

Introduction to IPOPT:

A tutorial for downloading, installing, and using IPOPT.

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Abstract

This document is a guide to using IPOPT 3.0 (the new C++ version of IPOPT). It includes instructions on how to obtain and compile IPOPT, an description of the interface, user options, etc., as well as a tutorial on how to solve a nonlinear optimization problem with IPOPT.

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1 Introduction

IPOPT (Interior Point Optimizer, pronounced “I-P-Opt”) is an open source software package for large-scale nonlinear optimization. It can be used to solve general nonlinear programming problems of the form

$$\min_{x \in \mathbb{R}^n} f(x) \quad (1)$$

$$\text{s.t.} \quad g^L \leq g(x) \leq g^U \quad (2)$$

$$x^L \leq x \leq x^U, \quad (3)$$

where $x \in \mathbb{R}^n$ are the optimization variables (possibly with lower and upper bounds, $x^L \in (\mathbb{R} \cup \{-\infty\})^n$ and $x^U \in (\mathbb{R} \cup \{+\infty\})^n$), $f : \mathbb{R}^n \rightarrow \mathbb{R}$ is the objective function, and $g : \mathbb{R}^n \rightarrow \mathbb{R}^m$ are the general nonlinear constraints. The functions $f(x)$ and $g(x)$ can be linear or nonlinear and convex or non-convex (but should be twice continuously differentiable). The constraints, $g(x)$, have lower and upper bounds, $g^L \in (\mathbb{R} \cup \{-\infty\})^m$ and $g^U \in (\mathbb{R} \cup \{+\infty\})^m$. Note that equality constraints of the form $g_i(x) = \bar{g}_i$ can be specified by setting $g_i^L = g_i^U = \bar{g}_i$.

1.1 Mathematical Background

IPOPT implements an interior point line search filter method that aims to find a local solution of (1-3). The mathematical details of the algorithm can be found in several publications [?, ?, ?, ?, ?].

1.2 Availability

The IPOPT package is available from COIN-OR (www.coin-or.org) under the CPL (Common Public License) open-source license and includes the source code for IPOPT. This means, it is available free of

charge, also for commercial purposes. However, if you give away software including IPOPT code (in source code or binary form) and you made changes to the IPOPT source code, you are required to publicly and clearly indicate which modifications you made. After all, the goal of open source software is the continuous development and improvement of software. For details, please refer to the Common Public License.

Also, if you are using IPOPT to obtain results for a publication, we politely ask you to point out in your paper that you used IPOPT, and to cite the publication [?]. Writing high-quality numerical software takes a lot of time and effort, and does usually not translate into a high number of publications, therefore we believe this request is only fair :).

1.3 Prerequisites

In order to build IPOPT, some third party components are required:

- BLAS (Basic Linear Algebra Subroutines). Many vendors of compilers and operating systems provide precompiled and optimized libraries for these dense linear algebra subroutines. But you can also get the source code from www.netlib.org and have the IPOPT distribution compile it automatically.
- A sparse symmetric indefinite linear solver. The IPOPT needs to obtain the solution of sparse, symmetric, indefinite linear systems, and for this it relies on third-party code. Currently, your only option is the solver MA27 from the Harwell Subroutine Library (see <http://www.cse.clrc.ac.uk/nag/hsl/>). We are working on interfaces for IPOPT with some other linear solvers.

Furthermore, we also make use a subroutine for scaling the linear systems arising in IPOPT, using the HSL routine MC19. It is not required to have MC19 to compile IPOPT; if this routine is missing, the scaling is never performed.

- ASL (AMPL Solver Library). The source code available is at www.netlib.org, and the IPOPT makefiles will automatically compile it for you if you put the source code into a designated space. NOTE: This is only required if you want to use IPOPT from AMPL and want to compile the IPOPT AMPL solver executable. solver

For more information on third-party components and how to obtain them, see Section 2.2.

Since the IPOPT code is written in C++, you will need a C++ compiler to build the IPOPT library. We tried very hard to write the code as platform and compiler independent as possible.

In addition, since MA27 is currently the only linear solver interfaced with IPOPT, and MA27 is written in Fortran, you will also need a Fortran compiler in order to compile MA27. However, we believe that it is possible to use the Fortran-to-C compiler `f2c` from www.netlib.org to convert the MA27 source code to C, and then compile the resulting C files with a C compiler and create a library containing the required HSL code. But so far we haven't tested this, and currently the configuration script for IPOPT looks for a Fortran compiler.

1.4 How to use IPOPT

If desired, the IPOPT distribution generates an executable for the modeling environment AMPL. As well, you can link your problem statement with IPOPT using interfaces for Fortran, C, or C++. IPOPT can be used with most Linux/Unix environments, and on Windows using Visual Studio .NET or Cygwin. Below in Section 3 this document demonstrate how to solve problems using IPOPT. This includes installation and compilation of IPOPT for use with AMPL as well as linking with your own code.

Finally, the IPOPT distribution includes an interface for CUTEr (see <http://cutter.rl.ac.uk/cuter-www/>), if you want to use IPOPT to solve problems modeled in SIF.

The old version of IPOPT has been interface with Matlab, and is also available from NEOS, and the new version will be available through similar means in the future. Please check the IPOPT homepage for updates.

1.5 More Information

More and up-to-date information can be found at the IPOPT homepage,

<http://projects.coin-or.org/Ipoprt>.

Here, you can find FAQs, some (hopefully useful) hints, etc.

IPOPT is an open source project, and we encourage people to contribute code (such as interfaces to appropriate linear solvers, modeling environments, or even algorithmic features). If you are interested in contributing code, please have a look at the COIN page <http://www.coin-or.org/contributions.html>, and contact the IPOPT project leader.

There is also a mailing list for IPOPT, <http://list.coin-or.org/mailman/listinfo/coin-ipopt>, where you can subscribe to get notified of updates, and to ask general questions regarding installation and usage. (You might want to look at the archives before posting a question.)

We try to answer questions posted to the mailing list in a reasonable manner. But you will understand that we cannot answer all questions in detail, and we will not be able to help you model and debug your particular optimization problem. However, if you have a challenging optimization problem and might be interested in consulting services by IBM Research, please contact the IPOPT project leader, Andreas Wächter..

1.6 History of IPOPT

The original IPOPT (Fortran version) was a product of the dissertation research of Andreas Wächter [?], under Lorenz T. Biegler at the Chemical Engineering Department at Carnegie Mellon University. The code was made open source and distributed by the COIN-OR initiative, which is now a non-profit corporation. IPOPT has been actively developed under COIN-OR since 2002.

To continue natural extension of the code and allow easy addition of new features, IBM Research decided to invest in an open source re-write of IPOPT in C++. The new C++ version of the IPOPT optimization code (IPOPT 3.0 and beyond) is currently developed at IBM Research and remains part of the COIN-OR initiative. Future development on the Fortran version will cease with the exception of occasional bug fix releases.

2 Installing IPOPT

The following sections describe the installation procedures on UNIX/Linux systems. For installation instructions on Windows see Section 2.4.

2.1 Getting the IPOPT Code

IPOPT is available from the COIN-OR subversion repository. You can either download the code using *svn* (the *subversion*⁴ client similar to CVS) or simply retrieve a tarball (compressed archive file). While

⁴see <http://subversion.tigris.org/>

the tarball is an easy method to retrieve the code, using the *subversion* system allows users the benefits of the version control system, including easy updates and revision control.

2.1.1 Getting the IPOPT code via subversion

Of course, the *subversion* client must be installed on your system if you want to obtain the code this way (the executable it called `svn`). Information about *subversion* and how to download it can be found at <http://subversion.tigris.org/>.

To obtain the IPOPT source code via subversion, follow the steps below:

1. Create a directory to store the code

```
$ mkdir Ipop
```

Note the `$` indicates the command line prompt, do not type `$`, only the text following it.

2. Download the code to the new directory

```
$ cd Ipop
```

```
$ svn co https://www.coin-or.org/svn/Ipop/trunk
```

3. Change into the root directory of the IPOPT distribution

```
$ cd trunk
```

In the following, “`$IPOPTDIR`” will refer to the directory in which you are right now (output of `pwd`).

2.1.2 Getting the IPOPT code as tarball

To use the tarball, follow the steps below:

1. Download the latest tarball from <http://www.coin-or.org/Tarballs>. The file you should look for has the form `ipop-3.x.x.tar.gz` (where “3.x.x.” is the version number). Put this file in a directory under which you want to put the IPOPT installation.

2. Issue the following commands to unpack the archive file:

```
$ gunzip ipop-3.x.x.tar.gz
```

```
$ tar xvf ipop-3.x.x.tar
```

Note the `$` indicates the command line prompt, do not type `$`, only the text following it.

3. Change into the root directory of the IPOPT distribution

```
$ cd ipop-3.x.x
```

In the following, “`$IPOPTDIR`” will refer to the directory in which you are right now (output of `pwd`).

2.2 Download External Code

IPOPT uses a few external packages that are not included in the distribution, namely ASL (the Ampl Solver Library), BLAS, and some sparse linear algebra routines from the Harwell Subroutine Library.

Note that you only need to obtain the ASL if you intend to use IPOPT from AMPL. It is not required if you want to specify your optimization problem in a programming language (C++, C, or Fortran).

2.2.1 Download BLAS and ASL

If you have the download utility `wget` installed on your system, retrieving ASL, and BLAS is straightforward using scripts included with the `ipopt` distribution. These scripts download the required files from the Netlib Repository (www.netlib.org).

```
$ cd $IPOPTDIR/Extern/blas
$ ./get.blas
$ cd ../ASL
$ ./get.asl
```

If you don't have `wget` installed on your system, please read the `INSTALL.*` files in the `$IPOPTDIR/Extern/blas` and `$IPOPTDIR/Extern/ASL` directories for alternative instructions.

2.2.2 Download HSL Subroutines

In addition to the IPOPT source code, two additional subroutines have to be downloaded from the Harwell Subroutine Library (HSL). The required routines are freely available for non-commercial, academic use, but it is your responsibility to investigate the licensing of all third party code.

