[])
def Populate_In_and_Out(model):
    # loop over the arcs and put the end points in the appropriate places
    for (i, j) in model.Arcs:
        model.NodesIn[j].add(i)
       model.NodesOut[i].add(j)
model.In_n_Out = BuildAction(rule = Populate_In_and_Out)
model.Flow = Var(model.Arcs, domain=NonNegativeReals)
model.FlowCost = Param(model.Arcs)
model.Demand = Param(model.Nodes)
model.Supply = Param(model.Nodes)
def Obj_rule(model):
```

```
return summation(model.FlowCost, model.Flow)
model.Obj = Objective(rule=Obj_rule, sense=minimize)

def FlowBalance_rule(model, node):
    return model.Supply[node] \
    + sum(model.Flow[i, node] for i in model.NodesIn[node]) \
    - model.Demand[node] \
    - sum(model.Flow[node, j] for j in model.NodesOut[node]) \
    == 0

model.FlowBalance = Constraint(model.Nodes, rule=FlowBalance_rule)
```

for this model, the same data file can be used as for Isinglecomm.py such as the toy data file:

```
# Isinglecomm.dat: data for Isinglecomm.py
set Nodes := CityA CityB CityC ;
set Arcs :=
CityA CityB
CityA CityC
CityC CityB
param : FlowCost :=
CityA CityB 1.4
CityA CityC 2.7
CityC CityB 1.6
param Demand :=
CityA 0
CityB 1
CityC 1
param Supply :=
CityA 2
CityB 0
CityC 0
```

Build actions can also be a way to implement data validation, particularly when multiple Sets or Parameters must be analyzed. However, the the BuildCheck component is prefered for this purpose. It executes its rule just like a BuildAction but will terminate the construction of the model instance if the rule returns False.

## **Chapter 12**

# The pyomo Command

The pyomo command is issued to the DOS prompt or a Unix shell. To see a list of Pyomo command line options, use:

```
pyomo solve --help
```

#### Note

There are two dashes before help.

In this section we will detail some of the options.

## 12.1 Passing Options to a Solver

To pass arguments to a solver when using the pyomo solve command, appned the Pyomo command line with the argument – solver-options= followed by an argument that is a string to be sent to the solver (perhaps with dashes added by Pyomo). So for most MIP solvers, the mip gap can be set using

```
--solver-options= "mipgap=0.01 "
```

Multiple options are separated by a space. Options that do not take an argument should be specified with the equals sign followed by either a space or the end of the string.

For example, to specify that the solver is GLPK, then to specify a mipgap of two percent and the GLPK cuts option, use

```
--solver=glpk --solver-options="mipgap=0.02 cuts="
```

If there are multiple "levels" to the keyword, as is the case for some Gurobi and CPLEX options, the tokens are separated by underscore. For example, mip cuts all would be specified as mip\_cuts\_all. For another example, to set the solver to be CPLEX, then to set a mip gap of one percent and to specify y for the sub-option numerical to the option emphasis use

```
--solver=cplex --solver-options="mipgap=0.001 emphasis_numerical=y"
```

See Solver Options for a discusion of passing options in a script.

### 12.2 Troubleshooting

Many of things that can go wrong are covered by error messages, but sometimes they can be confusing or do not provide enough information. Depending on what the troubles are, there might be ways to get a little additional information.

If there are syntax errors in the model file, for example, it can occasionally be helpful to get error messages directly from the Python interpreter rather than through Pyomo. Suppose the name of the model file is scuc.py, then

```
python scuc.py
```

can sometimes give useful information for fixing syntax errors.

When there are no syntax errors, but there troubles reading the data or generating the information to pass to a solver, then the -verbose option provides a trace of the execution of Pyomo. The user should be aware that for some models this option can generate a lot of output.

If there are troubles with solver (i.e., after Pyomo has output "Applying Solver"), it is often helpful to use the option <code>--stream-solver</code> that causes the solver output to be displayed rather than trapped. (See Solver Display for information about getting this output in a script). Advanced users may wish to examine the files that are generated to be passed to a solver. The type of file generated is controlled by the <code>--solver-io</code> option and the <code>--keepfiles</code> option instructs pyomo to keep the files and output their names. However, the <code>--symbolic-solver-labels</code> option should usually also be specified so that meaningful names are used in these files.

When there seem to be troubles expressing the model, it is often useful to embed print commands in the model in places that will yield helpful information. Consider the following snippet:

```
def ax_constraint_rule(model, i):
    # return the expression for the constraint for i
    print "ax_constraint_rule was called for i=",i
    return sum(model.a[i,j] * model.x[j] for j in model.J) >= model.b[i]

# the next line creates one constraint for each member of the set model.I
model.AxbConstraint = Constraint(model.I, rule=ax_constraint_rule)
```

The effect will be to output every member of the set model. I at the time the constraint named model. AxbConstraint is constructed.

#### 12.3 Direct Interfaces to Solvers

In many applications, the default solver interface works well. However, in some cases it is useful to specify the interface using the solver-io option. For example, if the solver supports a direct Python interface, then the option would be specified on the command line as

```
--solver-io=python
```

Here are some of the choices:

- lp: generate a standard linear programming format file with filename extension lp
- nlp: generate a file with a standard format that supports linear and nonlinear optimization with filename extension nllp
- os: generate an OSiL format XML file.
- python: use the direct Python interface.

Note that not all solvers support all interfaces.

## **Chapter 13**

# **PySP Overview**

This chapter describes PySP: (Pyomo Stochastic Programming), where parameters are allowed to be uncertain.

### 13.1 Overview of Modeling Components and Processes

The sequence of activities is typically the following:

- Create a deterministic model and declare components
- Develop base-case data for the deterministic model
- Test, verify and validate the deterministic model
- Model the stochastic processes
- Develop a way to generate scenarios (in the form of a tree if there are more than two stages)
- Create the data files need to describe the stochastics
- Use PySP to solve stochastic problem

When viewed from the standpoint of file creation, the process is

- Create an abstract model for the deterministic problem in a file called ReferenceModel.py
- Specify the stochastics in a file called ScenarioStructure.dat
- · Specify scenario data

## 13.2 Birge and Louveaux's Farmer Problem

Birge and Louveaux [BirgeLouveauxBook] make use of the example of a farmer who has 500 acres that can be planted in wheat, corn or sugar beets, at a per acre cost of 150, 230 and 260 (Euros, presumably), respectively. The farmer needs to have at least 200 tons of wheat and 240 tons of corn to use as feed, but if enough is not grown, those crops can be purchased for 238 and 210, respectively. Corn and wheat grown in excess of the feed requirements can be sold for 170 and 150, respectively. A price of 36 per ton is guaranteed for the first 6000 tons grown by any farmer, but beets in excess of that are sold for 10 per ton. The yield is 2.5, 3, and 20 tons per acre for wheat, corn and sugar beets, respectively.

#### 13.2.1 ReferenceModel.py

So far, this is a deterministic problem because we are assuming that we know all the data. The Pyomo model for this problem shown here is in the file ReferenceModel.py in the sub-directory examples/pysp/farmer/models that is distributed with Pyomo.

```
# Farmer: rent out version has a scalar root node var
# note: this will minimize
#
# Imports
from __future__ import division
from pyomo.environ import *
# Model
model = AbstractModel()
# Parameters
model.CROPS = Set()
model.TOTAL_ACREAGE = Param(within=PositiveReals)
model.PriceQuota = Param(model.CROPS, within=PositiveReals)
model.SubQuotaSellingPrice = Param(model.CROPS, within=PositiveReals)
def super_quota_selling_price_validate (model, value, i):
    return model.SubQuotaSellingPrice[i] >= model.SuperQuotaSellingPrice[i]
model.SuperQuotaSellingPrice = Param(model.CROPS, validate= \leftrightarrow
   super_quota_selling_price_validate)
model.CattleFeedRequirement = Param(model.CROPS, within=NonNegativeReals)
model.PurchasePrice = Param(model.CROPS, within=PositiveReals)
model.PlantingCostPerAcre = Param(model.CROPS, within=PositiveReals)
model.Yield = Param(model.CROPS, within=NonNegativeReals)
# Variables
model.DevotedAcreage = Var(model.CROPS, bounds=(0.0, model.TOTAL_ACREAGE))
model.QuantitySubQuotaSold = Var(model.CROPS, bounds=(0.0, None))
model.QuantitySuperQuotaSold = Var(model.CROPS, bounds=(0.0, None))
model.QuantityPurchased = Var(model.CROPS, bounds=(0.0, None))
model.FirstStageCost = Var()
model.SecondStageCost = Var()
```

```
# Constraints
def ConstrainTotalAcreage_rule(model):
        return summation(model.DevotedAcreage) <= model.TOTAL_ACREAGE
model.ConstrainTotalAcreage = Constraint(rule=ConstrainTotalAcreage_rule)
def EnforceCattleFeedRequirement_rule(model, i):
        return model.CattleFeedRequirement[i] <= (model.Yield[i] * model.DevotedAcreage[i]) + ←
                 \verb|model.QuantityPurchased[i]| - \verb|model.QuantitySubQuotaSold[i]| - \verb|model.| \leftarrow
                 QuantitySuperQuotaSold[i]
model.EnforceCattleFeedRequirement = Constraint(model.CROPS, rule= \leftrightarrow
        EnforceCattleFeedRequirement_rule)
def LimitAmountSold_rule(model, i):
         return\ model.QuantitySubQuotaSold[i]\ +\ model.QuantitySuperQuotaSold[i]\ -\ (model.Yield[i\ \leftrightarrow\ Colored by the colored by t
                 ] * model.DevotedAcreage[i]) <= 0.0</pre>
model.LimitAmountSold = Constraint(model.CROPS, rule=LimitAmountSold_rule)
def EnforceQuotas_rule(model, i):
        return (0.0, model.QuantitySubQuotaSold[i], model.PriceQuota[i])
model.EnforceQuotas = Constraint(model.CROPS, rule=EnforceQuotas_rule)
#
   Stage-specific cost computations
def ComputeFirstStageCost_rule(model):
        return model.FirstStageCost - summation(model.PlantingCostPerAcre, model.DevotedAcreage ↔
                 ) == 0.0
model.ComputeFirstStageCost = Constraint(rule=ComputeFirstStageCost_rule)
def ComputeSecondStageCost_rule(model):
        expr = summation(model.PurchasePrice, model.QuantityPurchased)
        expr -= summation(model.SubQuotaSellingPrice, model.QuantitySubQuotaSold)
        expr -= summation(model.SuperQuotaSellingPrice, model.QuantitySuperQuotaSold)
        return (model.SecondStageCost - expr) == 0.0
model.ComputeSecondStageCost = Constraint(rule=ComputeSecondStageCost_rule)
   Objective
def Total_Cost_Objective_rule(model):
        return model.FirstStageCost + model.SecondStageCost
model.Total_Cost_Objective = Objective(sense=minimize, rule=Total_Cost_Objective_rule)
```

#### 13.2.2 Example Data

The data introduced here are in the file ReferenceModel.dat in the sub-directory examples/pysp/farmer/scenariodata that is distributed with Pyomo. These data are given for illustration. The file ReferenceModel.dat is not required by PySP.

```
set CROPS := WHEAT CORN SUGAR_BEETS ;
```

```
param TOTAL_ACREAGE := 500 ;
# no quotas on wheat or corn
param PriceQuota :=
           WHEAT 100000 CORN 100000 SUGAR_BEETS 6000;
param SubQuotaSellingPrice :=
           WHEAT 170 CORN 150 SUGAR_BEETS 36;
param SuperQuotaSellingPrice :=
           WHEAT 0 CORN 0 SUGAR_BEETS 10;
param CattleFeedRequirement :=
            WHEAT 200 CORN 240 SUGAR_BEETS 0;
# can't purchase beets (no need, as cattle don't eat them)
param PurchasePrice :=
           WHEAT 238 CORN 210 SUGAR_BEETS 100000;
param PlantingCostPerAcre :=
           WHEAT 150 CORN 230 SUGAR_BEETS 260;
param Yield := WHEAT 3.0 CORN 3.6 SUGAR_BEETS 24;
```

Any of these data could be modeled as uncertain, but we will consider only the possibility that the yield per acre could be higher or lower than expected. Assume that there is a probability of 1/3 that the yields will be the average values that were given (i.e., wheat 2.5; corn 3; and beets 20). Assume that there is a 1/3 probability that they will be lower (2, 2.4, 16) and 1/3 probability they will be higher (3, 3.6, 24). We refer to each full set of data as a *scenario* and collectively we call them a *scenario tree*. In this case the scenario tree is very simple: there is a root node and three leaf nodes: one corresponding to each scenario. The acreage-to-plant decisions are root node decisions because they must be made without knowing what the yield will be. The other variables are so-called *second stage* decisions, because they will depend on which scenario is realized.

#### 13.2.3 ScenarioStructure.dat

PySP requires that users describe the scenario tree using specific constructs in a file named <code>ScenarioStructure.dat</code>; for the farmer problem, this file can be found in the pyomo sub-directory <code>examples/pysp/farmer/scenariodata</code> that is distributed with Pyomo.