1. Go to <http://hsl.rl.ac.uk/archive/hslarchive.html>
2. Follow the instruction on the website, read the license, and submit the registration form.
3. Go to *HSL Archive Programs*, and find the package list.
4. In your browser window, click on *MA27*.
5. Make sure that *Double precision:* is checked. Click *Download package (comments removed)*
6. Save the file as `ma27ad.f` in `$IPOPTDIR/Extern/HSL/`
Note: Some browsers append a file extension (`.txt`) when you save the file, so you may have to rename it.
7. Go back to the package list using the back button of your browser.
8. In your browser window, click on *MC19*.
9. Make sure *Double precision:* is checked. Click *Download package (comments removed)*
10. Save the file as `mc19ad.f` in `$IPOPTDIR/Extern/HSL/`
Note: Some browsers append a file extension (`.txt`) when you save the file, so you may have to rename it.

2.3 Compiling and Installing IPOPT

IPOPT can be easily compiled and installed with the usual `configure`, `make`, `make install` commands. Below are the basic steps that should work on most systems. For special compilations and some for troubleshooting see Appendix D and consult the IPOPT webpage before sending a message to the mailing list.

1. Go to the main directory of IPOPT:
\$ cd \$IPOPTDIR

2. Run the configure script

```
$ ./configure
```

If the last output line of the script reads “configure: Configuration successful” then everything worked fine. Otherwise, look at the screen output, have a look at the `config.log` output file and/or consult Appendix D.

The default configure (without any options) is sufficient for most users. If you want to see the configure options, consult Appendix D.

3. Build the code

```
$ make
```

4. Install IPOPT

```
$ make install
```

This installs

- the IPOPT AMPL solver executable (if ASL source was downloaded) in `$IPOPTDIR/bin`,
- the IPOPT library (`libipopt.a`) in `$IPOPTDIR/lib`,
- and the necessary header files in `$IPOPTDIR/include/ipopt`.

You can change the default installation directory (here `$IPOPTDIR`) to something else (such as `/usr/local`) by using the `--prefix` switch for `configure`.

2.4 Installation on Windows

There are two ways to install IPOPT on Windows systems. The first option, described in section 2.4.1, is to use Cygwin (see www.cygwin.com), which offers a UNIX-like environment on Windows and in which the installation procedure described earlier in this section can be used. The IPOPT distribution also includes projects files for the Microsoft Visual Studio (see Section 2.4.2).

2.4.1 Installation with Cygwin

Cygwin is a Linux-like environment for Windows; if you don’t know what it is you might want to have a look at the Cygwin homepage, www.cygwin.com.

It is possible to build the IPOPT AMPL solver executable in Cygwin for general use in Windows. You can also hook up IPOPT to your own program if you compile it in the Cygwin environment⁵.

If you want to compile IPOPT under Cygwin, you first have to install Cygwin in your Windows system. This is pretty straight forward; you simply download the “setup” program from www.cygwin.com and start it.

Then you do the following steps (assuming here that you don’t have any complications with firewall settings etc - in that case you might have to choose some connection settings differently):

1. Click next
2. Select “install from the internet” (default) and click next
3. Select a directory where Cygwin is to be installed (you can leave the default) and choose all other things to your liking, then click next

⁵It might also be possible to build an IPOPT DLL that can be used from non-cygwin compilers, but this is not (yet?) supported.

4. Select a temp dir for Cygwin setup to store some files (if you put it on your desktop you will later remember to delete it)
5. Select “direct connection” (default) and click next
6. Select some mirror site that seems close by to you and click next
7. OK, now comes the complicated part:

You need to select the packages that you want to have installed. By default, there are already selections, but the compilers are usually not pre-chosen. You need to make sure that you select the GNU compilers (for Fortran, C, and C++ — together with the MinGW options), the GNU Make, and Subversion. For this, click on the “Devel” branch (which opens a subtree) and select:

- gcc
- gcc-core
- gcc-g77
- gcc-g++
- gcc-mingw
- gcc-mingw-core
- gcc-mingw-g77
- gcc-mingw-g++
- make
- subversion

Then, in the “Web” branch, please select “wget” (which will make the installation of third party dependencies for IPOPT easier)

This will automatically also select some other packages.

8. Then you click on next, and Cygwin will be installed (follow the rest of the instructions and choose everything else to your liking). At a later point you can easily add/remove packages with the setup program.
9. Now that you have Cygwin, you can open a Cygwin window, which is like a UNIX shell window.
10. Now you just follow the instructions in the beginning of Sections 2: You download the IPOPT code into your Cygwin home directory (from the Windows explorer that is usually something like C:\Cygwin\home\your_user_name). After that you obtain the third party code (like on Linux/UNIX), type

```
./configure CXX=g++ CC=gcc F77=g77
```

and

```
make install
```

in the correct directories, and hopefully that will work. The IPOPT AMPL solver executable will be in the subdirectory bin (called “ipopt.exe”).

2.4.2 Using the Visual Studio

The IPOPT distribution includes project files that can be used to compile the IPOPT library and a Fortran and C++ example within the Microsoft Visual Studio. The project files have been created with Microsoft Visual C++ .NET 2003 Standard, and the Intel Visual Fortran Compiler 8.

In order to use those project files, download the IPOPT source code, as well as the required third party code (put it into the Extern/blas and Extern/HSL directories—ASL is not required for the Fortran and C examples). Then open the solution file

```
$IPOPTDIR\Windows\VisualStudio_dotNET\Ipopt\Ipopt.sln
```

Note: Since the project files were created only with the Standard edition of the C++ compiler, code optimization might be disabled; for fast performance make sure you enable code optimization.

3 Interfacing your NLP to IPOPT: A tutorial example.

IPOPT has been designed to be flexible for a wide variety of applications, and there are a number of ways to interface with IPOPT that allow specific data structures and linear solver techniques. Nevertheless, the authors have included a standard representation that should meet the needs of most users.

This tutorial will discuss four interfaces to IPOPT, namely the AMPL modeling language interface, and the C, C++, and Fortran code interfaces. AMPL is a 3rd party modeling language tool that allows users to write their optimization problem in a syntax that resembles the way the problem would be written mathematically. Once the problem has been formulated in AMPL, the problem can be easily solved using the (already compiled) executable. Interfacing your problem by directly linking code requires more effort to write, but can be far more efficient for large problems.

We will illustrate how to use each of the four interfaces using an example problem, number 71 from the Hock-Schittkowsky test suite [?],

$$\min_{x \in \mathbb{R}^4} \quad x_1 x_4 (x_1 + x_2 + x_3) + x_3 \quad (4)$$

$$\text{s.t.} \quad x_1 x_2 x_3 x_4 \geq 25 \quad (5)$$

$$x_1^2 + x_2^2 + x_3^2 + x_4^2 = 40 \quad (6)$$

$$1 \leq x_1, x_2, x_3, x_4 \leq 5, \quad (7)$$

with the starting point

$$x = (1, 5, 5, 1) \quad (8)$$

and the optimal solution

$$x^* = (1.00000000, 4.74299963, 3.82114998, 1.37940829).$$

3.1 Using the AMPL interface

Interfacing through the AMPL interface is by far the easiest way to solve a problem with IPOPT. The user must simply formulate the problem in AMPL syntax, and solve the problem through the AMPL environment. There are drawbacks, however. AMPL is a 3rd party package and, as such, must be appropriately licensed (a free, student version for limited problem size is available from the AMPL website,

```

# tell ampl to use the ipopt executable as a solver
# make sure ipopt is in the path!
option solver ipopt;

# declare the variables and their bounds,
# set notation could be used, but this is straightforward
var x1 >= 1, <= 5;
var x2 >= 1, <= 5;
var x3 >= 1, <= 5;
var x4 >= 1, <= 5;

# specify the objective function
minimize obj:
    x1 * x4 * (x1 + x2 + x3) + x3;

# specify the constraints
s.t.
    inequality:
        x1 * x2 * x3 * x4 >= 25;

    equality:
        x1^2 + x2^2 + x3^2 + x4^2 = 40;

# specify the starting point
let x1 := 1;
let x2 := 5;
let x3 := 5;
let x4 := 1;

# solve the problem
solve;

# print the solution
display x1;
display x2;
display x3;
display x4;

```

Figure 1: AMPL model file `hs071_ampl.mod`

www.ampl.com). Furthermore, the AMPL environment may be prohibitive for very large problems. Nevertheless, formulating the problem in AMPL is straightforward and even for large problems, it is often used as a prototyping tool before using one of the code interfaces.

This tutorial is not intended as a guide to formulating models in AMPL. If you are not already familiar with AMPL, please consult [?].

The problem presented in equations (4)–(8) can be solved with IPOPT with the AMPL model file given in Figure 1.

The line, “`option solver ipopt;`” tells AMPL to use IPOPT as the solver. The IPOPT executable (installed in Section 2.3) must be in the path for AMPL to find it. The remaining lines specify the problem in AMPL format. The problem can now be solved by starting AMPL and loading the mod file:

```

$ ampl
> model hs071_ampl.mod;
.

```

1. Problem dimensions
 - number of variables
 - number of constraints
2. Problem bounds
 - variable bounds
 - constraint bounds
3. Initial starting point
 - Initial values for the primal x variables
 - Initial values for the multipliers (only required for a warm start option)
4. Problem Structure
 - number of nonzeros in the Jacobian of the constraints
 - number of nonzeros in the Hessian of the Lagrangian
 - Structure of the Jacobian of the constraints
 - Structure of the Hessian of the Lagrangian
5. Evaluation of Problem Functions

Information evaluated using a given point $(x_k, \lambda_k, \sigma_f)$ coming from IPOPT

 - Objective function, $f(x_k)$
 - Gradient of the objective $\nabla f(x_k)$
 - Constraint residuals, $g(x_k)$
 - Jacobian of the constraints, $\nabla g(x_k)$
 - Hessian of the Lagrangian, $\sigma_f \nabla^2 f(x_k) + \sum_{i=1}^m \lambda_i \nabla^2 g_i(x_k)$

Figure 2: Information Required By IPOPT

.

The problem will be solved using IPOPT and the solution will be displayed.

At this point, AMPL users may wish to skip the sections about interfacing with code, but should read Section 4 concerning IPOPT options, and Section 5 which explains the output displayed by IPOPT.

3.2 Interfacing with IPOPT through code

In order to solve a problem, IPOPT needs more information than just the problem definition (for example, the derivative information). If you are using a modeling language like AMPL, the extra information is provided by the modeling tool and the IPOPT interface. When interfacing with IPOPT through your own code, however, you must provide this additional information.

The information required by IPOPT is shown in Figure 2. The problem dimensions and bounds are

straightforward and come solely from the problem definition. The initial starting point is used by the algorithm when it begins iterating to solve the problem. If IPOPT has difficulty converging, or if it converges to a locally infeasible point, adjusting the starting point may help.

Providing the problem structure is a bit more involved. IPOPT is a nonlinear programming solver that is designed for solving large scale, sparse problems. While IPOPT can be customized for a variety of matrix formats, the triplet format is used for the standard interfaces in this tutorial. For an overview of the triplet format for sparse matrices, see Appendix A. Before solving the problem, IPOPT needs to know the number of nonzeros and the structure (row and column indices of each of the nonzeros) of the Jacobian and the Hessian. Once defined, this nonzero structure **MUST** remain constant for the entire problem. This means that the structure needs to include entries for any element that could ever be nonzero, not only those that are nonzero at the starting point.

As IPOPT iterates, it will need the values for the items in (5) evaluated at particular points. Before we can begin coding the interface, however, we need to work out the details of these equations symbolically for example problem (4-7).

The gradient of the objective $f(x)$ is given by

$$\begin{bmatrix} x_1 x_4 + x_4(x_1 + x_2 + x_3) \\ x_1 x_4 \\ x_1 x_4 + 1 \\ x_1(x_1 + x_2 + x_3) \end{bmatrix}, \quad (9)$$

the Jacobian of the constraints $g(x)$ is,

$$\begin{bmatrix} x_2 x_3 x_4 & x_1 x_3 x_4 & x_1 x_2 x_4 & x_1 x_2 x_3 \\ 2x_1 & 2x_2 & 2x_3 & 2x_4 \end{bmatrix}. \quad (10)$$

We need to determine the Hessian of the Lagrangian. The Lagrangian is given by $f(x) + g(x)^T \lambda$ and the Hessian of the Lagrangian is technically, $\nabla^2 f(x_k) + \sum_{i=1}^m \lambda_i \nabla^2 g_i(x_k)$. However, so that IPOPT can ask for the Hessian of the objective or the constraints independently if required, we introduce a factor (σ_f) in front of the objective term. The value for σ_f is generally 1, although it may include scaling factors or even be set to zero to retrieve the Hessians of the constraints alone.

For our implementation then, the symbolic form of the Hessian of the Lagrangian

$$\sigma_f \nabla^2 f(x_k) + \sum_{i=1}^m \lambda_i \nabla^2 g_i(x_k)$$

(with the σ_f parameter) is,

$$\sigma_f \begin{bmatrix} 2x_4 & x_4 & x_4 & 2x_1 + x_2 + x_3 \\ x_4 & 0 & 0 & x_1 \\ x_4 & 0 & 0 & x_1 \\ 2x_1 + x_2 + x_3 & x_1 & x_1 & 0 \end{bmatrix} + \lambda_1 \begin{bmatrix} 0 & x_3 x_4 & x_2 x_4 & x_2 x_3 \\ x_3 x_4 & 0 & x_1 x_4 & x_1 x_3 \\ x_2 x_4 & x_1 x_4 & 0 & x_1 x_2 \\ x_2 x_3 & x_1 x_3 & x_1 x_2 & 0 \end{bmatrix} + \lambda_2 \begin{bmatrix} 2 & 0 & 0 & 0 \\ 0 & 2 & 0 & 0 \\ 0 & 0 & 2 & 0 \\ 0 & 0 & 0 & 2 \end{bmatrix} \quad (11)$$

where the first term comes from the Hessian of the objective function, and the second and third term from the Hessian of (5) and (6) respectively. Therefore, the dual variables λ_1 and λ_2 are then the multipliers for constraints (5) and (6) respectively.

The remaining sections of the tutorial will lead you through the coding required to solve example problem (4)–(7) using, first C++, then C, and finally Fortran. Completed versions of these examples can be found in \$IPOPTDIR/Examples under `hs071_c.cpp`, `hs071_c`, `hs071.f`.

As a user, you are responsible for coding two sections of the program that solves a problem using IPOPT: the main executable (e.g., `main`) and the problem representation. Typically, you will write an executable that prepares the problem, and then passes control over to IPOPT through an `Optimize` or `Solve` call. In this call, you will give IPOPT everything that it requires to call back to your code whenever it needs functions evaluated (like the objective, the Jacobian, etc.). In each of the three sections that follow (C++, C, and Fortran), we will first discuss how to code the problem representation, and then how to code the executable.

3.3 The C++ Interface

This tutorial assumes that you are familiar with the C++ programming language, however, we will lead you through each step of the implementation. For the problem representation, we will create a class that inherits off of the pure virtual base class, `TNLP` (`IpTNLP.hpp`). For the executable (the `main` function) we will make the call to IPOPT through the `IpoptApplication` class (`IpIpoptApplication.hpp`). In addition, we will also be using the `SmartPtr` class (`IpSmartPtr.hpp`) which implements a reference counting pointer that takes care of memory management (object deletion) for you.

After “make install” (see Section 2.3), the header files are installed in `$IPOPTDIR/include/ipopt` (or in `$PREFIX/include/ipopt` if the switch `--prefix=$PREFIX` was used for configure).

3.3.1 Coding the Problem Representation

We provide the information required in Figure 2 by coding the `HS071_NLP` class, a specific implementation of the `TNLP` base class. In the executable, we will create an instance of the `HS071_NLP` class and give this class to IPOPT so it can evaluate the problem functions through the `TNLP` interface. If you have any difficulty as the implementation proceeds, have a look at the completed example in the `Example/hs071.cpp` directory.

Start by creating a new directory under Examples, called `MyExample` and create the files `hs071_nlp.hpp` and `hs071_nlp.cpp`. In `hs071_nlp.hpp`, include `IpTNLP.hpp` (the base class), tell the compiler that we are using the IPOPT namespace, and create the declaration of the `HS071_NLP` class, inheriting off of `TNLP`. Have a look at the `TNLP` class in `IpTNLP.hpp`; you will see eight pure virtual methods that we must implement. Declare these methods in the header file. Implement each of the methods in `HS071_NLP.cpp` using the descriptions given below. In `hs071_nlp.cpp`, first include the header file for your class and tell the compiler that you are using the IPOPT namespace. A full version of these files can be found in the `Examples/hs071.cpp` directory.

Method `get_nlp_info` with Prototype

```
virtual bool get_nlp_info(Index& n, Index& m, Index& nnz_jac_g,
                        Index& nnz_h_lag, IndexStyleEnum& index_style)
```

Give IPOPT the information about the size of the problem (and hence, the size of the arrays that it needs to allocate).

- `n`: (out), the number of variables in the problem (dimension of `x`).
- `m`: (out), the number of constraints in the problem (dimension of `g`).
- `nnz_jac_g`: (out), the number of nonzero entries in the Jacobian.
- `nnz_h_lag`: (out), the number of nonzero entries in the Hessian.

- `index_style`: (out), the style used for row/col entries in the sparse matrix format (C_STYLE: 0-based, FORTRAN_STYLE: 1-based).

IPOPT uses this information when allocating the arrays that it will later ask you to fill with values. Be careful in this method since incorrect values will cause memory bugs which may be very difficult to find.

Our example problem has 4 variables (n), and two constraints (m). The Jacobian for this small problem is actually dense and has 8 nonzeros (we will still represent this Jacobian using the sparse matrix triplet format). The Hessian of the Lagrangian has 10 “symmetric” nonzeros. Keep in mind that the number of nonzeros is the total number of elements that may *ever* be nonzero, not just those that are nonzero at the starting point. This information is set once for the entire problem.

```
bool HS071_NLP::get_nlp_info(Index& n, Index& m, Index& nnz_jac_g,
                             Index& nnz_h_lag, IndexStyleEnum& index_style)
{
    // The problem described in HS071_NLP.hpp has 4 variables, x[0] through x[3]
    n = 4;

    // one equality constraint and one inequality constraint
    m = 2;

    // in this example the Jacobian is dense and contains 8 nonzeros
    nnz_jac_g = 8;

    // the Hessian is also dense and has 16 total nonzeros, but we
    // only need the lower left corner (since it is symmetric)
    nnz_h_lag = 10;

    // use the C style indexing (0-based)
    index_style = TNLP::C_STYLE;

    return true;
}
```

Method `get_bounds_info` with Prototype

```
virtual bool get_bounds_info(Index n, Number* x_l, Number* x_u,
                             Index m, Number* g_l, Number* g_u)
```

Give IPOPT the value of the bounds on the variables and constraints.

- `n`: (in), the number of variables in the problem (dimension of \mathbf{x}).
- `x_l`: (out) the lower bounds for \mathbf{x} .
- `x_u`: (out) the upper bounds for \mathbf{x} .
- `m`: (in), the number of constraints in the problem (dimension of \mathbf{g}).
- `g_l`: (out) the lower bounds for \mathbf{g} .
- `g_u`: (out) the upper bounds for \mathbf{g} .

The values of n and m that you specified in `get_nlp_info` are passed to you for debug checking. Setting a lower bound to a value less than or equal to the value of the option `nlp_lower_bound_inf` (data member of TNLP) will cause IPOPT to assume no lower bound. Likewise, specifying the upper bound above or

equal to the value of the option `nlp_upper_bound_inf` will cause IPOPT to assume no upper bound. These options, `nlp_lower_bound_inf` and `nlp_upper_bound_inf`, are set to -10^{19} and 10^{19} respectively, by default, but may be modified by changing the options (see Section 4).