```
# IMPORTANT - THE STAGES ARE ASSUMED TO BE IN TIME-ORDER.
set Stages := FirstStage SecondStage ;
set Nodes := RootNode
            BelowAverageNode
            AverageNode
            AboveAverageNode ;
param NodeStage := RootNode
                                  FirstStage
                  BelowAverageNode SecondStage
                  AverageNode SecondStage
                  AboveAverageNode SecondStage ;
set Children[RootNode] := BelowAverageNode
                         AverageNode
                         AboveAverageNode ;
param ConditionalProbability := RootNode
                                                 1.0
                               BelowAverageNode 0.33333333
                               AverageNode 0.33333334
```

This data file is verbose and somewhat redundant, but in most applications it is generated by software rather than by a person, so this is not an issue. Generally, the left-most part of each expression (e.g. "set Stages :=") is required and uses reserved words (e.g., Stages) and the other names are supplied by the user (e.g., "FirstStage" could be any name). Every assignment is terminated with a semi-colon. We will now consider the assignments in this file one at a time.

The first assignments provides names for the stages and the words "set Stages" are required, as are the := symbols. Any names can be used. In this example, we used "FirstStage" and "SecondStage" but we could have used "EtapPrimero" and "ZweiteEtage" if we had wanted to. Whatever names are given here will continue to be used to refer to the stages in the rest of the file. The order of the names is important. A simple way to think of it is that generally, the names must be in time order (technically, they need to be in order of information discovery, but that is usually time-order). Stages refers to decision stages, which may, or may not, correspond directly with time stages. In the farmer example, decisions about how much to plant are made in the first stage and "decisions" (which are pretty obvious, but which are decision variables nonetheless) about how much to sell at each price and how much needs to be bought are second stage decisions because they are made after the yield is known.

```
set Stages := FirstStage SecondStage ;
```

Node names are constructed next. The words "set Nodes" are required, but any names may be assigned to the nodes. In two stage stochastic problems there is a root node, which we chose to name "RootNode" and then there is a node for each scenario.

Nodes are associated with time stages with an assignment beginning with the required words "param Nodestage." The assignments must make use of previously defined node and stage names. Every node must be assigned a stage.

The structure of the scenario tree is defined using assignment of children to each node that has them. Since this is a two stage problem, only the root node has children. The words "param Children" are required for every node that has children and the name of the node is in square brackets before the colon-equals assignment symbols. A list of children is assigned.

The probability for each node, conditional on observing the parent node is given in an assignment that begins with the required words "param ConditionalProbability." The root node always has a conditional probability of 1, but it must always be given anyway. In this example, the second stage nodes are equally likely.

Scenario names are given in an assignment that begins with the required words "set Scenarios" and provides a list of the names of the scenarios. Any names may be given. In many applications they are given unimaginative names generated by software such as "Scen1" and the like. In this example, there are three scenarios and the names reflect the relative values of the yields.

Leaf nodes, which are nodes with no children, are associated with scenarios. This assignment must be one-to-one and it is initiated with the words "param ScenarioLeafNode" followed by the colon-equals assignment characters.

Variables are associated with stages using an assignment that begins with the required words "set Stage Variables" and the name of a stage in square brackets followed by the colon-equals assignment characters. Variable names that have been defined in the file ReferenceModel.py can be assigned to stages. Any variables that are not assigned are assumed to be in the last stage. Variable indexes can be given explicitly and/or wildcards can be used. Note that the variable names appear without the prefix "model." In the farmer example, DevotedAcreage is the only first stage variable.

### Note

Variable names appear without the prefix "model."

#### Note

Wildcards can be used, but fully general Python slicing is not supported.

For reporting purposes, it is useful to define auxiliary variables in <code>ReferenceModel.py</code> that will be assigned the cost associated with each stage. This variables do not impact algorithms, but the values are output by some software during execution as well as upon completion. The names of the variables are assigned to stages using the "param StageCostVariable" assignment. The stages are previously defined in <code>ScenarioStructure.dat</code> and the variables are previously defined in <code>ReferenceModel.py</code>.

#### 13.2.4 Scenario data specification

So far, we have given a model in the file named <code>ReferenceModel.py</code>, a set of deterministic data in the file named <code>ReferenceModel.py</code>, and a description of the stochastics in the file named <code>ScenarioStructure.dat</code>. All that remains is to give the data for each scenario. There are two ways to do that in <code>PySP</code>: <code>scenario-based</code> and <code>node-based</code>. The default is scenario-based so we will describe that first.

For scenario-based data, the full data for each scenario is given in a .dat file with the root name that is the name of the scenario. So, for example, the file named AverageScenario.dat must contain all the data for the model for the scenario named "AvererageScenario." It turns out that this file can be created by simply copying the file ReferenceModel.dat as shown above because it contains a full set of data for the "AverageScenario" scenario. The files BelowAverageScenario.dat and AboveAverageScenario.dat will differ from this file and from each other only in their last line, where the yield is specified. These three files are distributed with Pyomo and are in the pyomo sub-directory examples/pysp/farmer/scenariodata along with ScenarioStructure.dat and ReferenceModel.dat.

Scenario-based data wastes resources by specifying the same thing over and over again. In many cases, that does not matter and it is convenient to have full scenario data files available (for one thing, the scenarios can easily be run independently using the pyomo command). However, in many other settings, it is better to use a node-based specification where the data that is unique to each node is specified in a .dat file with a root name that matches the node name. In the farmer example, the file RootNode.dat will be the same as ReferenceModel.dat except that it will lack the last line that specifies the yield. The files BelowAverageNode.dat, AverageNode.dat, and AboveAverageNode.dat will contain only one line each to specify the yield. If node-based data is to be used, then the ScenarioStructure.dat file must contain the following line:

```
param ScenarioBasedData := False ;
```

An entire set of files for node-based data for the farmer problem are distributed with Pyomo in the sub-directory examples/pysp/farmer/nodedata

### 13.3 Finding Solutions for Stochastic Models

PySP provides a variety of tools for finding solutions to stochastic programs.

#### 13.3.1 runef

The runef command puts together the so-called *extensive form* version of the model. It creates a large model that has constraints to ensure that variables at a node have the same value. For example, in the farmer problem, all of the DevotedAcres variables must have the same value regardless of which scenario is ultimately realized. The objective can be the expected value of the objective function, or the CVaR, or a weighted combination of the two. Expected value is the default. A full set of options for runef can be obtained using the command:

```
runef --help
```

The pyomo distribution contains the files need to run the farmer example in the sub-directories to the sub-directory examples/pysp/farmer so if this is the current directory and if CPLEX is installed, the following command will cause formation of the EF and its solution using CPLEX.

```
runef -m models -i nodedata --solver=cplex --solve
```

The option -m models has one dash and is short-hand for the option --model-directory=models and note that the full option uses two dashes. The -i is equivalent to --instance-directory= in the same fashion. The default solver is CPLEX, so the solver option is not really needed. With the --solve option, runef would simply write an .lp data file that could be passed to a solver.

#### 13.3.2 runph

The runph command executes an implementation of Progressive Hedging (PH) that is intended to support scripting and extension.

The pyomo distribution contains the files need to run the farmer example in the sub-directories to the sub-directory examples/pysp/farmer so if this is the current directory and if CPLEX is installed, the following command will cause PH to execute using the default sub-problem solver, which is CPLEX.

```
runph -m models -i nodedata
```

The option -m models has one dash and is short-hand for the option --model-directory=models and note that the full option uses two dashes. The -i is equivalent to --instance-directory= in the same fashion.

After about 33 iterations, the algorithm will achieve the default level of convergence and terminate. A lot of output is generated and among the output is the following solution information:

```
Variable=DevotedAcreage
       Index: [CORN]
                                 (Scenarios:
                                             BelowAverageScenario
                                                                    AverageScenario
           AboveAverageScenario
                                 )
                                          80.0000
                           79.9844
                                                       79.9768
                                                                                 0.0232 ←
               Values:
                                                                   Max-Min=
                                79.9871
                       Avg=
       Index: [SUGAR_BEETS]
                                         (Scenarios: BelowAverageScenario
           AverageScenario
                             AboveAverageScenario
                                                    )
               Values:
                            249.9848
                                         249.9770
                                                      250.0000
                                                                   Max-Min=
                                                                                 0.0230
                       Avg=
                               249.9873
        Index: [WHEAT]
                                (Scenarios: BelowAverageScenario
                                                                    AverageScenario
           AboveAverageScenario
                                  )
               Values:
                          170.0308
                                         170.0230
                                                      170.0232
                                                                   Max-Min=
                                                                                 0.0078 ↔
                       Avg=
                             170.0256
Cost Variable=FirstStageCost
       Tree Node=RootNode
                                         (Scenarios: BelowAverageScenario
           AverageScenario AboveAverageScenario
                                                   )
       Values: 108897.0836 108897.4725 108898.1476
                                                           Max-Min=
                                                                         1.0640
                                                                                    Ava=
           108897.5679
```

For problems with no, or few, integer variables, the default level of convergence leaves root-node variables almost converged. Since the acreage to be planted cannot depend on the scenario that will be realized in the future, the average, which is labeled "Avg" in this output, would be used. A farmer would probably interpret acreages of 79.9871, 249.9873, and 170.0256 to be 80, 250, and 170. In real-world applications, PH is embedded in scripts that produce output in a format desired by a decision maker.

But in real-world applications, the default settings for PH seldom work well enough. In addition to post-processing the output, a number of parameters need to be adjusted and sometimes scripting to extend or augment the algorithm is needed to improve convergence rates. A full set of options can be obtained with the command:

```
runph --help
```

Note that there are two dashes before help.

By default, PH uses quadratic objective functions after iteration zero; in some settings it may be desirable to linearize the quadratic terms. This is required to use a solver such as glpk for MIPs because it does not support quadratic MIPs. The directive ——linearize—nonbinary—penalty—terms=n causes linearization of the penalty terms using n pieces. For example, to use glpk on the farmer, assuming glpk is installed and the command is given when the current directory is the examples/pysp/farmer, the following command will use default settings for most parameters and four pieces to approximate quadratic terms in sub-problems:

```
runph -i nodedata -m models --solver=glpk --linearize-nonbinary-penalty-terms=4
```

Use of the linearize-nonbinary-penalty-terms option requires that all variables not in the final stage have bounds.

#### **13.3.3 Solution Output Control**

To get the full solution, including leaf node solution values, use the runph --output-scenario-tree-solution option. In both runph and runef the solution can be written in csv format using the --solution-writer=pyomo.pysp. plugins.csvsolutionwriter option.

## 13.4 Summary of PySP File Names

PySP scripts such as runef and runph require files that specify the model and data using files with specific names. All files can be in the current directory, but typically, the file ReferenceModel.py is in a directory that is specified using --model-

directory= option (the short version of this option is -i +) and the data files are in a directory specified in the +--instance-directory= option (the short version of this option is +-m +).

#### Note

A file name other than ReferenceModel.py can be used if the file name is given in addition to the directory name as an argument to the --instance-directory option. For example, on a Windows machine --instance-directory=models\MyModel.py would specify the file MyModel.py in the local directory models.

- ReferenceModel.py: A full Pyomo model for a singe scenario. There should be no scenario indexes in this model because they are implicit.
- ScenarioStructure.dat: Specifies the nature of the stochastics. It also specifies whether the rest of the data is node-based or scenario-based. It is scenario-based unless ScenarioStructure.dat contains the line

```
param ScenarioBasedData := False ;
```

If scenario-based, then there is a data file for each scenario that specifies a full set of data for the scenario. The name of the file is the name of the scenario with .dat appended. The names of the scenarios are given in the ScenarioStructure.dat file.

If node-based, then there is a file with data for each node that specifies only that data that is unique for the node. The name of the file is the name of the node with .dat appended. The names of the nodes are given in the ScenarioStructure.dat file.

### 13.5 Solving Sub-problems in Parallel and/or Remotely

The Python package called Pyro provides capalities that are used to enable PH to make use of multiple solver processes for sub-problems and allows both runef and runph to make use remote solvers. We will focus on PH in our discussion here.

There are two solver management systems available for runph, one is based on a pyro\_mip\_server and the other is based on a phsolverserver. Regardless of which is used, a name server and a dispatch server must be running and accessible to the runph process. The name server is launched using the command pyomo\_ns and then the dispatch server is launched with dispatch\_srvr. Note that both commands contain an underscore. Both programs keep running until terminated by an external signal, so it is common to pipe their output to a file.

Solvers are controlled by solver servers. The pyro mip solver server is launched with the command pyro\_mip\_server. This command may be repeated to launch as many solvers as are desired. The runph then needs a --solver-manager=pyro option to signal that runph should not launch its own solver, but should send subproblems to be dispatched to parallel solvers. To summarize the commands:

```
• Once: pyomo_ns
```

• Once: dispatch\_srvr

• Multiple times: pyro\_mip\_server

• Once: runph ... --solver-manager=pyro ...

#### Note

The runph option —-shutdown-pryo will cause a shutdown signal to be sent to pyomo\_ns, dispatch\_srvr and all pyro\_mip\_server programs upon termination of runph.

Instead of using pyro\_mip\_server, one can use phsolverserver in its place. You can get a list of arguments using pyrosolverserver —help, which does not launch a solver server (it just displays help and terminates). If you use the phsolversover, then use —solver—manager=phpyro as an argument to runph rather than —solver—manager=pyro.



#### Warning

Unlike the normal pyro\_mip\_server, there must be one phsolverserver for each sub-problem. One can use fewer phsolverservers than there are scenarios by adding the command-line option "--phpyro-required-workers=X". This will partition the jobs among the available workers,

## **Chapter 14**

## **Suffixes**

Suffixes provide a mechanism for declaring extraneous model data, which can be used in a number of contexts. Most commonly, suffixes are used by solver plugins to store extra information about the solution of a model. This and other suffix functionality is made available to the modeler through the use of the Suffix component class. Uses of Suffix include:

- Importing extra information from a solver about the solution of a mathematical program (e.g., constraint duals, variable reduced costs, basis information).
- Exporting information to a solver or algorithm to aid in solving a mathematical program (e.g., warm-starting information, variable branching priorities).
- · Tagging modeling components with local data for later use in advanced scripting algorithms.

## 14.1 Suffix Notation and the Pyomo NL File Interface

The Suffix component used in Pyomo has been adapted from the suffix notation used in the modeling language AMPL [AMPL]. Therefore, it follows naturally that AMPL style suffix functionality is fully available using Pyomo's NL file interface. For information on AMPL style suffixes the reader is referred to the AMPL website:

http://www.ampl.com

A number of scripting examples that highlight the use AMPL style suffix functionality are available in the examples/pyomo/suffixes directory distributed with Pyomo.

#### 14.2 Declaration

The effects of declaring a Suffix component on a Pyomo model are determined by the following traits:

- direction: This trait defines the direction of information flow for the suffix. A suffix direction can be assigned one of four possible values:
  - LOCAL suffix data stays local to the modeling framework and will not be imported or exported by a solver plugin (default)
  - IMPORT suffix data will be imported from the solver by its respective solver plugin
  - EXPORT suffix data will be exported to a solver by its respective solver plugin
  - IMPORT\_EXPORT suffix data flows in both directions between the model and the solver or algorithm
- datatype: This trait advertises the type of data held on the suffix for those interfaces where it matters (e.g., the NL file interface). A suffix datatype can be assigned one of three possible values:

- FLOAT the suffix stores floating point data (default)
- INT the suffix stores integer data
- None the suffix stores any type of data

#### Note

Exporting suffix data through Pyomo's NL file interface requires all active export suffixes have a strict datatype (i.e., datatype=None is not allowed).

The following code snippet shows examples of declaring a Suffix component on a Pyomo model:

```
from pyomo.environ import *

model = ConcreteModel()

# Export integer data
model.priority = Suffix(direction=Suffix.EXPORT, datatype=Suffix.INT)

# Export and import floating point data
model.dual = Suffix(direction=Suffix.IMPORT_EXPORT)

# Store floating point data
model.junk = Suffix()
```

Declaring a Suffix with a non-local direction on a model is not guaranteed to be compatible with all solver plugins in Pyomo. Whether a given Suffix is acceptable or not depends on both the solver and solver interface being used. In some cases, a solver plugin will raise an exception if it encounters a Suffix type that it does not handle, but this is not true in every situation. For instance, the NL file interface is generic to all AMPL-compatible solvers, so there is no way to validate that a Suffix of a given name, direction, and datatype is appropriate for a solver. One should be careful in verifying that Suffix declarations are being handled as expected when switching to a different solver or solver interface.

### 14.3 Operations

The Suffix component class provides a dictionary interface for mapping Pyomo modeling components to arbitrary data. This mapping functionality is captured within the ComponentMap base class, which is also available within Pyomo's modeling environment. The ComponentMap can be used as a more lightweight replacement for Suffix in cases where a simple mapping from Pyomo modeling components to arbitrary data values is required.

#### Note

ComponentMap and Suffix use the built-in id() function for hashing entry keys. This design decision arises from the fact that most of the modeling components found in Pyomo are either not hashable or use a hash based on a mutable numeric value, making them unacceptable for use as keys with the built-in dict class.



#### Warning

The use of the built-in id() function for hashing entry keys in ComponentMap and Suffix makes them inappropriate for use in situations where built-in object types must be used as keys. It is strongly recommended that only Pyomo modeling components be used as keys in these mapping containers (Var, Constraint, etc.).



#### Warning

Do not attempt to pickle or deepcopy instances of ComponentMap or Suffix unless doing so along with the components for which they hold mapping entries. As an example, placing one of these objects on a model and then cloning or pickling that model is an acceptable scenario.

In addition to the dictionary interface provided through the ComponentMap base class, the Suffix component class also provides a number of methods whose default semantics are more convenient for working with indexed modeling components. The easiest way to highlight this functionality is through the use of an example.

```
from pyomo.environ import *

model = ConcreteModel()
model.x = Var()
model.y = Var([1,2,3])
model.foo = Suffix()
```

In this example we have a concrete Pyomo model with two different types of variable components (indexed and non-indexed) as well as a Suffix declaration (foo). The next code snippet shows examples of adding entries to the suffix foo.

```
# Assign a suffix value of 1.0 to model.x
model.foo.setValue(model.x, 1.0)

# Same as above with dict interface
model.foo[model.x] = 1.0

# Assign a suffix value of 0.0 to all indices of model.y
# By default this expands so that entries are created for
# every index (y[1], y[2], y[3]) and not model.y itself
model.foo.setValue(model.y, 0.0)

# The same operation using the dict interface results in an entry only
# for the parent component model.y
model.foo[model.y] = 50.0

# Assign a suffix value of -1.0 to model.y[1]
model.foo.setValue(model.y[1], -1.0)
# Same as above with the dict interface
model.foo[model.y[1]] = -1.0
```

In this example we highlight the fact that the <code>setitem</code> and <code>setValue</code> entry methods can be used interchangeably except in the case where indexed components are used (model.y). In the indexed case, the <code>setitem</code> approach creates a single entry for the parent indexed component itself, whereas the <code>setValue</code> approach by default creates an entry for each index of the component. This behavior can be controlled using the optional keyword <code>expand</code>, where assigning it a value of <code>False</code> results in the same behavior as <code>setitem</code>.

Other operations like accessing or removing entries in our mapping can performed as if the built-in dict class is in use.

```
# -> 1.0
print (model.foo.get (model.x))
print (model.foo[model.x])
                                        # -> 1.0
print (model.foo.get (model.y[1]))
                                        \# -> -1.0
print (model.foo[model.y[1]])
                                        \# -> -1.0
                                        # -> 0.0
print (model.foo.get (model.y[2]))
print (model.foo[model.y[2]])
                                        # -> 0.0
                                        # -> 50.0
print (model.foo.get (model.y))
print (model.foo[model.y])
                                        # -> 50.0
del model.foo[model.y]
print (model.foo.get (model.y))
                                        # -> None
                                        # -> raise KeyError
print (model.foo[model.y])
```

The non-dict method clearValue can be used in place of delitem to remove entries, where it inherits the same default behavior as setValue for indexed components and does not raise a KeyError when the argument does not exist as a key in the mapping.

```
model.foo.clearValue(model.y)

print(model.foo[model.y[1]]) # -> raise KeyError

del model.foo[model.y[1]] # -> raise KeyError

model.foo.clearValue(model.y[1]) # -> does nothing
```

A summary non-dict Suffix methods is provided here:

```
clearAllValues()
    Clears all suffix data.
clearValue(component, expand=True)
    Clears suffix information for a component.
setAllValues(value)
    Sets the value of this suffix on all components.
setValue(component, value, expand=True)
    Sets the value of this suffix on the specified component.
updateValues(data_buffer, expand=True)
    Updates the suffix data given a list of component, value tuples. Provides
    an improvement in efficiency over calling setValue on every component.
getDatatype()
    Return the suffix datatype.
setDatatype (datatype)
    Set the suffix datatype.
getDirection()
    Return the suffix direction.
setDirection(direction)
    Set the suffix direction.
importEnabled()
   Returns True when this suffix is enabled for import from solutions.
exportEnabled()
    Returns True when this suffix is enabled for export to solvers.
```

## 14.4 Importing Suffix Data

Importing suffix information from a solver solution is achieved by declaring a Suffix component with the appropriate name and direction. Suffix names available for import may be specific to third-party solvers as well as individual solver interfaces within Pyomo. The most common of these, available with most solvers and solver interfaces, is constraint dual multipliers. Requesting that duals be imported into suffix data can be accomplished by declaring a Suffix component on the model.

```
from pyomo.environ import *
```

```
model = ConcreteModel()
model.dual = Suffix(direction=Suffix.IMPORT)
model.x = Var()
model.obj = Objective(expr=model.x)
model.con = Constraint(expr=model.x>=1.0)
```

The existence of an active suffix with the name dual that has an import style suffix direction will cause constraint dual information to be collected into the solver results (assuming the solver supplies dual information). In addition to this, after loading solver results into a problem instance (using a python script or Pyomo callback functions in conjunction with the pyomo command), one can access the dual values associated with constraints using the dual Suffix component.

```
print(instance.dual[instance.con]) # -> 1.0
```

Alternatively, the pyomo option —solver—suffixes can be used to request suffix information from a solver. In the event that suffix names are provided via this command-line option, the pyomo script will automatically declare these Suffix components on the constructed instance making these suffixes available for import.

### 14.5 Exporting Suffix Data

Exporting suffix data is accomplished in a similar manner as to that of importing suffix data. One simply needs to declare a Suffix component on the model with an export style suffix direction and associate modeling component values with it. The following example shows how one can declare a special ordered set of type 1 using AMPL-style suffix notation in conjunction with Pyomo's NL file interface.

```
from pyomo.environ import *

model = ConcreteModel()
model.y = Var([1,2,3],within=NonNegativeReals)

model.sosno = Suffix(direction=Suffix.EXPORT)
model.ref = Suffix(direction=Suffix.EXPORT)

# Add entry for each index of model.y
model.sosno.setValue(model.y,1)
model.ref[model.y[1]] = 0
model.ref[model.y[2]] = 1
model.ref[model.y[3]] = 2
```

Most AMPL-compatible solvers will recognize the suffix names sosno and ref as declaring a special ordered set, where a positive value for sosno indicates a special ordered set of type 1 and a negative value indicates a special ordered set of type 2.

#### Note

Pyomo provides the SOSConstraint component for declaring special ordered sets, which is recognized by all solver interface, including the NL file interface.

Pyomo's NL file interface will recognize an EXPORT style Suffix component with the name *dual* as supplying initializations for constraint multipliers. As such it will be treated separately than all other EXPORT style suffixes encountered in the NL writer, which are treated as AMPL-style suffixes. The following example script shows how one can warmstart the interior-point solver Ipopt by supplying both primal (variable values) and dual (suffixes) solution information. This dual suffix information can be both imported and exported using a single Suffix component with an IMPORT\_EXPORT direction.