In our example, the first constraint has a lower bound of 25 and no upper bound, so we set the lower bound of constraint [0] to 25 and the upper bound to some number greater than 10^{19} . The second constraint is an equality constraint and we set both bounds to 40. IPOPT recognizes this as an equality constraint and does not treat it as two inequalities.

```
bool HS071_NLP::get_bounds_info(Index n, Number* x_l, Number* x_u,
                                Index m, Number* g_l, Number* g_u)
{
    // here, the n and m we gave IPOPT in get_nlp_info are passed back to us.
    // If desired, we could assert to make sure they are what we think they are.
    assert(n == 4);
    assert(m == 2);

    // the variables have lower bounds of 1
    for (Index i=0; i<4; i++) {
        x_l[i] = 1.0;
    }

    // the variables have upper bounds of 5
    for (Index i=0; i<4; i++) {
        x_u[i] = 5.0;
    }

    // the first constraint g1 has a lower bound of 25
    g_l[0] = 25;
    // the first constraint g1 has NO upper bound, here we set it to 2e19.
    // Ipopt interprets any number greater than nlp_upper_bound_inf as
    // infinity. The default value of nlp_upper_bound_inf and nlp_lower_bound_inf
    // is 1e19 and can be changed through ipopt options.
    g_u[0] = 2e19;

    // the second constraint g2 is an equality constraint, so we set the
    // upper and lower bound to the same value
    g_l[1] = g_u[1] = 40.0;

    return true;
}
```

Method `get_starting_point` with Prototype

```
virtual bool get_starting_point(Index n, bool init_x, Number* x,
                                bool init_z, Number* z_L, Number* z_U,
                                Index m, bool init_lambda, Number* lambda)
```

Give IPOPT the starting point before it begins iterating.

- `n`: (in), the number of variables in the problem (dimension of `x`).
- `init_x`: (in), if true, this method must provide an initial value for `x`.
- `x`: (out), the initial values for the primal variables, `x`.
- `init_z`: (in), if true, this method must provide an initial value for the bound multipliers `z_L`, and `z_U`.

- `z_L`: (out), the initial values for the bound multipliers, `z_L`.
- `z_U`: (out), the initial values for the bound multipliers, `z_U`.
- `m`: (in), the number of constraints in the problem (dimension of `g`).
- `init_lambda`: (in), if true, this method must provide an initial value for the constraint multipliers, `lambda`.
- `lambda`: (out), the initial values for the constraint multipliers, `lambda`.

The index variables n , and m are passed in only to help your debugging. These variables will have the same values you specified in `get_nlp_info`.

Depending on the options that have been set, IPOPT may or may not require bounds for the primal variables x , the bound multipliers z , and the equality multipliers λ . The boolean flags `init_x`, `init_z`, and `init_lambda` tell you whether or not you should provide initial values for x , z , or λ respectively. The default options only require an initial value for the primal variables x .

In our example, we provide initial values for x as specified in the example problem. We do not provide any initial values for the dual variables, but use an assert to immediately let us know if we are ever asked for them.

```
bool HS071_NLP::get_starting_point(Index n, bool init_x, Number* x,
                                   bool init_z, Number* z_L, Number* z_U,
                                   Index m, bool init_lambda,
                                   Number* lambda)
{
    // Here, we assume we only have starting values for x, if you code
    // your own NLP, you can provide starting values for the dual variables
    // if you wish to use a warmstart option
    assert(init_x == true);
    assert(init_z == false);
    assert(init_lambda == false);

    // initialize to the given starting point
    x[0] = 1.0;
    x[1] = 5.0;
    x[2] = 5.0;
    x[3] = 1.0;

    return true;
}
```

Method `eval_f` with Prototype

```
virtual bool eval_f(Index n, const Number* x,
                    bool new_x, Number& obj_value)
```

Return the value of the objective function as calculated using x .

- `n`: (in), the number of variables in the problem (dimension of x).
- `x`: (in), the current values for the primal variables, x .
- `new_x`: (in), false if any evaluation method was previously called with the same values in x , true otherwise.

- `obj_value`: (out) the value of the objective function ($f(x)$).

The boolean variable `new_x` will be false if the last call to any of the evaluation methods used the same x values. This can be helpful when users have efficient implementations that calculate multiple outputs at once. IPOPT internally caches results from the TNLP and generally, this flag can be ignored.

The index variable n is passed in only to help your debugging. This variable will have the same value you specified in `get_nlp_info`.

For our example, we ignore the `new_x` flag and calculate the objective.

```
bool HS071_NLP::eval_f(Index n, const Number* x, bool new_x, Number& obj_value)
{
    assert(n == 4);

    obj_value = x[0] * x[3] * (x[0] + x[1] + x[2]) + x[2];

    return true;
}
```

Method `eval_grad_f` with Prototype

```
virtual bool eval_grad_f(Index n, const Number* x, bool new_x,
                        Number* grad_f)
```

Return the gradient of the objective to IPOPT, as calculated by the values in x .

- `n`: (in), the number of variables in the problem (dimension of x).
- `x`: (in), the current values for the primal variables, x .
- `new_x`: (in), false if any evaluation method was previously called with the same values in x , true otherwise.
- `grad_f`: (out) the array of values for the gradient of the objective function ($\nabla f(x)$).

The gradient array is in the same order as the x variables (i.e. the gradient of the objective with respect to $x[2]$ should be put in `grad_f[2]`).

The boolean variable `new_x` will be false if the last call to any of the evaluation methods used the same x values. This can be helpful when users have efficient implementations that calculate multiple outputs at once. IPOPT internally caches results from the TNLP and generally, this flag can be ignored.

The index variable n is passed in only to help your debugging. This variable will have the same value you specified in `get_nlp_info`.

In our example, we ignore the `new_x` flag and calculate the values for the gradient of the objective.

```
bool HS071_NLP::eval_grad_f(Index n, const Number* x, bool new_x, Number* grad_f)
{
    assert(n == 4);

    grad_f[0] = x[0] * x[3] + x[3] * (x[0] + x[1] + x[2]);
    grad_f[1] = x[0] * x[3];
    grad_f[2] = x[0] * x[3] + 1;
    grad_f[3] = x[0] * (x[0] + x[1] + x[2]);

    return true;
}
```

Method `eval_g` with Prototype

```
virtual bool eval_g(Index n, const Number* x,
                    bool new_x, Index m, Number* g){}
```

Give IPOPT the value of the constraints as calculated by the values in **x**.

- **n**: (in), the number of variables in the problem (dimension of **x**).
- **x**: (in), the current values for the primal variables, **x**.
- **new_x**: (in), false if any evaluation method was previously called with the same values in **x**, true otherwise.
- **m**: (in), the number of constraints in the problem (dimension of **g**).
- **g**: (out) the array of constraint residuals.

The values returned in **g** should be only the $g(x)$ values, do not add or subtract the bound values g_l or g_u .

The boolean variable **new_x** will be false if the last call to any of the evaluation methods used the same x values. This can be helpful when users have efficient implementations that calculate multiple outputs at once. IPOPT internally caches results from the TNLP and generally, this flag can be ignored.

The index variables n , and m are passed in only to help your debugging. These variables will have the same values you specified in `get_nlp_info`.

In our example, we ignore the **new_x** flag and calculate the values for the gradient of the objective.

```
bool HS071_NLP::eval_g(Index n, const Number* x, bool new_x, Index m, Number* g)
{
    assert(n == 4);
    assert(m == 2);

    g[0] = x[0] * x[1] * x[2] * x[3];
    g[1] = x[0]*x[0] + x[1]*x[1] + x[2]*x[2] + x[3]*x[3];

    return true;
}
```

Method `eval_jac_g` with Prototype

```
virtual bool eval_jac_g(Index n, const Number* x, bool new_x,
                        Index m, Index nele_jac, Index* iRow,
                        Index *jCol, Number* values)
```

Return either the structure of the Jacobian of the constraints, or the values for the Jacobian of the constraints as calculated by the values in **x**.

- **n**: (in), the number of variables in the problem (dimension of **x**).
- **x**: (in), the current values for the primal variables, **x**.
- **new_x**: (in), false if any evaluation method was previously called with the same values in **x**, true otherwise.
- **m**: (in), the number of constraints in the problem (dimension of **g**).

- `n_ele1_jac`: (in), the number of nonzero elements in the Jacobian (dimension of `iRow`, `jCol`, and `values`).
- `iRow`: (out), the row indices of entries in the Jacobian of the constraints.
- `jCol`: (out), the column indices of entries in the Jacobian of the constraints.
- `values`: (out), the values of the entries in the Jacobian of the constraints.

The Jacobian is the matrix of derivatives where the derivative of constraint i with respect to variable j is placed in row i and column j . See Appendix A for a discussion of the sparse matrix format used in this method.

If the `iRow` and `jCol` arguments are not NULL, then IPOPT wants you to fill in the structure of the Jacobian (the row and column indices only). At this time, the `x` argument and the `values` argument will be NULL.

If the `x` argument and the `values` argument are not NULL, then IPOPT wants you to fill in the values of the Jacobian as calculated from the array `x` (using the same order as you used when specifying the structure). At this time, the `iRow` and `jCol` arguments will be NULL;

The boolean variable `new_x` will be false if the last call to any of the evaluation methods used the same `x` values. This can be helpful when users have efficient implementations that calculate multiple outputs at once. IPOPT internally caches results from the TNLP and generally, this flag can be ignored.

The index variables `n`, `m`, and `nele_jac` are passed in only to help your debugging. These arguments will have the same values you specified in `get_nlp_info`.

In our example, the Jacobian is actually dense, but we still specify it using the sparse format.

```
bool HS071_NLP::eval_jac_g(Index n, const Number* x, bool new_x,
                          Index m, Index nele_jac, Index* iRow, Index *jCol,
                          Number* values)
{
    if (values == NULL) {
        // return the structure of the Jacobian

        // this particular Jacobian is dense
        iRow[0] = 0; jCol[0] = 0;
        iRow[1] = 0; jCol[1] = 1;
        iRow[2] = 0; jCol[2] = 2;
        iRow[3] = 0; jCol[3] = 3;
        iRow[4] = 1; jCol[4] = 0;
        iRow[5] = 1; jCol[5] = 1;
        iRow[6] = 1; jCol[6] = 2;
        iRow[7] = 1; jCol[7] = 3;
    }
    else {
        // return the values of the Jacobian of the constraints

        values[0] = x[1]*x[2]*x[3]; // 0,0
        values[1] = x[0]*x[2]*x[3]; // 0,1
        values[2] = x[0]*x[1]*x[3]; // 0,2
        values[3] = x[0]*x[1]*x[2]; // 0,3

        values[4] = 2*x[0]; // 1,0
        values[5] = 2*x[1]; // 1,1
        values[6] = 2*x[2]; // 1,2
        values[7] = 2*x[3]; // 1,3
    }
}
```

```

    return true;
}

```

Method `eval_h` with Prototype

```

virtual bool eval_h(Index n, const Number* x, bool new_x,
                    Number obj_factor, Index m, const Number* lambda,
                    bool new_lambda, Index nele_hess, Index* iRow,
                    Index* jCol, Number* values)

```

Return the structure of the Hessian of the Lagrangian or the values of the Hessian of the Lagrangian as calculated by the values in `obj_factor`, `x`, and `lambda`.

- `n`: (in), the number of variables in the problem (dimension of `x`).
- `x`: (in), the current values for the primal variables, `x`.
- `new_x`: (in), false if any evaluation method was previously called with the same values in `x`, true otherwise.
- `obj_factor`: (in), factor in front of the objective term in the Hessian.
- `m`: (in), the number of constraints in the problem (dimension of `g`).
- `lambda`: (in), the current values of the equality multipliers to use for each constraint in the evaluation of the Hessian.
- `n_elel_jac`: (in), the number of nonzero elements in the Jacobian (dimension of `iRow`, `jCol`, and `values`).
- `new_lambda`: (in), false if any evaluation method was previously called with the same values in `lambda`, true otherwise.
- `nele_hess`: (in), the number of nonzero elements in the Hessian (dimension of `iRow`, `jCol`, and `values`).
- `iRow`: (out), the row indices of entries in the Hessian.
- `jCol`: (out), the column indices of entries in the Hessian.
- `values`: (out), the values of the entries in the Hessian.

If the `iRow` and `jCol` arguments are not NULL, then IPOPT wants you to fill in the structure of the Hessian (the row and column indices only). In this case, the `x`, `lambda`, and `values` arrays will be NULL.

If the `x`, `lambda`, and `values` arrays are not NULL, then IPOPT wants you to fill in the values of the Hessian as calculated using `x` and `lambda` (using the same order as you used when specifying the structure). In this case, the `iRow` and `jCol` arguments will be NULL.

The boolean variables `new_x` and `new_lambda` will both be false if the last call to any of the evaluation methods used the same values. This can be helpful when users have efficient implementations that calculate multiple outputs at once. IPOPT internally caches results from the TNLP and generally, this flag can be ignored.

The index variables `n`, `m`, and `nele_hess` are passed in only to help your debugging. These arguments will have the same values you specified in `get_nlp_info`.

In our example, the Hessian is dense, but we still specify it using the sparse matrix format. Because the Hessian is symmetric, we only need to specify the lower left corner.

```
bool HS071_NLP::eval_h(Index n, const Number* x, bool new_x,
                      Number obj_factor, Index m, const Number* lambda,
                      bool new_lambda, Index nele_hess, Index* iRow,
                      Index* jCol, Number* values)
{
    if (values == NULL) {
        // return the structure. This is a symmetric matrix, fill the lower left
        // triangle only.

        // the Hessian for this problem is actually dense
        Index idx=0;
        for (Index row = 0; row < 4; row++) {
            for (Index col = 0; col <= row; col++) {
                iRow[idx] = row;
                jCol[idx] = col;
                idx++;
            }
        }

        assert(idx == nele_hess);
    }
    else {
        // return the values. This is a symmetric matrix, fill the lower left
        // triangle only

        // fill the objective portion
        values[0] = obj_factor * (2*x[3]); // 0,0

        values[1] = obj_factor * (x[3]); // 1,0
        values[2] = 0; // 1,1

        values[3] = obj_factor * (x[3]); // 2,0
        values[4] = 0; // 2,1
        values[5] = 0; // 2,2

        values[6] = obj_factor * (2*x[0] + x[1] + x[2]); // 3,0
        values[7] = obj_factor * (x[0]); // 3,1
        values[8] = obj_factor * (x[0]); // 3,2
        values[9] = 0; // 3,3

        // add the portion for the first constraint
        values[1] += lambda[0] * (x[2] * x[3]); // 1,0

        values[3] += lambda[0] * (x[1] * x[3]); // 2,0
        values[4] += lambda[0] * (x[0] * x[3]); // 2,1

        values[6] += lambda[0] * (x[1] * x[2]); // 3,0
        values[7] += lambda[0] * (x[0] * x[2]); // 3,1
        values[8] += lambda[0] * (x[0] * x[1]); // 3,2

        // add the portion for the second constraint
        values[0] += lambda[1] * 2; // 0,0
    }
}
```

```

    values[2] += lambda[1] * 2; // 1,1

    values[5] += lambda[1] * 2; // 2,2

    values[9] += lambda[1] * 2; // 3,3
}

return true;
}

```

Method finalize_solution with Prototype

```

virtual void finalize_solution(SolverReturn status, Index n,
                             const Number* x, const Number* z_L,
                             const Number* z_U, Index m, const Number* g,
                             const Number* lambda, Number obj_value)

```

This is the only method that is not mentioned in Figure 2. This method is called by IPOPT after the algorithm has finished (successfully or even with most errors).

- **status:** (in), gives the status of the algorithm as specified in `IpAlgTypes.hpp`,
 - **SUCCESS:** Algorithm terminated successfully at a locally optimal point.
 - **MAXITER_EXCEEDED:** Maximum number of iterations exceeded (can be specified by an option).
 - **STOP_AT_TINY_STEP:** Algorithm proceeds with very little progress.
 - **STOP_AT_ACCEPTABLE_POINT:** Algorithm stopped at a point that was converged, not to “desired” tolerances, but to “acceptable” tolerances (see the ??? options).
 - **LOCAL_INFEASIBILITY:** Algorithm converged to a point of local infeasibility. Problem may be infeasible.
 - **RESTORATION_FAILURE:** Restoration phase was called, but failed to find a more feasible point.
 - **INTERNAL_ERROR:** An unknown internal error occurred. Please contact the IPOPT Authors through the mailing list.
- **n:** (in), the number of variables in the problem (dimension of **x**).
- **x:** (in), the current values for the primal variables, **x**.
- **z_L:** (in), the current values for the lower bound multipliers.
- **z_U:** (in), the current values for the upper bound multipliers.
- **m:** (in), the number of constraints in the problem (dimension of **g**).
- **g:** (in), the current value of the constraint residuals.
- **lambda:** (in), the current values of the equality multipliers.
- **obj_value:** (in), the current value of the objective.

This method gives you the return status of the algorithm (`SolverReturn`), and the values of the variables, the objective and constraint residuals when the algorithm exited.

In our example, we will print the values of some of the variables to the screen.

```

void HS071_NLP::finalize_solution(SolverReturn status,
                                Index n, const Number* x, const Number* z_L,
                                const Number* z_U, Index m, const Number* g,
                                const Number* lambda, Number obj_value)
{
    // here is where we would store the solution to variables, or write to a file, etc
    // so we could use the solution.

    // For this example, we write the solution to the console
    printf("\n\nSolution of the primal variables, x\n");
    for (Index i=0; i<n; i++) {
        printf("x[%d] = %e\n", i, x[i]);
    }

    printf("\n\nSolution of the bound multipliers, z_L and z_U\n");
    for (Index i=0; i<n; i++) {
        printf("z_L[%d] = %e\n", i, z_L[i]);
    }
    for (Index i=0; i<n; i++) {
        printf("z_U[%d] = %e\n", i, z_U[i]);
    }

    printf("\n\nObjective value\n");
    printf("f(x*) = %e\n", obj_value);
}

```

This is all that is required for our HS071_NLP class and the coding of the problem representation.

3.3.2 Coding the Executable (main)

Now that we have a problem representation, the HS071_NLP class, we need to code the main function that will call IPOPT and ask IPOPT to find a solution.

Here, we must create an instance of our problem (HS071_NLP), create an instance of the IPOPT solver (IpoptApplication), and ask the solver to find a solution. We always use the SmartPtr template class instead of raw C++ pointers when creating and passing IPOPT objects. To find out more information about smart pointers and the SmartPtr implementation used in IPOPT, see Appendix B.

Create the file MyExample.cpp in the MyExample directory. Include HS071_NLP.hpp and IpIpoptApplication.hpp, tell the compiler to use the Ipopt namespace, and implement the main function.

```

#include "IpIpoptApplication.hpp"
#include "hs071_nlp.hpp"

using namespace Ipopt;

int main(int argv, char* argc[])
{
    // Create a new instance of your nlp
    // (use a SmartPtr, not raw)
    SmartPtr<TNLP> mynlp = new HS071_NLP();

    // Create a new instance of IpoptApplication
    // (use a SmartPtr, not raw)
    SmartPtr<IpoptApplication> app = new IpoptApplication();

    // Change some options
    app->Options()->SetNumericValue("tol", 1e-9);
    app->Options()->SetStringValue("mu_strategy", "adaptive");
}

```

```

// Ask Ipopt to solve the problem
ApplicationReturnStatus status = app->OptimizeTNLP(mynlp);

if (status == Solve_Succeeded) {
    printf("\n\n*** The problem solved!\n");
}
else {
    printf("\n\n*** The problem FAILED!\n");
}

// As the SmartPtrs go out of scope, the reference count
// will be decremented and the objects will automatically
// be deleted.

return (int) status;
}

```

The first line of code in main creates an instance of HS071_NLP. We then create an instance of the IPOPT solver, IpoptApplication. The call to app->OptimizeTNLP(...) will run IPOPT and try to solve the problem. By default, IPOPT will write to its progress to the console, and return the SolverReturn status.