```
from pyomo.environ import *
from pyomo.opt import SolverFactory

### Create the ipopt solver plugin using the ASL interface
```

```
solver = 'ipopt'
solver_io = 'nl'
                         # True prints solver output to screen
stream_solver = True
keepfiles = False # True prints intermediate file names (.nl,.sol,...)
opt = SolverFactory(solver, solver_io=solver_io)
if opt is None:
    print("")
    print("ERROR: Unable to create solver plugin for %s "\
          "using the %s interface" % (solver, solver_io))
    print("")
    exit(1)
### Create the example model
model = ConcreteModel()
model.x1 = Var(bounds=(1,5),initialize=1.0)
model.x2 = Var(bounds=(1,5),initialize=5.0)
model.x3 = Var(bounds=(1,5), initialize=5.0)
model.x4 = Var(bounds=(1,5), initialize=1.0)
\verb|model.obj| = Objective(expr=model.x1*model.x4*(model.x1+model.x2+model.x3) + model.x3)|
model.inequality = Constraint(expr=model.x1*model.x2*model.x3*model.x4 >= 25.0)
model.equality = Constraint(expr=model.x1**2 + model.x2**2 + model.x3**2 + model.x4**2 == \leftrightarrow
    40.0)
###
### Declare all suffixes
# Ipopt bound multipliers (obtained from solution)
model.ipopt_zL_out = Suffix(direction=Suffix.IMPORT)
model.ipopt_zU_out = Suffix(direction=Suffix.IMPORT)
# Ipopt bound multipliers (sent to solver)
model.ipopt_zL_in = Suffix(direction=Suffix.EXPORT)
model.ipopt_zU_in = Suffix(direction=Suffix.EXPORT)
# Obtain dual solutions from first solve and send to warm start
model.dual = Suffix(direction=Suffix.IMPORT_EXPORT)
### Send the model to ipopt and collect the solution
print("")
print("INITIAL SOLVE")
results = opt.solve(model,keepfiles=keepfiles,tee=stream_solver)
# load the results (including any values for previously declared
# IMPORT / IMPORT_EXPORT Suffix components)
model.solutions.load_from(results)
###
### Set Ipopt options for warm-start
# The current values on the ipopt_zU_out and
# ipopt_zL_out suffixes will be used as initial
# conditions for the bound multipliers to solve
# the new problem
model.ipopt_zL_in.update(model.ipopt_zL_out)
model.ipopt_zU_in.update(model.ipopt_zU_out)
opt.options['warm_start_init_point'] = 'yes'
opt.options('warm_start_bound_push') = 1e-6
opt.options['warm_start_mult_bound_push'] = 1e-6
opt.options['mu_init'] = 1e-6
### Send the model and suffix information to ipopt and collect the solution
print("")
print("WARM-STARTED SOLVE")
```

```
# The solver plugin will scan the model for all active suffixes
# valid for importing, which it will store into the results object
results = opt.solve(model,keepfiles=keepfiles,tee=stream_solver)
# load the results (including any values for previously declared
# IMPORT / IMPORT_EXPORT Suffix components)
model.solutions.load_from(results)
###
```

The difference in performance can be seen by examining Ipopt's iteration log with and without warm starting:

• Without Warmstart:

```
inf_pr
iter
       objective
                           inf_du lg(mu) ||d|| lg(rg) alpha_du alpha_pr ls
  0 1.6109693e+01 1.12e+01 5.28e-01 -1.0 0.00e+00 - 0.00e+00 0.00e+00
    1.6982239e+01 7.30e-01 1.02e+01 -1.0 6.11e-01
                                                       7.19e-02 1.00e+00f
     1.7318411e+01 3.60e-02 5.05e-01
                                    -1.0 1.61e-01
                                                       1.00e+00 1.00e+00h
                                    -1.7 2.85e-01
     1.6849424e+01 2.78e-01 6.68e-02
                                                       7.94e-01 1.00e+00h
     1.7051199e+01 4.71e-03 2.78e-03 -1.7 6.06e-02
                                                       1.00e+00 1.00e+00h
     1.7011979e+01 7.19e-03 8.50e-03 -3.8 3.66e-02
                                                    - 9.45e-01 9.98e-01h 1
   6 1.7014271e+01 1.74e-05 9.78e-06 -3.8 3.33e-03 - 1.00e+00 1.00e+00h 1
     1.7014021e+01 1.23e-07 1.82e-07 -5.7 2.69e-04 - 1.00e+00 1.00e+00h 1
  7
  8 1.7014017e+01 1.77e-11 2.52e-11 -8.6 3.32e-06 - 1.00e+00 1.00e+00h 1
Number of Iterations...: 8
```

• With Warmstart:

```
iter    objective    inf_pr    inf_du lg(mu) ||d|| lg(rg) alpha_du alpha_pr ls
    0 1.7014032e+01 2.00e-06 4.07e-06 -6.0 0.00e+00 - 0.00e+00 0.00e+00 0
    1 1.7014019e+01 3.65e-12 1.00e-11 -6.0 2.50e-01 - 1.00e+00 1.00e+00h 1
    2 1.7014017e+01 4.48e-12 6.43e-12 -9.0 1.92e-06 - 1.00e+00 1.00e+00h 1
Number of Iterations...: 2
```

## 14.6 Using Suffixes With an AbstractModel

In order to allow the declaration of suffix data within the framework of an AbstractModel, the Suffix component can be initialized with an optional construction rule. As with constraint rules, this function will be executed at the time of model construction. The following simple example highlights the use of the rule keyword in suffix initialization. Suffix rules are expected to return an iterable of (component, value) tuples, where the expand=True semantics are applied for indexed components.

```
from pyomo.environ import *

model = AbstractModel()
model.x = Var()
model.c = Constraint(expr= model.x >= 1)

def foo_rule(m):
    return ((m.x, 2.0), (m.c, 3.0))
model.foo = Suffix(rule=foo_rule)

# Instantiate the model
inst = model.create()
print(inst.foo[model.x]) # -> raise KeyError
print(inst.foo[inst.x]) # -> 2.0
print(inst.foo[inst.c]) # -> 3.0
```

The next example shows an abstract model where suffixes are attached only to the variables:

```
from pyomo.environ import *

model = AbstractModel()
model.I = RangeSet(1,4)
model.x = Var(model.I)
def c_rule(m, i):
    return m.x[i] >= i
model.c = Constraint(model.I, rule=c_rule)

def foo_rule(m):
    return ((m.x[i], 3.0*i) for i in m.I)
model.foo = Suffix(rule=foo_rule)

# instantiate the model
inst = model.create_instance()
for i in inst.I:
    print (i, inst.foo[inst.x[i]])
```

## **Chapter 15**

## **DAE Toolbox**

The DAE toolbox allows users to incorporate differential equations in a Pyomo model. The modeling components in this toolbox are able to represent ordinary or partial differential equations. The differential equations do not have to be written in a particular format and the components are flexible enough to represent higher-order derivatives or mixed partial derivatives. The toolbox also includes model transformations which use a simultaneous discretization approach for transforming a DAE model into an algebraic model.

## 15.1 DAE Modeling Components

The DAE toolbox introduces three new modeling components to Pyomo:

#### ContinuousSet

Used to represent bounded continuous domains

#### DerivativeVar

Defines how a Var will be differentiated or the derivatives to be included in the model

#### Integral

Defines an integral over a continous domain

As will be shown later, differential equations can be declared using using these new DAE modeling components along with the standard Pyomo Var and Constraint components.

#### 15.1.1 ContinuousSet

This component is used to define continuous bounded domains (for example *spatial* or *time* domains). It is similar to a Pyomo Set component and can be used to index things like variables and constraints. In the current implementation, models with ContinuousSet components may not be solved until every ContinuousSet has been discretized. Minimally, a ContinuousSet must be initialized with two numeric values representing the upper and lower bounds of the continuous domain. A user may also specify additional points in the domain to be used as finite element points in the discretization.

The following code snippet shows examples of declaring a ContinuousSet component on a concrete Pyomo model:

```
# Required imports
from pyomo.environ import *
from pyomo.dae import *

model = ConcreteModel()

# declare by providing bounds
model.t = ContinuousSet(bounds=(0,5))
```

```
# declare by initializing with desired discretization points
model.x = ContinuousSet(initialize=[0,1,2,5])
```

The following code snippet shows an example of declaring a ContinuousSet component on an abstract Pyomo model using the example data file.

```
set t := 0 0.5 2.25 3.75 5;
```

```
# Required imports
from pyomo.environ import *
from pyomo.dae import *

model = AbstractModel()

# The ContinuousSet below will be initialized using the points
# in the data file when a model instance is created.
model.t = ContinuousSet()
```

#### Note

A ContinuousSet may not be constructed unless two numeric bounding points are provided.

#### Note

If a separate data file is used to initialize a ContinuousSet, it is done using the set command and not continuousset

Most valid ways to declare and initialize a Set can be used to declare and initialize a ContinuousSet. See the documentation for Set for additional options.



#### Warning

Be careful using a ContinuousSet as an implicit index in an expression, i.e. sum(m.v[i] for i in m.myContinuousSet). The expression will be generated using the discretization points contained in the ContinuousSet at the time the expression was constructed and will not be updated if additional points are added to the set.

SUMMARY OF NON-DICT CONTINUOUS SET METHODS

#### get finite elements()

If the ContinuousSet has been discretizaed using a collocation scheme, this method will return a list of the finite element discretization points but not the collocation points over each finite element. Otherwise this method returns a list of all the discretization points in the ContinuousSet.

#### get\_discretization\_info()

Returns a dictionary containing information on the discretization scheme that has been applied to the ContinuousSet.

#### get\_changed()

Returns "True" if additional points were added to the ContinousSet while applying a discretization scheme

#### get\_upper\_element\_boundary(value)

Returns the first finite element point that is greater than or equal to the value sent to the function.

#### get\_lower\_element\_boundary(value)

Returns the first finite element point that is less than or equal to the value sent to the function.

#### 15.1.2 DerivativeVar

The DerivativeVar component is used to declare a derivative of a Var. A Var may only be differentiated with respect to a ContinuousSet that it is indexed by. The indexing sets of a DerivativeVar are identical to those of the Var it is differentiating.

The code snippet below shows examples of declaring DerivativeVar components on a Pyomo model. In each case, the variable being differentiated is supplied as the only positional argument and the type of derivative is specified using the *wrt* (or the more verbose *withrespectto*) keyword argument. Any keyword argument that is valid for a Pyomo Var component may also be specified.

```
# Required imports
from pyomo.environ import *
from pyomo.dae import *
model = ConcreteModel()
model.s = Set(initialize=['a','b'])
model.t = ContinuousSet(bounds=(0,5))
model.l = ContinuousSet(bounds=(-10,10))
model.x = Var(model.t)
model.y = Var(model.s, model.t)
model.z = Var(model.t, model.l)
# Declare the first derivative of model.x with respect to model.t
model.dxdt = DerivativeVar(model.x, withrespectto=model.t)
# Declare the second derivative of model.y with respect to model.t
# Note that this DerivativeVar will be indexed by both model.s and model.t
model.dydt2 = DerivativeVar(model.y, wrt=(model.t, model.t))
# Declare the partial derivative of model.z with respect to model.1
# Note that this DerivativeVar will be indexed by both model.t and model.1
model.dzdl = DerivativeVar(model.z, wrt=(model.l), initialize=0)
# Declare the mixed second order partial derivative of model.z with respect
# to model.t and model.l and set bounds
model.dz2 = DerivativeVar(model.z, wrt=(model.t, model.l), bounds=(-10,10))
```

#### Note

The *initialize* keyword argument will initialize the value of a derivative and is not the same as specifying an initial condition. Initial or boundary conditions should be specified using a Constraint or ConstraintList.

Another way to use derivatives without explicitly declaring <code>DerivativeVar</code> components is to use the .derivative() method on a variable within an expression or constraint. For example:

```
# Required imports
from pyomo.environ import *
from pyomo.dae import *

model = ConcreteModel()
model.t = ContinuousSet(bounds=(0,5))
model.x = Var(model.t)

# Create the first derivative of model.x with respect to model.t
# within a constraint rule.
def _diffeq_rule(m,i):
    return m.x[i].derivative(m.t) == m.x[i]**2
model.diffeq = Constraint(model.t,rule=_diffeq_rule)
```

In the above example a DerivatveVar component representing the desired derivative will automatically be added to the Pyomo model when the constraint is constructed. The .derivative() method accepts positional arguments representing what the derivative is being taken with respect to.

#### Note

If a variable is indexed by a single ContinuousSet then the .derivative() method with no positional arguments may be used to specify the first derivative of that variable with respect to the ContinuousSet.

### 15.2 Declaring Differential Equations

A differential equations is declared as a standard Pyomo Constraint and is not required to have any particular form. The following code snippet shows how one might declare an ordinary or partial differential equation.