3.3.3 Compiling and Testing the Example

Our next task is to compile and test the code. If you are familiar with the compiler and linker used on your system, you can build the code, including the IPOPT library libipopt.a (and other necessary libraries). If you are using Linux/UNIX, then a sample makefile exists already that was created by configure. Copy Examples/hs071_cpp/Makefile into your MyExample directory. This makefile was created for the hs071_cpp code, but it can be easily modified for your example problem. Edit the file, making the following changes,

- change the EXE variable
EXE = my_example
- change the OBJS variable
OBJS = HS071_NLP.o MyExample.o

and the problem should compile easily with,
\$ make Now run the executable,
\$./my_example and you should see output resembling the following,

```

Total number of variables.....:          4
      variables with only lower bounds:          0
      variables with lower and upper bounds:        4
      variables with only upper bounds:          0
Total number of equality constraints.....:          1
Total number of inequality constraints.....:          1
      inequality constraints with only lower bounds:          1
      inequality constraints with lower and upper bounds:          0
      inequality constraints with only upper bounds:          0

iter    objective    inf_pr  inf_du lg(mu)  ||d||  lg(rg) alpha_du alpha_pr  ls
  0    1.7159878e+01  2.01e-02  5.20e-01  -1.0  0.00e+00   -  0.00e+00  0.00e+00   0 y
  1    1.7146308e+01  1.63e-01  1.47e-01  -1.0  1.15e-01   -  9.86e-01  1.00e+00f  1
  2    1.7065508e+01  3.10e-02  8.47e-02  -1.7  1.99e-01   -  9.54e-01  1.00e+00h  1 Nhj

```



```

3  1.7002626e+01 4.10e-02 4.81e-03 -2.5 5.52e-02 - 1.00e+00 1.00e+00h 1
4  1.7019082e+01 1.20e-03 1.81e-04 -2.5 1.10e-02 - 1.00e+00 1.00e+00h 1
5  1.7014253e+01 1.80e-04 4.87e-05 -3.8 4.86e-03 - 1.00e+00 1.00e+00h 1
6  1.7014020e+01 9.25e-07 2.15e-07 -5.7 2.76e-04 - 1.00e+00 1.00e+00h 1
7  1.7014017e+01 1.01e-10 2.60e-11 -8.6 3.32e-06 - 1.00e+00 1.00e+00h 1

```

Number of Iterations.....: 7

	(scaled)	(unscaled)
Objective.....:	1.7014017145177885e+01	1.7014017145177885e+01
Dual infeasibility.....:	2.5980210027546616e-11	2.5980210027546616e-11
Constraint violation.....:	1.8175683180743363e-11	1.8175683180743363e-11
Complementarity.....:	2.5282956951655172e-09	2.5282956951655172e-09
Overall NLP error.....:	2.5282956951655172e-09	2.5282956951655172e-09

```

Number of objective function evaluations      = 8
Number of objective gradient evaluations     = 8
Number of equality constraint evaluations     = 8
Number of inequality constraint evaluations   = 8
Number of equality constraint Jacobian evaluations = 8
Number of inequality constraint Jacobian evaluations = 8
Number of Lagrangian Hessian evaluations    = 9

```

EXIT: Optimal Solution Found.

Solution of the primal variables, x

```

x[0] = 1
x[1] = 4.743
x[2] = 3.82115
x[3] = 1.37941

```

Solution of the bound multipliers, z_L and z_U

```

z_L[0] = 1.08787
z_L[1] = 6.69317e-10
z_L[2] = 8.8877e-10
z_L[3] = 6.57011e-09
z_U[0] = 6.26262e-10
z_U[1] = 9.78906e-09
z_U[2] = 2.12283e-09
z_U[3] = 6.92528e-10

```

Objective value
f(x*) = 17.014

*** The problem solved!

This completes the basic C++ tutorial, but see Section 5 which explains the standard console output of IPOPT and Section 4 for information about the use of options to customize the behavior of IPOPT.

3.4 The C Interface

The C interface for IPOPT is declared in the header file `IpStdCInterface.h`, which is found in `$IPOPTDIR/include/ipopt` (or in `$PREFIX/include/ipopt` if the switch `--prefix=$PREFIX` was used

for `configure`); while reading this section, it will be helpful to have a look at this file.

In order to solve an optimization problem with the C interface, one has to create an `IpoptProblem`⁶ with the function `CreateIpoptProblem`, which later on has to be passed to the `IpoptSolve` function.

The `IpoptProblem` created by `CreateIpoptProblem` contains the problem dimensions, the variable and constraint bounds, and the function pointers for callbacks that will be used to evaluate the NLP problem functions and their derivatives (see also the discussion of the C++ methods `get_nlp_info` and `get_bounds_info` in Section 3.3.1 for information about the arguments of `CreateIpoptProblem`).

The prototypes for the callback functions, `Eval_F_CB`, `Eval_Grad_F_CB`, etc., are defined in the header file `IpStdCInterface.h`. Their arguments correspond one-to-one to the arguments for the C++ methods discussed in Section 3.3.1; for example, for the meaning of `n`, `x`, `new_x`, `obj_value` in the declaration of `Eval_F_CB` see the discussion of “`eval_f`”. The callback functions should return `TRUE`, unless there was a problem doing the requested function/derivative evaluation at the given point `x` (then it should return `FALSE`).

Note the additional argument of type `UserDataPtr` in the callback functions. This pointer argument is available for you to communicate information between the main program that calls `IpoptSolve` and any of the callback functions. This pointer is simply passed unmodified by IPOPT among those functions. For example, you can use this to pass constants that define the optimization problem and are computed before the optimization in the main C program to the callback functions.

After an `IpoptProblem` has been created, you can set algorithmic options for IPOPT (see Section 4) using the `AddIpopt*Option` functions. Finally, the IPOPT algorithm is called with `IpoptSolve`, giving IPOPT the `IpoptProblem`, the starting point, and arrays to store the solution values (primal and dual variables), if desired. Finally, after everything is done, you should call `FreeIpoptProblem` to release internal memory that is still allocated inside IPOPT.

In the remainder of this section we discuss how the example problem (4)–(7) can be solved using the C interface. A completed version of this example can be found in `Examples/hs071.c`.

In order to implement the example problem on your own, create a new directory `MyCExample` and create a new file, `hs071.c`. Here, include the interface header file `IpStdCInterface.h`, along with other necessary header files, such as `stdlib.h` and `assert.h`. Add the prototypes and implementations for the five callback functions. Have a look at the C++ implementation for `eval_f`, `eval_g`, `eval_grad_f`, `eval_jac_g`, and `eval_h` in Section 3.3.1. The C implementations have somewhat different prototypes, but are implemented almost identically to the C++ code. See the completed example in `Examples/hs071_c/hs071.c` if you are not sure how to do this.

We now need to implement the main function, create the `IpoptProblem`, set options, and call `IpoptSolve`. The `CreateIpoptProblem` function requires the problem dimensions, the variable and constraint bounds, and the function pointers to the callback routines. The `IpoptSolve` function requires the `IpoptProblem`, the starting point, and allocated arrays for the solution. The main function from the example is shown next, and discussed below.

```
int main()
{
    Index n=-1;                /* number of variables */
    Index m=-1;                /* number of constraints */
    Number* x_L = NULL;        /* lower bounds on x */
    Number* x_U = NULL;        /* upper bounds on x */
    Number* g_L = NULL;        /* lower bounds on g */
}
```

⁶`IpoptProblem` is a pointer to a C structure; you should not access this structure directly, only through the functions provided in the C interface.

```

Number* g_U = NULL;           /* upper bounds on g */
IpoptProblem nlp = NULL;      /* IpoptProblem */
enum ApplicationReturnStatus status; /* Solve return code */
Number* x = NULL;             /* starting point and solution vector */
Number* mult_x_L = NULL;      /* lower bound multipliers
at the solution */
Number* mult_x_U = NULL;      /* upper bound multipliers
at the solution */
Number obj;                   /* objective value */
Index i;                      /* generic counter */

/* set the number of variables and allocate space for the bounds */
n=4;
x_L = (Number*)malloc(sizeof(Number)*n);
x_U = (Number*)malloc(sizeof(Number)*n);
/* set the values for the variable bounds */
for (i=0; i<n; i++) {
    x_L[i] = 1.0;
    x_U[i] = 5.0;
}

/* set the number of constraints and allocate space for the bounds */
m=2;
g_L = (Number*)malloc(sizeof(Number)*m);
g_U = (Number*)malloc(sizeof(Number)*m);
/* set the values of the constraint bounds */
g_L[0] = 25; g_U[0] = 2e19;
g_L[1] = 40; g_U[1] = 40;

/* create the IpoptProblem */
nlp = CreateIpoptProblem(n, x_L, x_U, m, g_L, g_U, 8, 10, 0,
    &eval_f, &eval_g, &eval_grad_f,
    &eval_jac_g, &eval_h);

/* We can free the memory now - the values for the bounds have been
copied internally in CreateIpoptProblem */
free(x_L);
free(x_U);
free(g_L);
free(g_U);

/* set some options */
AddIpoptNumOption(nlp, "tol", 1e-9);
AddIpoptStrOption(nlp, "mu_strategy", "adaptive");

/* allocate space for the initial point and set the values */
x = (Number*)malloc(sizeof(Number)*n);
x[0] = 1.0;
x[1] = 5.0;
x[2] = 5.0;
x[3] = 1.0;

/* allocate space to store the bound multipliers at the solution */
mult_x_L = (Number*)malloc(sizeof(Number)*n);
mult_x_U = (Number*)malloc(sizeof(Number)*n);

/* solve the problem */
status = IpoptSolve(nlp, x, NULL, &obj, NULL, mult_x_L, mult_x_U, NULL);

```

```

if (status == Solve_Succeeded) {
    printf("\n\nSolution of the primal variables, x\n");
    for (i=0; i<n; i++) {
        printf("x[%d] = %e\n", i, x[i]);
    }

    printf("\n\nSolution of the bound multipliers, z_L and z_U\n");
    for (i=0; i<n; i++) {
        printf("z_L[%d] = %e\n", i, mult_x_L[i]);
    }
    for (i=0; i<n; i++) {
        printf("z_U[%d] = %e\n", i, mult_x_U[i]);
    }

    printf("\n\nObjective value\n");
    printf("f(x*) = %e\n", obj);
}

/* free allocated memory */
FreeIpoptProblem(nlp);
free(x);
free(mult_x_L);
free(mult_x_U);

return 0;
}

```

Here, we declare all the necessary variables and set the dimensions of the problem. The problem has 4 variables, so we set `n` and allocate space for the variable bounds (don't forget to call `free` for each of your `malloc` calls before the end of the program). We then set the values for the variable bounds.