```
# Required imports
from pyomo.environ import *
from pyomo.dae import *
model = ConcreteModel()
model.s = Set(initialize=['a','b'])
model.t = ContinuousSet(bounds=(0,5))
model.l = ContinuousSet(bounds=(-10,10))
model.x = Var(model.s, model.t)
model.y = Var(model.t, model.l)
model.dydt = DerivativeVar(model.y, wrt=model.t)
model.dydl2 = DerivativeVar(model.y, wrt=(model.1, model.1))
# An ordinary differential equation
def _ode_rule(m,i,j):
    if j == 0:
        return Constraint.Skip
    return m.x[i].derivative(m.t) == m.x[i] **2
model.ode = Constraint(model.s, model.t, rule=_ode_rule)
# A partial differential equation
def _pde_rule(m,i,j):
    if i == 0 or j == -10 or j == 10:
        return Constraint.Skip
    return m.dydt[i,j] == m.dydl2[i,j]
model.pde = Constraint(model.t, model.l, rule=_pde_rule)
```

#### Note

Often a modeler does not want to apply a differential equation at one or both boundaries of a continuous domain. This must be addressed explicitly in the Constraint declaration using *Constraint.Skip* as shown above. By default, a Constraint declared over a ContinuousSet will be applied at every discretization point contained in the set.

## 15.3 Declaring Integrals

The Integral component is still under development but some basic functionality is available in the current Pyomo release. Integrals must be taken over the entire domain of a ContinuousSet. Once every ContinuousSet in a model has been discretized, any integrals in the model will be converted to algebraic equations using the trapezoid rule. Future releases of this tool will include more sophisticated numerical integration methods.

Declaring an Integral component is similar to declaring an Expression component. A simple example is shown below:

```
def __intX(m,i):
        return (m.X[i]-m.X_desired)**2
model.intX = Integral(model.time,wrt=model.time,rule=_intX)

def __obj(m):
        return m.scale*m.intX
model.obj = Objective(rule=_obj)
```

Notice that the positional arguments supplied to the Integral declaration must include all indices needed to evaluate the integral expression. The integral expression is defined in a function and supplied to the *rule* keyword argument. Finally, a user must specify a ContinuousSet that the integral is being evaluated over. This is done using the *wrt* keyword argument.

#### Note

The ContinousSet specified using the wrt keyword argument must be explicitly specified as one of the indexing sets (meaning it must be supplied as a positional argument)

After an Integral has been declared, it can be used just like a Pyomo Expression component and can be included in constraints or the objective function as shown above.

If an Integral is specified with multiple positional arguments, i.e. multiple indexing sets, the final component will be indexed by all of those sets except for the ContinuousSet that the integral was taken over. In other words, the ContinuousSet specified with the wrt keyword argument is removed from the indexing sets of the Integral even though it must be specified as a positional argument. The reason for this is to keep track of the order of the indexing sets. This logic should become more clear with the following example showing a double integral over the ContinuousSet components t1 and t2. In addition, the expression is also indexed by the Set s.

```
def __intX1(m,i,j,s):
    return (m.X[i,j,s]-m.X_desired[j,s])**2
model.intX1 = Integral(model.t1,model.t2,model.s,wrt=model.t1,rule=_intX1)

def __intX2(m,j,s):
    return (m.intX1[j,s]-m.X_desired[s])**2
model.intX2 = Integral(model.t2,model.s,wrt=model.t2,rule=_intX2)

def __obj(m):
    return sum(model.intX2[k] for k in m.s)
model.obj = Objective(rule=_obj)
```

#### 15.4 Discretization Transformations

Before a Pyomo model with DerivativeVar or Integral components can be sent to a solver it must first be sent through a discretization transformation. These transformations approximate any derivatives or integrals in the model by using some numerical method. The numerical methods currently included in this tool discretize the continuous domains in the problem and introduce equality constraints which approximate the derivatives and integrals at the discretization points. Two families of discretization schemes have been implemented in Pyomo, Finite Difference and Collocation. These schemes are described in more detail below.

#### Note

The schemes described here are for derivatives only. All integrals will be transformed using the trapezoid rule.

The user must write a Python script in order to use these discretizations, they have not been tested on the pyomo command line. Example scripts are shown below for each of the discretization schemes. The transformations return Pyomo models which can be further manipulated before being sent to a solver. Examples of this are also shown below.

#### 15.4.1 Finite Difference Transformation

This transformation includes implementations of several finite difference methods. For example, the Backward Difference method (also called Implicit or Backward Euler) has been implemented. The discretization equations for this method are shown below:

```
Given dx/dt = f(t,x) and x(t0) = x_0
discretize t and t and
```

where *h* is the step size between discretization points or the size of each finite element. These equations are generated automatically as Constraint components when the backward difference method is applied to a Pyomo model.

There are several discretization options available to a Finite\_Difference\_Transformation which can be specified as keyword arguments to the .apply() function of the transformation object. These keywords are summarized below:

KEYWORD ARGUMENTS FOR APPLYING A FINITE DIFFERENCE TRANSFORMATION.

nfe

The desired number of finite element points to be included in the discretization. The default value is 10.

wrt

Indicates which ContinuousSet the transformation should be applied to. If this keyword argument is not specified then the same scheme will be applied to all ContinuousSets.

#### scheme

Indicates which finite difference method to apply. Options are *BACKWARD*, *CENTRAL*, or *FORWARD*. The default scheme is the backward difference method.

#### inplace

Indicates whether the transformation should be applied to a copy of the model or the model itself. The default value is

If the existing number of finite element points in a ContinuousSet is less than the desired number, new discretization points will be added to the set. If a user specifies a number of finite element points which is less than the number of points already included in the ContinuousSet then the transformation will ignore the specified number and proceed with the larger set of points. Discretization points will never be removed from a ContinuousSet during the discretization.

The following code is a Python script applying the backward difference method. The code also shows how to add a constraint to a discretized model.

```
from pyomo.environ import *
from pyomo.dae import *
from pyomo.opt import SolverFactory
from pyomo.dae.plugins.finitedifference import Finite_Difference_Transformation

# Import Pyomo model
from pyomoExample import model

# Create model
instance = model.create()

# Discretize model using Backward Difference method
discretize = Finite_Difference_Transformation()
disc_instance = discretize.apply(instance,nfe=20,wrt=instance.time,scheme='BACKWARD')

# Add another constraint to discretized model
def _sum_limit(model):
    return sum(model.x1[i] for i in model.time) <= 50
disc_instance.con_sum_limit = Constraint(rule=_sum_limit)</pre>
```

```
# Solve discretized model
opt = SolverFactory('ipopt')
results = opt.solve(disc_instance)
# Load Results
disc_instance.load(results)
```

#### 15.4.2 Collocation Transformation

This transformation uses orthogonal collocation to discretize the differential equations in the model. Currently, two types of collocation have been implemented. They both use Lagrange polynomials with either Gauss-Radau roots or Gauss-Legendre roots. For more information on orthogonal collocation and the discretization equations associated with this method please see chapter 10 of the book "Nonlinear Programming: Concepts, Algorithms, and Applications to Chemical Processes" by L.T. Biegler.

The discretization options available to a Collocation\_Discretization\_Transformation are the same as those described above for the +Finite\_Difference\_Transformation with different available schemes and the addition of the *ncp* option.

ADDITIONAL KEYWORD ARGUMENTS FOR COLLOCATION DISCRETIZATIONS

#### scheme

The desired collocation scheme, either *LAGRANGE-RADAU* or *LAGRANGE-LEGENDRE*. The default is *LAGRANGE-RADAU*.

#### ncp

The number of collocation points within each finite element. The default value is 3.

#### Note

If the user's version of Python has access to the package Numpy then any number of collocation points may be specified, otherwise the maximum number is 10.

#### Note

Any points that exist in a ContinuousSet before discretization will be used as finite element boundaries and not as collocation points. The locations of the collocation points cannot be specified by the user, they must be generated by the transformation.

The following code is a Python script applying collocation with Lagrange polynomials and Radau roots. The code also shows how to add an objective function to a discretized model.

```
from pyomo.environ import *
from pyomo.dae import *
from pyomo.opt import SolverFactory
from pyomo.dae.plugins.colloc import Collocation_Discretization_Transformation

# Import Pyomo model
from pyomoExample2 import model

# Create model
instance = model.create()

# Discretize model using Radau Collocation
discretize = Collocation_Discretization_Transformation()
disc_instance = discretize.apply(instance,nfe=20,ncp=6,scheme='LAGRANGE-RADAU')

# Add objective function after model has been discretized
def obj_rule(model):
    return sum((model.x[i]-model.x_ref)**2 for i in model.time)
```

```
disc_instance.obj = Objective(rule=obj_rule)

# Solve discretized model
opt = SolverFactory('ipopt')
results = opt.solve(disc_instance)

# Load Results
disc_instance.load(results)
```

### 15.4.3 Applying Multiple Discretization Transformations

Discretizations can be applied independently to each ContinuousSet in a model. This allows the user great flexibility in discretizing their model. For example the same numerical method can be applied with different resolutions:

```
discretize = Finite_Difference_Transformation()
disc_instance = discretize.apply(instance,wrt=instance.t1,nfe=10)
disc_instance = discretize.apply(instance,wrt=instance.t2,nfe=100)
```

This also allows the user to combine different methods. For example, applying the forward difference method to one ContinuousSet and the central finite difference method to another ContinuousSet:

```
discretize = Finite_Difference_Transformation()
disc_instance = discretize.apply(instance,wrt=instance.t1,scheme='FORWARD')
disc_instance = discretize.apply(instance,wrt=instance.t2,scheme='CENTRAL')
```

In addition, the user may combine finite difference and collocation discretizations. For example:

```
disc_fe = Finite_Difference_Transformation()
disc_instance = disc_fe.apply(instance,wrt=instance.t1,nfe=10)
disc_col = Collocation_Discretization_Transformation()
disc_instance = disc_col.apply(instance,wrt=instance.t2,nfe=10,ncp=5)
```

If the user would like to apply the same discretization to all ContinuousSet components in a model, just specify the discretization once without the *wrt* keyword argument. This will apply that scheme to all ContinuousSet components in the model that haven't already been discretized.

#### 15.4.4 Custom Discretization Schemes

A transformation framework along with certain utility functions has been created so that advanced users may easily implement custom discretization schemes other than those listed above. The transformation framework consists of the following steps:

- 1. Specify Discretization Options
- 2. Discretize the ContinuousSet(s)
- 3. Update Model Components
- 4. Add Discretization Equations
- 5. Return Discretized Model

If a user would like to create a custom finite difference scheme then they only have to worry about step (4) in the framework. The discretization equations for a particular scheme have been isolated from of the rest of the code for implementing the transformation. The function containing these discretization equations can be found at the top of the source code file for the transformation. For example, below is the function for the forward difference method:

```
def _forward_transform(v,s):
    """
    Applies the Forward Difference formula of order O(h) for first derivatives
    """
    def _fwd_fun(i):
        tmp = sorted(s)
        idx = tmp.index(i)
        return 1/(tmp[idx+1]-tmp[idx])*(v(tmp[idx+1])-v(tmp[idx]))
    return _fwd_fun
```

In this function, v represents the continuous variable or function that the method is being applied to. s represents the set of discrete points in the continuous domain. In order to implement a custom finite difference method, a user would have to copy the above function and just replace the equation next to the first return statement with their method.

After implementing a custom finite difference method using the above function template, the only other change that must be made is to add the custom method to the *all\_schemes* dictionary in the Finite\_Difference\_Transformation class.

In the case of a custom collocation method, changes will have to be made in steps (2) and (4) of the transformation framework. In addition to implementing the discretization equations, the user would also have to ensure that the desired collocation points are added to the ContinuousSet being discretized.

## **Chapter 16**

## **Scripts**

There are two main ways to add scripting for Pyomo models: using Python scripts and using callbacks for the pyomo command that alter or supplement its workflow.

#### Note

The examples are written to conform with the Python version 3 print function. If executed with Python version 2, the output from print statements may not look as nice.

### 16.1 Python Scripts

#### 16.1.1 Iterative Example

To illustrate Python scripts for Pyomo we consider an example that is in the file iterative1.py and is executed using the command

python iterative1.py

#### Note

This is a Python script that contains elements of Pyomo, so it is executed using the python command. The pyomo command can be used, but then there will be some strange messages at the end when Pyomo finishes the script and attempts to send the results to a solver, which is what the pyomo command does.

This script creates a model, solves it, and then adds a constraint to preclude the solution just found. This process is repeated, so the script finds and prints multiple solutions. The particular model it creates is just the sum of four binary variables. One does not need a computer to solve the problem or even to iterate over solutions. This example is provided just to illustrate some elementary aspects of scripting.

#### Note

The built-in code for printing solutions prints only non-zero variable values. So if you run this code, no variable values will be output for the first solution found because all of the variables are zero. However, other information about the solution, such as the objective value, will be displayed.

```
# iterative1.py
from pyomo.environ import *
from pyomo.opt import SolverFactory
# Create a solver
opt = SolverFactory('glpk')
# A simple model with binary variables and
# an empty constraint list.
model = AbstractModel()
model.n = Param(default=4)
model.x = Var(RangeSet(model.n), within=Binary)
def o_rule(model):
   return summation(model.x)
model.o = Objective(rule=o_rule)
model.c = ConstraintList()
# Create a model instance and optimize
instance = model.create_instance()
results = opt.solve(instance)
print(results)
# Iterate to eliminate the previously found solution
for i in range(5):
    instance.solutions.load_from(results)
    expr = 0
    for j in instance.x:
        if instance.x[j].value == 0:
            expr += instance.x[j]
        else:
            expr += (1-instance.x[j])
    instance.c.add( expr >= 1 )
    results = opt.solve(instance)
    print (results)
```

Let us now analyze this script. The first line is a comment that happens to give the name of the file. This is followed by two lines that import symbols for Pyomo:

```
# iterative1.py
from pyomo.environ import *
from pyomo.opt import SolverFactory
```

An object to perform optimization is created by calling SolverFactory with an argument giving the name of the solver.t The argument would be <code>gurobi</code> if, e.g., Gurobi was desired instead of glpk:

```
# Create a solver
opt = SolverFactory('glpk')
```

The next lines after a comment create a model. For our discussion here, we will refer to this as the base model because it will be extended by adding constraints later. (The words "base model" are not reserved words, they are just being introduced for the discussion of this example). There are no constraints in the base model, but that is just to keep it simple. Constraints could be present in the base model. Even though it is an abstract model, the base model is fully specified by these commands because it requires no external data:

```
model = AbstractModel()
model.n = Param(default=4)
```

```
model.x = Var(RangeSet(model.n), within=Binary)
def o_rule(model):
    return summation(model.x)
model.o = Objective(rule=o_rule)
```

The next line is not part of the base model specification. It creates an empty constraint list that the script will use to add constraints.

```
model.c = ConstraintList()
```

The next non-comment line creates the instantiated model and refers to the instance object with a Python variable instance. Models run using the pyomo script do not typically contain this line because model instantiation is done by the pyomo script. In this example, the create function is called without arguments because none are needed; however, the name of a file with data commands is given as an argument in many scripts.

```
instance = model.create_instance()
```

The next line invokes the solver and refers to the object contain results with the Python variable results.

```
results = opt.solve(instance)
```

The print method of the results object is invoked by the Python print command:

```
print(results)
```

The next non-comment line is a Python iteration command that will successively assign the integers from 0 to 4 to the Python variable i, although that variable is not used in script. This loop is what causes the script to generate five more solutions:

```
for i in range(5):
```

The next line associates the results obtained with the instance. This then enables direct queries of solution values in subsequent lines using variable names contained in the instance:

```
instance.solutions.load_from(results)
```

An expression is built up in the Python variable named expr. The Python variable j will be iteratively assigned all of the indexes of the variable x. For each index, the value of the variable (which was loaded by the load method just described) is tested to see if it is zero and the expression in expr is augmented accordingly. Although expr is initialized to 0 (an integer), its type will change to be a Pyomo expression when it is assigned expressions involving Pyomo variable objects:

```
expr = 0
for j in instance.x:
   if instance.x[j].value == 0:
       expr += instance.x[j]
   else:
       expr += (1-instance.x[j])
```

During the first iteration (when i is 0), we know that all values of x will be 0, so we can anticipate what the expression will look like. We know that x is indexed by the integers from 1 to 4 so we know that y will take on the values from 1 to 4 and we also know that all value of x will be zero for all indexes so we know that the value of x will be something like

```
0 + instance.x[1] + instance.x[2] + instance.x[3] + instance.x[4]
```

The value of j will be evaluated because it is a Python variable; however, because it is a Pyomo variable, the value of instance.x[j] not be used, instead the variable object will appear in the expression. That is exactly what we want in this case. When we wanted to use the current value in the if statement, we used the value method to get it.

The next line adds to the constaint list called c the requirement that the expression be greater than or equal to one:

```
instance.c.add( expr >= 1 )
```

The proof that this precludes the last solution is left as an exerise for the reader.

The final lines in the outer for loop find a solution and display it:

```
results = opt.solve(instance)
print results
```

## 16.2 Changing the Model or Data and Re-solving

The iterative1.py example illustrates how a model can be changed and then re-solved. In that example, the model is changed by adding a constraint, but the model could also be changed by altering the values of parameters. Note, however, that in these examples, we make the changes to the instance object rather than the model object so that we do not have to create a new model object. Here is the basic idea:

- 1. Create an AbstractModel (suppose it is called model)
- 2. Call model.create\_instance() to create an instance (suppose it is called instance)
- 3. Solve instance
- 4. Change someting in instance
- 5. Call presolve
- 6. Solve instance again

If instance has a parameter whose name is in ParamName with an index that is in idx, the the value in NewVal can be assigned to it using

```
getattr(instance, ParamName)[idx] = NewVal
```

For a singleton parameter named ParamName (i.e., if it is not indexed), the assignment can be made using either

```
getattr(instance, ParamName)[None] = NewVal
```

or else

```
getattr(instance, ParamName).set_value(NewVal)
```

The function getattr is provided by Python. For more information about access to Pyomo parameters, see the section in this document on Param Access.

## 16.3 Fixing Variables and Re-solving

Instead of changing model data, scripts are often used to fix variable values. The following example illustrates this.

```
# iterative2.py

from pyomo.environ import *
from pyomo.opt import SolverFactory

# Create a solver
opt = SolverFactory('cplex')

# A simple model with binary variables and
# an empty constraint list.
#
```

```
model = AbstractModel()
model.n = Param(default=4)
model.x = Var(RangeSet(model.n), within=Binary)
def o_rule(model):
    return summation(model.x)
model.o = Objective(rule=o_rule)
model.c = ConstraintList()
# Create a model instance and optimize
instance = model.create_instance()
results = opt.solve(instance)
print(results)
# "flip" the value of x[2] (it is binary)
# then solve again
instance.solutions.load_from(results)
if instance.x[2] == 0:
    instance.x[2] = 1
else:
    instance.x[2] = 0
instance.x[2].fixed = True
results = opt.solve(instance)
print(results)
```

In this example, the variables are binary. The model is solved and then the value of model.x[2] is flipped to the opposite value before solving the model again. The main lines of interest are:

```
instance.solutions.load_from(results)

if instance.x[2] == 0:
    instance.x[2] = 1

else:
    instance.x[2] = 0
instance.x[2].fixed = True
results = opt.solve(instance)
```

This could also have been accomplished by setting the upper and lower bounds:

```
instance.solutions.load_from(results)

if instance.x[2] == 0:
    instance.x[2].setlb(1)
    instance.x[2].setub(1)

else:
    instance.x[2].setlb(0)
    instance.x[2].setub(0)

results = opt.solve(instance)
```

Notice that when using the bounds, we do not set fixed to True because that would fix the variable at whatever value it presently has and then the bounds would be ignored by the solver.

For more information about access to Pyomo variables, see the section in this document on Variable Access.

Note that instance.x.fix(2) is equivalent to

```
instance.x.value = 2
instance.x.fixed = True
```

and instance.x.fix() is equivalent to instance.x.fixed =True

## 16.4 Pyomo Callbacks

Pyomo enables altering or extending its workflow through the use of callbacks that are defined in the model file. Taken together, the callbacks allow for consruction of a rich set of workflows. However, many users might be interesting in making use of only one or two of the callbacks. They are executable Python functions with pre-defined names:

- pyomo\_preprocess: Preprocessing before model construction
- pyomo\_create\_model: Constructs and returns the model object
- pyomo\_create\_modeldata: Constructs and returns a ModelData object
- pyomo\_print\_model: Display model information
- pyomo\_modify\_instance: Modify the model instance
- pyomo\_print\_instance: Display instance information
- pyomo\_save\_instance: Write the model instance to a file
- pyomo\_print\_results: Display the results of optimization
- pyomo\_save\_results: Store the optimization results
- pyomo\_postprocess: Postprocessing after optimization

Many of these functions have arguments, which must be declared when the functions are declared. This can be done either by listing the arguments, as we will show below, or by providing a dictionary for arbitrary keyword arguments in the form \*\*kwds. If the abritrary keywords are used, then the arguments are access using the get method. For example the preprocess function takes one argument (as will be described below) so the following two function will produce the same output:

```
def pyomo_preprocess(options=None):
    if options == None:
        print "No command line options were given."
    else:
        print "Command line arguments were: %s" % options
```

```
def pyomo_preprocess(**kwds):
    options = kwds.get('options',None)
    if options == None:
        print "No command line options were given."
    else:
        print "Command line arguments were: %s" % options
```

To access the various arguments using the \*\*kwds argument, use the following strings:

- options for the command line arguments dictionary
- model-options for the --model-options dictionary
- mode1 for a model object
- instance for an instance object
- results for a results object

#### 16.4.1 pyomo\_preprocess

This function has one argument, which is an enhanced Python dictionary containing the command line options given to launch Pyomo. It is called before model construction so it augments the workflow. It is defined in the model file as follows:

```
def pyomo_preprocess(options=None):
```

#### 16.4.2 pyomo\_create\_model

This function is for experts who want to replace the model creation functionality provided by the pyomo script with their own. It takes two arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo and a dictionary with the options given in the --model-options argument to the pyomo command. The function must return the model object that has been created.

#### 16.4.3 pyomo\_create\_modeldata

Users who employ ModelData objects may want to give their own method for populating the object. This function returns returns a ModelData object that will be used to instantiate the model to form an instance. It takes two arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo and a model object.

#### 16.4.4 pyomo\_print\_model

This callback is executed between model creation and instance creation. It takes two arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo and a model object.

#### 16.4.5 pyomo\_modify\_instance

This callback is executed after instance creation. It takes three arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo, a model object, and an instance object.

#### 16.4.6 pyomo\_print\_instance

This callback is executed after instance creation (and after the pyomo\_modify\_instance callback). It takes two arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo and an instance object.

#### 16.4.7 pyomo\_save\_instance

This callback also takes place after instance creation and takes It takes two arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo and an instance object.

#### 16.4.8 pyomo\_print\_results

This callback is executed after optimization. It takes three arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo, an instance object, and a results object. Note that the --print-results option provides a way to print results; this callback is intended for users who want to customize the display.

#### 16.4.9 pyomo\_save\_results

This callback is executed after optimization. It takes three arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo, an instance object, and a results object. Note that the --save-results option provides a way to store results; this callback is intended for users who want to customize the format or contents.

#### 16.4.10 pyomo\_postprocess

This callback is also executed after optimization. It also takes three arguments: an enhanced Python dictionary containing the command line options given to launch Pyomo, an instance object, and a results object.

### 16.5 Accessing Variable Values

#### 16.5.1 Primal Variable Values

Often, the point of optimization is to get optimal values of variables. The pyomo script automatically outputs the values to a file and optionally displays the non-zero values on the standard output device (usually the computer screen). Some user may want to process the values in a script. We will describe how to access a particular variable from a Python script as well as how to access all variables from a Python script and from a callback. This should enable the reader to understand how to get the access that they desire. The Iterative example given above also illustrates access to variable values.

#### 16.5.2 One Variable from a Python Script

Assuming the model has been instantiated and solved and the results have been loded back into the instance object, then we can make use of the fact that the variable is a member of the instance object and its value can be accessed using its value member. For example, suppose the model contains a variable named quant that is a singleton (has no indexes) and suppose further that the name of the instance object is instance. Then the value of this variable can be accessed using instance. quant. value. Variables with indexes can be referenced by supplying the index.

Consider the following very simple example, which is similar to the iterative example. This is a very simple model and there are no parameter values to be read from a data file, so the  $model.create_instance()$  call does not specify a file name. In this example, the value of x[2] is accessed.