The problem has 2 constraints, so we set `m` and allocate space for the constraint bounds. The first constraint has a lower bound of 25 and no upper bound. Here we set the upper bound to `2e19`. IPOPT interprets any number greater than `nlp_upper_bound_inf` as infinity. The default value of `nlp_lower_bound_inf` and `nlp_upper_bound_inf` is `-1e19` and `1e19`, respectively, and can be changed through IPOPT options. The second constraint is an equality, so we set both the upper and the lower bound to 40.

We next create an instance of the `IpoptProblem` by calling `CreateIpoptProblem`, giving it the problem dimensions and the variable and constraint bounds. The arguments `nele_jac` and `nele_hess` are the number of elements in Jacobian and the Hessian, respectively. See Appendix A for a description of the sparse matrix format. The `index_style` argument specifies whether we want to use C style indexing for the row and column indices of the matrices or Fortran style indexing. Here, we set it to 0 to indicate C style. We also include the references to each of our callback functions. IPOPT uses these function pointers to ask for evaluation of the NLP when required.

After freeing the bound arrays that are no longer required, the next two lines illustrate how you can change the value of options through the interface. IPOPT options can also be changed by creating a `PARAMS.DAT` file (see Section 4). We next allocate space for the initial point and set the values as given in the problem definition.

The call to `IpoptSolve` can provide us with information about the solution, but most of this is optional. Here, we want values for the bound multipliers at the solution and we allocate space for these.

We can now make the call to `IpoptSolve` and find the solution of the problem. We pass in the `IpoptProblem`, the starting point `x` (IPOPT will use this array to return the solution as well). The next 5 arguments are pointers so IPOPT can fill in values at the solution. If these pointers are set to `NULL`,

IPOPT will ignore that entry. For example, here, we do not want the constraint residuals at the solution or the equality multipliers, so we set those entries to NULL. We do want the value of the objective, and the multipliers for the variable bounds. The last argument is a `void*` for user data. Any pointer you give here will also be passed to you in the callback functions.

The return code is an `ApplicationReturnStatus` enumeration, see the header file `ReturnCodes_inc.h` which is installed along `IpStdCInterface.h` in the IPOPT include directory.

After the problem has solved, we check the status and print the solution if successful. Finally, we free the `IpoptProblem` and the remaining memory, and return from `main`.

3.5 The Fortran Interface

The Fortran interface is essentially a wrapper of the C interface discussed in Section 3.4. The way to hook up IPOPT in a Fortran program is very similar to how it is done for the C interface, and the functions of the Fortran interface correspond one-to-one to the those of the C interface, including their arguments. You can find an implementation of the example problem (4)–(7) in `$IPOPTDIR/Examples/hs071.f`.

The only special things to consider are:

- The return value of the function `IPCREATE` is of an `INTEGER` type that must be large enough to capture a pointer on the particular machine. This means, that you have to declare the “handle” for the `IpoptProblem` as `INTEGER*8` if your program is compiled in 64-bit mode. All other `INTEGER`-type variables must be of the regular type.
- For the call of `IPSOLVE` (which is the function that is to be called to run IPOPT), all arrays, including those for the dual variables, must be given (in contrast to the C interface). The return value `IERR` of this function indicates the outcome of the optimization (see the include file `IpReturnCodes.inc` in the IPOPT include directory).
- The return `IERR` value of the remaining functions is zero, unless there was a problem during execution of the function call.
- The callback functions (`EV_*` in the example) include the arguments `IDAT` and `DAT`, which are `INTEGER` and `DOUBLE PRECISION` arrays that are passed unmodified between the main program calling `IPSOLVE` and the evaluation subroutines `EV_*` (similarly to `UserDataPtr` arguments in the C interface). Those can be used to pass “private” data between the user-provided Fortran subroutines.

The last argument of the `EV_*` subroutines, `IERR`, is to be set to 0 by the users on return, unless there was a problem during the evaluation of the optimization problem function/derivative for the given point `X` (then it should return a non-zero value).

4 Ipopt Options

Ipopt has many (maybe too many) options that can be adjusted for the algorithm. Options are all identified by a string name and their values can be of one of three types, Number (real), Integer, or String. Number options are used for things like tolerances, integer options are used for things like maximum number of iterations, and string options are used for setting algorithm details, like the NLP scaling method. Options can be set through code, through the AMPL interface if you are using AMPL, or by creating a `PARAMS.DAT` file in the directory you are executing Ipopt.

The `PARAMS.DAT` file is read line by line and each line should contain the option name, followed by whitespace, and then the value. Comments can be included with the `#` symbol. Don't forget to ensure you have a newline at the end of the file. For example,

```
# This is a comment

# Turn off the NLP scaling
nlp_scaling_method none

# Change the initial barrier parameter
mu_init 1e-2

# Set the max number of iterations
max_iter 500
```

is a valid `PARAMS.DAT` file.

Options can also be set in code. Have a look at the examples to see how this is done.

A subset of IPOPT options are available through AMPL. To set options through AMPL, use the internal AMPL command `options`. For example,

```
options ipopt 'nlp_scaling_method=none mu_init=1e-2 max_iter=500'
```

is a valid options command in AMPL. The most common options are referenced in Appendix C. These are also the options that are available through AMPL using the `options` command. To specify other options when using AMPL, you can always create `PARAMS.DAT`. Note, the `PARAMS.DAT` file is given preference when setting options. This way, you can easily override any options set in a particular executable or AMPL model by specifying new values in `PARAMS.DAT`.

For a short list of the valid options, see the Appendix C. You can print the documentation for all Ipopt options by adding the option,

```
print_options_documentation yes
```

and running Ipopt (like the `Ampl` executable, for instance). This will output all of the options documentation to the console.

5 Ipopt Output

This section describes the standard Ipopt console output with the default setting for `print_level`. The output is designed to provide a quick summary of each iteration as Ipopt solves the problem.

Before Ipopt starts to solve the problem, it displays the problem statistics (number of variables, etc.). Note that if you have fixed variables (both upper and lower bounds are equal), Ipopt may remove these variables from the problem.

Following the problem statistics, Ipopt will begin to solve the problem and you will see output resembling the following,

iter	objective	inf_pr	inf_du	lg(mu)	d	lg(rg)	alpha_du	alpha_pr	ls
0	1.6109693e+01	1.12e+01	5.28e-01	0.0	0.00e+00	-	0.00e+00	0.00e+00	0
1	1.8029749e+01	9.90e-01	6.62e+01	0.1	2.05e+00	-	2.14e-01	1.00e+00f	1
2	1.8719906e+01	1.25e-02	9.04e+00	-2.2	5.94e-02	2.0	8.04e-01	1.00e+00h	1

and the columns of output are defined as,

iter The current iteration count. This includes regular iterations and iterations while in restoration phase. If in restoration phase, the letter **r**' will be appended to the iteration number.

objective The unscaled objective value at the current point. In restoration, this value remains the unscaled objective value for the original problem.

inf_pr The scaled primal infeasibility at the current point. In restoration, this value is the primal infeasibility of the original problem at the current point.

inf_du The scaled dual infeasibility at the current point. In restoration, this is the value of the dual infeasibility for the restoration problem.

lg(mu) The value of the barrier parameter mu (\log_{10}).

||d|| The infinity norm (max) of the primal step (in **x** and **s**). In restoration, this value includes the values of the additional variables, *p* and *n*.

lg(rg) The value of the regularization term for the Hessian of the Lagrangian (\log_{10}).

alpha_du The stepsize for the dual variables.

alpha_pr The stepsize for the primal variables.

ls The number of backtracking line search steps.

A Triplet Format for Sparse Matrices

IPOPT was designed for optimizing large sparse nonlinear programs. Because of problem sparsity, the required matrices (like the Jacobian or Hessian) are not stored as traditional dense matrices, but rather in a sparse matrix format. For the tutorials in this document, we use the triplet format. Consider the matrix,

$$\begin{bmatrix} 1.1 & 0 & 0 & 0 & 0 & 0 & 0.5 \\ 0 & 1.9 & 0 & 0 & 0 & 0 & 0.5 \\ 0 & 0 & 2.6 & 0 & 0 & 0 & 0.5 \\ 0 & 0 & 7.8 & 0.6 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1.5 & 2.7 & 0 & 0 \\ 1.6 & 0 & 0 & 0 & 0.4 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0.9 & 1.7 \end{bmatrix} \quad (12)$$

A standard dense matrix representation would need to store $7 \cdot 7 = 49$ floating point numbers, where many entries would be zero. In triplet format, however, only the nonzero entries are stored. The triplet format records the row number, the column number, and the value of all nonzero entries in the matrix. For the matrix above, this means storing 14 integers for the rows, 14 integers for the columns, and 14 floating point numbers for the values. While this does not seem like a huge space savings over the 49 floating point numbers stored in the dense representation, for larger matrices, the space savings are very dramatic⁷

The option `index_style` in `get_nlp_info` tells Ipopt if you prefer to use C style indexing (0-based) for the row and column indices or Fortran style (1-based). Tables 1 and 2 below show the triplet format for both indexing styles, using the example matrix (12).