```
# noiteration1.py
from pyomo.environ import *
from pyomo.opt import SolverFactory
# Create a solver
opt = SolverFactory('glpk')
# A simple model with binary variables and
# an empty constraint list.
model = AbstractModel()
model.n = Param(default=4)
model.x = Var(RangeSet(model.n), within=Binary)
def o_rule(model):
    return summation(model.x)
model.o = Objective(rule=o_rule)
model.c = ConstraintList()
# Create a model instance and optimize
instance = model.create_instance()
results = opt.solve(instance)
instance.solutions.load_from(results)
if instance.x[2].value == 0:
    print("The second index has a zero")
else:
    print("x[2]=",instance.x[2].value)
```

#### Note

If this script is run without modification, Pyomo is likely to issue a warning because there are no constraints. The warning is because some solvers may fail if given a problem instance that does not have any constraints.

#### 16.5.3 All Variables from a Python Script

As with one variable, we assume that the model has been instantiated and solved and the results have been loded back into the instance object using instance.solutions.load\_from(results) and the code includes the line from pyomo. core import Var, then we can make use of the fact that the variable is a member of the instance object and its value can be accessed using its value member. Assuming the instance object has the name instance, the following code snippet displays all variables and their values:

```
for v in instance.component_objects(Var, active=True):
    print ("Variable",v)
    varobject = getattr(instance, v)
    for index in varobject:
        print (" ",index, varobject[index].value)
```

This code could be improved by checking to see if the variable is not indexed (i.e., the only index value is None), then the code could print the value without the word None next to it.

Assuming again that the model has been instantiated and solved and the results have been loded back into the instance object using instance.solutions.load\_from(results) and that the code includes the line from pyomo.core import Var here is a code snippet for fixing all integers at their current value:

```
for v in instance.component_objects(Var, active=True):
    varobject = getattr(instance, v)
    if isinstance(varobject.domain, IntegerSet) or isinstance(varobject.domain, BooleanSet) 
        :
        print ("fixing "+str(v))
        for index in varobject:
            varobject[index].fixed = True # fix the current value
```

#### 16.5.4 All Variables from Workflow Callbacks

The pyomo\_print\_results, pyomo\_save\_results, and pyomo\_postprocess callbacks from the pyomo script take the instance as one of their arguments and the instance has the solver results at the time of the callback so the body of the callback matches the code snipped given for a Python script.

For example, if the following defintion were included in the model file, then the pyomo command would output all variables and their values (including those variables with a value of zero):

```
def pyomo_print_results(options, instance, results):
    from pyomo.core import Var
    for v in instance.component_objects(Var, active=True):
        print ("Variable "+str(v))
        varobject = getattr(instance, v)
        for index in varobject:
            print (" ",index, varobject[index].value)
```

### 16.6 Accessing Parameter Values

Access to paramaters is completely analgous to access to variables. For example, here is a code snippet to print the name and value of every Parameter:

```
from pyomo.core import Param
for p in instance.component_objects(Param, active=True):
    print ("Parameter "+str(p))
    parmobject = getattr(instance, p)
    for index in parmobject:
        print (" ",index, parmobject[index].value)
```

NOTE: The value of a Param can be returned as None+ if no data was specified for it. This will be true even if a default value was given. To inspect the default value of a Param, replace .value with .default () but note that the default might be a function.

## 16.7 Accessing Duals

Access to dual values in scripts is similar to accessing primal variable values, except that dual values are not captured by default so additional directives are needed before optimization to signal that duals are desired.

To get duals without a script, use the pyomo option --solver-suffixes=dual which will cause dual values to be included in output. Note: In addition to duals (dual), reduced costs (rc) and slack values (slack) can be requested. All suffixes can be requested using the pyomo option --solver-suffixes=.\*



#### Warning

Some of the duals may have the value None, rather than 0.

#### 16.7.1 Access Duals in a Python Script

To signal that duals are desired, declare a Suffix component with the name "dual" on the model or instance with an IMPORT or IMPORT\_EXPORT direction.

```
# Create a 'dual' suffix component on the instance
# so the solver plugin will know which suffixes to collect
instance.dual = Suffix(direction=Suffix.IMPORT)
```

See the section on [?simpara] for more information on Pyomo's Suffix component. After the results are obtained and loaded into an instance, duals can be accessed in the following fashion.

```
# display all duals
print "Duals"
from pyomo.core import Constraint
for c in instance.component_objects(Constraint, active=True):
    print (" Constraint "+str(c))
    cobject = getattr(instance, c)
    for index in cobject:
        print (" ", index, instance.dual[cobject[index]])
```

The following snippet will only work, of course, if there is a constraint with the name AxbConstraint that has and index, which is the string Film.

```
# access (display, this case) one dual
print ("Dual for Film=", instance.dual[instance.AxbConstraint['Film']])
```

Here is a complete example that relies on the file abstract2.py to provide the model and the file abstract2.dat to provide the data. Note that the model in abstract2.py does contain a constraint named AxbConstraint and abstract2.dat does specify an index for it named Film.

```
# driveabs2.py
from __future__ import division
from pyomo.environ import *
from pyomo.opt import SolverFactory

# Create a solver
opt = SolverFactory('cplex')
```

```
# get the model from another file
from abstract2 import model
# Create a model instance and optimize
instance = model.create_instance('abstract2.dat')
# Create a 'dual' suffix component on the instance
# so the solver plugin will know which suffixes to collect
instance.dual = Suffix(direction=Suffix.IMPORT)
results = opt.solve(instance)
# also puts the results back into the instance for easy access
# display all duals
print ("Duals")
from pyomo.core import Constraint
for c in instance.component_objects(Constraint, active=True):
    print (" Constraint",c)
    cobject = getattr(instance, str(c))
    for index in cobject:
        print ("
                    ", index, instance.dual[cobject[index]])
# access one dual
print ("Dual for Film=", instance.dual[instance.AxbConstraint['Film']])
```

#### 16.7.2 All Duals from Workflow Callbacks

The pyomo script needs to be instructed to obtain duals, either by using a command line option such as --solver-suffixes=dual or by adding code in the pyomo\_preprocess callback to add solver-suffixes to the list of command line arguments if it is not there and to add dual to its list of arguments if it is there, but dual is not. If a suffix with the name dual has been declared on the model the use of the command line option or pyomo\_preprocess callback is not required.

The pyomo\_print\_results, pyomo\_save\_results, and pyomo\_postprocess callbacks from the pyomo script take the instance as one of their arguments and the instance has the solver results at the time of the callback so the body of the callback matches the code snipped given for a Python script.

For example, if the following definition were included in the model file, then the pyomo command would output all constraints and their duals.

```
def pyomo_print_results(options, instance, results):
    # display all duals
    print ("Duals")
    from pyomo.core import Constraint
    for c in instance.component_objects(Constraint, active=True):
        print (" Constraint",c)
        cobject = getattr(instance, c)
        for index in cobject:
            print (" ", index, instance.dual[cobject[index]])
```

#### Note

If the --solver-suffixes command line option is used to request constraint duals, an IMPORT style Suffix component will be added to the model by the pyomo command.

## 16.8 Accessing Slacks

The functions lslack() and uslack() return the upper and lower slacks, respectively, for a constraint.

## 16.9 Accessing Solver Status

After a solve, the results object has a member Solution. Status that contains the solver status. The following snippet shows an example of access via a print statement:

```
instance = model.create()
results = opt.solve(instance)
print ("The solver returned a status of:"+str(results.Solution.Status))
```

The use of the Python str function to cast the value to a be string makes it easy to test it. In particular, the value *optimal* indicates that the solver succeeded. It is also possible to access Pyomo data that can be compared with the solver status as in the following code snippet:

## 16.10 Display of Solver Output

To see the output of the solver, use the option tee=True as in

```
results = opt.solve(instance, tee=True)
```

This can be useful for troubleshooting solver difficulties.

## 16.11 Sending Options to the Solver

Most solvers accept options and Pyomo can pass options through to a solver. In scripts or callbacks, the options can be attached to the solver object by adding to its options dictionary as illustrated by this snippet:

```
opt = SolverFactory['cbc']
opt.options["threads"] = 4
```

If multiple options are needed, then multiple dictionary entries should be added.

Sometime it is desirable to pass options as part of the call to the solve function as in this snippet:

```
results = opt.solve(instance, options="threads=4", tee=True)
```

The quoted string is passed directly to the solver. If multiple options need to be passed to the solver in this way, they should be separated by a space within the quoted string. Notice that tee is a Pyomo option and is solver-independent, while the string argument to options is passed to the solver without very little processing by Pyomo. If the solver does not have a "threads" option, it will probably complain, but Pyomo will not.

#### 16.12 Warm Starts

Some solvers support a warm start based on current values of variables. To use this feature, set the values of variables in the instance and pass warmstart=True to the solve() method. E.g.,

```
instance = model.create()
instance.y[0] = 1
instance.y[1] = 0

opt = SolverFactory("cplex")

results = opt.solve(instance, warmstart=True)
```

#### Note

The Cplex and Gurobi LP file (and Python) interfaces will generate an MST file with the variable data and hand this off to the solver in addition to the LP file.



#### Warning

Solvers using the NL file interface (e.g., "gurobi\_ampl", "cplexamp") do not accept warmstart as a keyword to the solve() method as the NL file format, by default, includes variable initialization data (drawn from the current value of all variables).

## **16.13** Solving Multiple Instances in Parallel

Use of parallel solvers for PySP is discussed in the section on Parallel PySP.

Solvers are controlled by solver servers. The pyro mip solver server is launched with the command pyro\_mip\_server. This command may be repeated to launch as many solvers as are desired. A name server and a dispatch server must be running and accessible to the process that runs the script that will use the mip servers as well as to the mip servers. The name server is launched using the command pyomo\_ns and then the dispatch server is launched with dispatch\_srvr. Note that both commands contain an underscore. Both programs keep running until terminated by an external signal, so it is common to pipe their output to a file. The commands are:

• Once: pyomo\_ns

• Once: dispatch\_srvr

• Multiple times: pyro\_mip\_server

This example demonstrates how to use these services to solve two instances in parallel.

```
# parallel.py
from __future__ import division
from pyomo.environ import *
from pyomo.opt import SolverFactory
from pyomo.opt.parallel import SolverManagerFactory
import sys

action_handle_map = {} # maps action handles to instances

# Create a solver
optsolver = SolverFactory('cplex')

# create a solver manager
```

```
# 'pyro' could be replaced with 'serial'
solver_manager = SolverManagerFactory('pyro')
if solver_manager is None:
    print "Failed to create solver manager."
    sys.exit(1)
# A simple model with binary variables and
# an empty constraint list.
model = AbstractModel()
model.n = Param(default=4)
model.x = Var(RangeSet(model.n), within=Binary)
def o_rule(model):
    return summation (model.x)
model.o = Objective(rule=o_rule)
model.c = ConstraintList()
# Create two model instances
instance1 = model.create()
instance2 = model.create()
instance2.x[1] = 1
instance2.x[1].fixed = True
# send them to the solver(s)
action_handle = solver_manager.queue(instance1, opt=optsolver, warmstart=False, tee=True,
   verbose=False)
action_handle_map[action_handle] = "Original"
action_handle = solver_manager.queue(instance2, opt=optsolver, warmstart=False, tee=True,
   verbose=False)
action_handle_map[action_handle] = "One Var Fixed"
# retrieve the solutions
for i in range(2): # we know there are two instances
    this_action_handle = solver_manager.wait_any()
    solved_name = action_handle_map[this_action_handle]
    results = solver_manager.get_results(this_action_handle)
    print "Results for", solved_name
    print results
```

This example creates two instances that are very similar and then sends them to be dispatched to solvers. If there are two solvers, then these problems could be solved in parallel (we say "could" because for such trivial problems to be actually solved in parallel, the solvers would have to be very, very slow). This example is non-sensical; the goal is simply to show <code>solver\_manager.queue</code> to submit jobs to a name server for dispatch to solver servers and <code>solver\_manager.wait\_any</code> to recover the results. The <code>wait\_all</code> function is similar, but it takes a list of action handles (returned by <code>queue</code>) as an argument and returns all of the results at once.

## 16.14 Changing the temporary directory

A "temporary" directory is used for many intermediate files. Normally, the name of the directory for temporary files is provided by the operating system, but the user can specify their own directory name. The pyomo command-line "--tempdir" option propagates through to the TempFileManager service. One can accomplish the same through the following few lines of code in a script:

```
from pyutilib.services import TempFileManager
TempfileManager.tempdir = YourDirectoryNameGoesHere
```

# **Chapter 17**

# **Pyomo Solver Interfaces**

This chapter describes how Pyomo interfaces with different solvers.

## **Chapter 18**

## **Using Black-Box Optimizers with Pyomo.Opt**

Many optimization software packages contain *black-box* optimizers, which perform optimization without using detailed knowledge of the structure of an optimization problem. Thus, black-box optimizers require a generic interface for optimization problems that defines key features of problems, like objectives and constraints.

The pyomo.opt package contains the pyomo.opt.blackbox subpackage, which provides facilities for (a) integrating Pyomo solvers with blackbox optimization applications and (b) wrapping Pyomo models for use by external blackbox optimizers. We illustrate these capabilities in this chapter with simple examples that illustrate the use of pyomo.opt.blackbox.

## 18.1 Defining and Optimizing Simple Black-Box Applications

Many black-box optimizers interact with an optimization problem by executing a separate process that computes properties of the optimization problem. This process typically reads an input file that defines the requested properties and writes an output file that contains the computed values. Unfortunately, no standards have emerged for black-box optimizers that interact with problems in this manner. Thus, different file formats are used by different optimizer software packages.

#### 18.1.1 Defining an Optimization Problem

The pyomo.opt.blackbox package provides several Python classes for optimization problems that coordinates file I/O for the user and simplifies the definition of simple black-box problems. The RealOptProblem class provides a generic interface for continuous optimization problems (i.e. with real variables). The following example defines a simple continuous optimization problem:

```
class RealProblem1 (RealOptProblem):

    def __init__(self):
        RealOptProblem.__init__(self)
        self.lower=[0.0, -1.0, 1.0, None]
        self.upper=[None, 0.0, 2.0, -1.0]
        self.nvars=4

def function_value(self, point):
        self.validate(point)
        return point.vars[0] - point.vars[1] + (point.vars[2]-1.5)**2 + (point.vars[3]+2) \leftarrow
        **4
```

This problem is equivalent to the following problem definition:

min 
$$x_0 - x_1 + (x_2 - 1.5)^2 + (x_3 + 2)^4$$
  
s.t.  $0 \le x_0$   
 $-1 \le x_1 \le 0$   
 $0 \le x_2 \le 2$   
 $x_3 \le -1$ 

Note that the problem class does *not* specify the sense of the optimization problem. These problem classes are not a complete specification of an optimization problem. Rather, an instance of a problem class can compute information about the problem that is used during optimization.

Similarly, the MixedIntOptProblem class provides a generic interface for mixed-integer optimization problems, which may contain real variables, integer variables and binary variables. The following example defines a simple mixed-integer optimization problem:

```
class MixedIntProblem1 (MixedIntOptProblem):

    def __init__(self):
        MixedIntOptProblem.__init__(self)
        self.real_lower=[0.0]*4
        self.real_upper=[2.0]*4
        self.int_lower=[-2]*3
        self.int_upper=[0]*3
        self.nnt=3
        self.nint=3
        self.nbinary=2

def function_value(self, point):
        self.validate(point)
        return sum((x-1)**2 for x in self.reals) + \
              sum((y+1)**2 for y in self.ints) + \
              sum(b for b in self.bits)
```

This problem is equivalent to the following problem definition:

$$\begin{aligned} & \min & & \sum_{i=1}^4 (x_i - 1)^2 + \sum_{i=1}^3 (y_i + 1)^2 + \sum_{i=1}^2 z_i \\ & \text{s.t.} & & 0 \leq x_i \leq 2 \\ & & & -2 \leq y_i \leq 0 \\ & & & z_i \in \{0, 1\} \end{aligned}$$

#### 18.1.2 Optimizating with Coliny Solvers

The Coliny software library supports interfaces to a variety of black-box optimizers <Coliny>. The coliny executable reads an XML specification of the optimization problem and solver, as well as a specification of how the optimizer is applied. Consider the following XML specification:

```
<!--- RealProblem1.xml
    This Coliny XML specification illustrates the execution of the
    colin:ls solver on the RealProblem1 problem.
<ColinInput>
  <Problem type="MINLP0">
     <Domain>
        <RealVars num="4">
          <Lower index="1" value="0.0"/>
          <Lower index="2" value="-1.0"/>
          <Lower index="3" value="1.0"/>
          <Upper index="2" value="0.0"/>
          <Upper index="3" value="2.0"/>
          <Upper index="4" value="-1.0"/>
        </RealVars>
     </Domain>
     <Driver>
```

This XML specification defines a MINLPO problem, which indicated that this is a mixed-integer problem that supports zero-order derivatives (i.e. no derivatives). This problem has four real variables with lower and upper bounds specified. The problem values are computed with the RealProblem1.py command-line, which defines and uses the RealProblem1 class defined above:

Note that this command is a Python script that includes the shebang character sequence on the first line. On Linux and Unix systems, this line indicates that this is a script that is executed using the python command that is found in the user environment. Thus, this example assumes that the python command has pyomo.opt installed. Since multiple versions of Python can be installed on a single computer, the XML Command element may need to be defined with an explicitly Python version. For example, if Python 2.6 is installed in /usr/local with pyomo.opt, then the Command element would look like:

```
<Command>/usr/local/bin/python26 RealProblem1.py</Command>
```

Additionally, the duplication of bounds information between RealProblem1.py and RealProblem1.xml is not strictly necessary in this example. The bounds information in RealProblem1.py is used in the validate method to verify that the point being evaluated is consistent with the bounds information. We can generally assume that the Coliny solver will only evaluate feasible points, so a simpler problem definition can be used:

```
#!/usr/bin/env python
#
RealProblem2.py
import sys
from pyomo.opt.blackbox import RealOptProblem
```

```
class RealProblem2 (RealOptProblem):

    def __init__(self):
        RealOptProblem.__init__(self)
        self.nvars=4

    def function_value(self, point):
        return point.vars[0] - point.vars[1] + (point.vars[2]-1.5)**2 + (point.vars[3]+2) \leftarrow
        **4

problem = RealProblem2()
problem.main(sys.argv)
```

The last two lines of RealProblem1.py create a problem instance and then call the main method to parse the command-line arguments. This script has the following command-line syntax:

```
RealProblem1.py <input-file> <output-file>
```

The first argument is the name of an XML input file, and the second argument is the name of an XML output file. The optimization problem class manages the parsing of the input and generation of the output file. For example, consider the following input file:

The RealProblem1.py script creates the following output file:

## 18.2 Diving Deeper

The previous section provided an overview of the how the pyomo.opt.blackbox package supports the definition of optimization problems that are solved with black-box optimizers. In this section we provide more detail about how the Python problem class can be customized, as well as details about the XML file format used to communicate with Coliny optimizers. The Dakota User Manual <Dakota> provides documentation of the file format of the input and output files used with Dakota optimizers.

Table Table 18.1 summarizes the methods of the OptProblem class that a user is likely to either use or redefine when declaring a subclass. The MixedIntOptProblem class is a convenient base class for the problems solved by most black-box optimizers, and this class provides the definition of the main, create\_point and validate methods. However, any of the remaining methods may need to be defined, depending on the problem.

| Table 18.1: Methods in the | OptProblem class. |
|----------------------------|-------------------|
|----------------------------|-------------------|

| Method | Description                                        |
|--------|----------------------------------------------------|
| _init_ | The constructor, which may be redefined to specify |
|        | problem properties.                                |

Table 18.1: (continued)

| main                        | Method that processes command-line options to create a       |
|-----------------------------|--------------------------------------------------------------|
|                             | results file from an input file.                             |
| create_point                | Create an instances of the class that defines a point in the |
|                             | search domain.                                               |
| function_value              | Returns the value of the objective function.                 |
| function_values             | Returns a list of objective function values.                 |
| gradient                    | Returns a list that represents the gradient vector at the    |
|                             | given point.                                                 |
| hessian                     | Returns a Hessian matrix.                                    |
| nonlinear_constraint_values | Returns a list of values for the constraint functions.       |
| jacobian                    | Returns a Jacobian matrix.                                   |
| validate                    | Returns True if the given point is feasible, and False       |
|                             | otherwise.                                                   |

The following detailed example illustrates the use of all of these methods in a simple application:

```
class RealProblem3 (RealOptProblem):
    def __init__(self):
        RealOptProblem.__init__(self)
        self.nvars=4
        self.ncons=4
        self.response_types = [response_enum.FunctionValue,
                                response_enum.Gradient,
                                response_enum.Hessian,
                                response_enum.NonlinearConstraintValues,
                                 response_enum.Jacobian]
    def function_value(self, point):
        return point.vars[0] - point.vars[1] + (point.vars[2]-1.5) **2 + (point.vars[3]+2) ←
            * * 4
    def gradient(self, point):
        return [1, -1, 2*(point.vars[2]-1.5), 4*(point.vars[3]+2)**3]
    def hessian(self, point):
        H = []
        H.append((2,2,2))
        \text{H.append}((3,3,12*(point.vars[3]+2)**2))
        return H
    def nonlinear_constraint_values(self, point):
        C.append( sum(point.vars) )
        C.append( sum(x**2 for x in point.vars))
        return C
    def jacobian(self, point):
        J = []
        for j in range(self.nvars):
            J.append((0,j,1))
        for j in range(self.nvars):
            J.append( (1, j, 2*point.vars[j]) )
        return J
```

The response\_types attribute defined in the constructor specifies the type of information that this class can compute. For example, consider the following input XML file:

This input file requests that the class compute all of the response values, and thus the following output is generated:

```
<?xml version="1.0" encoding="UTF-8"?>
<ColinResponse>
 <Gradient>
 </Gradient>
 <NonlinearConstraintValues>
 -0.889999999999999 4.830099999999998
 </NonlinearConstraintValues>
 <FunctionValue>
 0.2701
 </FunctionValue>
 <Hessian>
 (2, 2, 2) (3, 3, 0.12000000000000022)
 </Hessian>
 <Jacobian>
 </Jacobian>
</ColinResponse>
```

Note that the values for Jacobian and Hessian matrices are represented in a sparse manner. Currently, these are represented with a list of tuple values, though a sparse matrix representation might be supported in the future.

## **Chapter 19**

## **Distributed Optimization with Pyro**

Pyomo supports distributed computing via the Python "PYRO" package. PYRO stands for PYthon Remote Objects. Full documentation of PYRO is available from: <a href="http://pyro.sourceforge.net/">http://pyro.sourceforge.net/</a>.

The following describes a "quick-start" process for creating and using a client and multiple solvers on a single, presumably multi-core compute server. For example, an institution may have an 8-core workstation with numerous CPLEX licenses. With distributed solves under PYRO, Pyomo algorithms can take advantage of the full set of resources on a machine.

The following example assumes a unix/linux platform. The steps for Windows are qualitatively identical - the sole difference is that you can't (or at least we haven't figured out how to) put processes in the background on Windows. The work-around is simply (albeit painfully) to launch the various processes in distinct shells.

## 19.1 Step 1: Starting a Name Server

All PYRO objects communicate via a name server, which provides a well-defined point of contact through which distributed objects can interact. You can think of the name server as a phone directory.

To start the name server, type:

```
pyomo-ns
```

In general, we suggest that the output be redirected to a file, with the entire process being placed in the background:

```
pyomo-ns >& ns.out &
```

## 19.2 Step 2: Starting a Dispatch Server

With the name server up and running, the next step is to create a dispatch server. The function of the dispatch server is to route work from clients to servers - both of the latter will be established in the immediately following steps. We again assume the process is executed in the background, with the output redirected:

```
dispatch_srvr >& dispatch_srvr.out &
```

## 19.3 Step 3: Starting a MIP server

With the work dispatcher up, the next step is to create servers to do real work! Pyomo ships with a pyro\_mip\_server script, which launches a server capable of solving a single MIP at a time. This server can be invoked as follows:

```
pyro_mip_server >& pyro_mip_server1.out &
```

We can also create multiple instances of the pyro\_mip\_server, e.g., to take advantage of multiple solver licenses:

```
pyro_mip_server >& pyro_mip_server2.out &
```

With this configuration, the dispatch server "sees" two mip servers, and can route work to both.

### 19.4 Step 4: Running a Client

To take advantage of the distributed MIP servers, a Pyomo user only needs to change the type of the solver manager supplied to the various solver scripts.

For example, one can run pyomo as follows, considering the PySP example found in: pyomo/examples/pysp/farmer:

```
pyomo solve --solver-cplex --solver-manager-pyro farmer_lp.py farmer_lp.dat
```

This will execute the LP solve using one of the two mip servers established in Step 3, which might be useful if they are on remote servers.

To take advantage of parallelism, we can solve the farmer example using progressive hedging, as follows:

```
\mbox{runph } -\mbox{--solver-manager=pyro } -\mbox{--model-directory=models } -\mbox{--instance-directory=} \longleftrightarrow \mbox{scenariodata}
```

## 19.5 Moving from Multi-Core to Distributed Computation

Truly distributed computation, i.e., with the client and server components on different hosts, is only incrementally more difficult than what is outlined above. If multiple hosts are involved in the computation, the only real issue is making sure the various hosts can all find a common nameserver. After starting pyomo-ns on some host (presumably a server-class machine), the other components (dispatch\_srvr and the pyro\_mip\_server) can be pointed to the nameserver by simply setting the environment variable PYRO\_NS\_HOSTNAME to the name (or IP address) of the host running the nameserver. The same process should be followed on the client prior to executing either pyomo, runph, or some other client solver script.

We have tested this on linux clusters with success. The only issues encountered involve overly aggressive firewalls on the host running the nameserver, which was easily corrected. In theory, Pyro should also work on Windows clusters, and linux-Window hybrid clusters via the same mechanism.

## 19.6 Cleaning Up After Yourself

It is important to remember that the name server, the dispatch server, and the mip server processes are persistent, and need to be terminated when a user has completed computational experiments. Actually, that is not entirely correct - the server processes can live forever, and continue to receive work. The only issue is when multiple users are attempting to use the same compute platform, are running their own servers, etc. While this may work, we have not tested it fully yet!

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## Colophon

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