Table 1: Triplet Format of Matrix (12) with `index_style=FORTRAN_STYLE`

row	col	value
iRow[0] = 1	jCol[0] = 1	values[0] = 1.1
iRow[1] = 1	jCol[1] = 7	values[1] = 0.5
iRow[2] = 2	jCol[2] = 2	values[2] = 1.9
iRow[3] = 2	jCol[3] = 7	values[3] = 0.5
iRow[4] = 3	jCol[4] = 3	values[4] = 2.6
iRow[5] = 3	jCol[5] = 7	values[5] = 0.5
iRow[6] = 4	jCol[6] = 3	values[6] = 7.8
iRow[7] = 4	jCol[7] = 4	values[7] = 0.6
iRow[8] = 5	jCol[8] = 4	values[8] = 1.5
iRow[9] = 5	jCol[9] = 5	values[9] = 2.7
iRow[10] = 6	jCol[10] = 1	values[10] = 1.6
iRow[11] = 6	jCol[11] = 5	values[11] = 0.4
iRow[12] = 7	jCol[12] = 6	values[12] = 0.9
iRow[13] = 7	jCol[13] = 7	values[13] = 1.7

The individual elements of the matrix can be listed in any order, and if there are multiple items for the same nonzero position, the values provided for those positions are added.

⁷For an $n \times n$ matrix, the dense representation grows with the the square of n , while the sparse representation grows linearly in the number of nonzeros.

Table 2: Triplet Format of Matrix (12) with `index_style=C_STYLE`

row	col	value
iRow[0] = 0	jCol[0] = 0	values[0] = 1.1
iRow[1] = 0	jCol[1] = 6	values[1] = 0.5
iRow[2] = 1	jCol[2] = 1	values[2] = 1.9
iRow[3] = 1	jCol[3] = 6	values[3] = 0.5
iRow[4] = 2	jCol[4] = 2	values[4] = 2.6
iRow[5] = 2	jCol[5] = 6	values[5] = 0.5
iRow[6] = 3	jCol[6] = 2	values[6] = 7.8
iRow[7] = 3	jCol[7] = 3	values[7] = 0.6
iRow[8] = 4	jCol[8] = 3	values[8] = 1.5
iRow[9] = 4	jCol[9] = 4	values[9] = 2.7
iRow[10] = 5	jCol[10] = 0	values[10] = 1.6
iRow[11] = 5	jCol[11] = 4	values[11] = 0.4
iRow[12] = 6	jCol[12] = 5	values[12] = 0.9
iRow[13] = 6	jCol[13] = 6	values[13] = 1.7

The Hessian of the Lagrangian is a symmetric matrix. In the case of a symmetric matrix, you only need to specify the lower left corner. For example, given the matrix,

$$\begin{bmatrix} 1.0 & 0 & 3.0 & 0 & 2.0 \\ 0 & 1.1 & 0 & 0 & 5.0 \\ 3.0 & 0 & 1.2 & 6.0 & 0 \\ 0 & 0 & 6.0 & 1.3 & 9.0 \\ 2.0 & 5.0 & 0 & 9.0 & 1.4 \end{bmatrix} \quad (13)$$

the triplet format is shown in Tables 3 and 4.

Table 3: Triplet Format of Matrix (12) with `index_style=FORTRAN_STYLE`

row	col	value
iRow[0] = 1	jCol[0] = 1	values[0] = 1.0
iRow[1] = 2	jCol[1] = 1	values[1] = 1.1
iRow[2] = 3	jCol[2] = 1	values[2] = 3.0
iRow[3] = 3	jCol[3] = 3	values[3] = 1.2
iRow[4] = 4	jCol[4] = 3	values[4] = 6.0
iRow[5] = 4	jCol[5] = 4	values[5] = 1.3
iRow[6] = 5	jCol[6] = 1	values[6] = 2.0
iRow[7] = 5	jCol[7] = 2	values[7] = 5.0
iRow[8] = 5	jCol[8] = 4	values[8] = 9.0
iRow[9] = 5	jCol[9] = 5	values[9] = 1.4

Table 4: Triplet Format of Matrix (12) with `index_style=C_STYLE`

row	col	value
iRow[0] = 0	jCol[0] = 0	values[0] = 1.0
iRow[1] = 1	jCol[1] = 0	values[1] = 1.1
iRow[2] = 2	jCol[2] = 0	values[2] = 3.0
iRow[3] = 2	jCol[3] = 2	values[3] = 1.2
iRow[4] = 3	jCol[4] = 2	values[4] = 6.0
iRow[5] = 3	jCol[5] = 3	values[5] = 1.3
iRow[6] = 4	jCol[6] = 0	values[6] = 2.0
iRow[7] = 4	jCol[7] = 1	values[7] = 5.0
iRow[8] = 4	jCol[8] = 3	values[8] = 9.0
iRow[9] = 4	jCol[9] = 4	values[9] = 1.4

B The Smart Pointer Implementation: SmartPtr

The SmartPtr class is described in `IpSmartPtr.hpp`. It is a template class that takes care of deleting objects for us so we need not be concerned about memory leaks. Instead of pointing to an object with a raw C++ pointer (e.g. `HS071_NLP*`), we use a SmartPtr. Every time a SmartPtr is set to reference an object, it increments a counter in that object (see the ReferencedObject base class if you are interested). If a SmartPtr is done with the object, either by leaving scope or being set to point to another object, the counter is decremented. When the count of the object goes to zero, the object is automatically deleted. SmartPtr's are very simple, just use them as you would a standard pointer.

It is very important to use SmartPtr's instead of raw pointers when passing objects to IPOPT. Internally, IPOPT uses smart pointers for referencing objects. If you use a raw pointer in your executable, the object's counter will NOT get incremented. Then, when IPOPT uses smart pointers inside its own code, the counter will get incremented. However, before IPOPT returns control to your code, it will decrement as many times as it incremented and the counter will return to zero. Therefore, IPOPT will delete the object. When control returns to you, you now have a raw pointer that points to a deleted object.

This might sound difficult to anyone not familiar with the use of smart pointers, but just follow one simple rule; always use a SmartPtr when creating or passing an IPOPT object.

C Options Reference

Options can be set using `PARAMS.DAT`, through your own code, or through the AMPL `options` command. See Section 4 for an explanation of how to use these commands. Shown here is a short list of the most common options for Ipopt. To view the full list of options, run the `ipopt` executable with the option,

```
print_options_documentation yes
```

The most common options are:

print_level: Output verbosity level.

Sets the default verbosity level for console output. The larger this value the more detailed is the output. The valid range for this integer option is $0 \leq \text{print_level} \leq 10$ and its default value is 3.

pivotol: Pivot tolerance for the linear solver.

A smaller number pivots for sparsity, a larger number pivots for stability. The valid range for this real option is $0 < \text{pivotol} < 1$ and its default value is $1 \cdot 10^{-08}$.

pivotolmax: Maximum pivot tolerance.

IPOPT may increase pivotol as high as pivotolmax to get a more accurate solution to the linear system. The valid range for this real option is $0 < \text{pivotolmax} < 1$ and its default value is 0.0001.

tol: Desired convergence tolerance (relative).

Determines the convergence tolerance for the algorithm. The algorithm terminates successfully, if the (scaled) NLP error becomes smaller than this value, and if the (absolute) criteria according to "dual_inf_tol", "primal_inf_tol", and "compl_inf_tol" are met. (This is `epsilon_tol` in Eqn. (6) in implementation paper). Note, some other algorithmic features also use this quantity. The valid range for this real option is $0 < \text{tol} < +\text{inf}$ and its default value is $1 \cdot 10^{-08}$.

compl_inf_tol: Desired threshold for the complementarity conditions.

Absolute tolerance on the complementarity. Successful termination requires that the (unscaled) complementarity is less than this threshold. The valid range for this real option is $0 < \text{compl_inf_tol} < +\text{inf}$ and its default value is 0.0001.

dual_inf_tol: Desired threshold for the dual infeasibility.

Absolute tolerance on the dual infeasibility. Successful termination requires that the (unscaled) dual infeasibility is less than this threshold. The valid range for this real option is $0 < \text{dual_inf_tol} < +\text{inf}$ and its default value is 0.0001.

constr_mult_init_max: Maximum allowed least-square guess of constraint multipliers.

Determines how large the initial least-square guesses of the constraint multipliers are allowed to be (in max-norm). If the guess is larger than this value, it is discarded and all constraint multipliers are set to zero. This options is also used when initializing the restoration phase. By default, "resto.constr_mult_init_max" (the one used in `RestoIterateInitializer`) is set to zero. The valid range for this real option is $0 \leq \text{constr_mult_init_max} < +\text{inf}$ and its default value is 1000.

constr_viol_tol: Desired threshold for the constraint violation.

Absolute tolerance on the constraint violation. Successful termination requires that the (unscaled) constraint violation is less than this threshold. The valid range for this real option is $0 < \text{constr_viol_tol} < +\text{inf}$ and its default value is 0.0001.

mu_strategy: Update strategy for barrier parameter.

Determines which barrier parameter update strategy is to be used.

Possible values:

- **monotone:** use the monotone (Fiacco-McCormick) strategy
- **adaptive:** use the adaptive update strategy

mu_init: Initial value for the barrier parameter.

This option determines the initial value for the barrier parameter (μ). It is only relevant in the monotone, Fiacco-McCormick version of the algorithm. (i.e., if "mu_strategy" is chosen as "monotone") The valid range for this real option is $0 < \text{mu_init} < +\text{inf}$ and its default value is 0.1.

mu_oracle: Oracle for a new barrier parameter in the adaptive strategy

Determines how a new barrier parameter is computed in each "free-mode" iteration of the adaptive barrier parameter strategy. (Only considered if "adaptive" is selected for option "mu_strategy").

Possible values:

- **probing:** Mehrotra's probing heuristic
- **loqo:** LOQO's centrality rule
- **quality_function:** minimize a quality function

corrector_type: The type of corrector steps that should be taken.

If "mu_strategy" is "adaptive", this option determines what kind of corrector steps should be tried.

Possible values:

- **none:** no corrector
- **affine:** corrector step towards $\mu=0$
- **primal-dual:** corrector step towards current μ

obj_scaling_factor: Scaling factor for the objective function.

This option sets a scaling factor for the objective function. The scaling is seen internally by Ipopt but the unscaled objective is reported in the console output. If additional scaling parameters are computed (e.g. user-scaling or gradient-based), both factors are multiplied. If this value is chosen to be negative, Ipopt will maximize the objective function instead of minimizing it. The valid range for this real option is $-\text{inf} < \text{obj_scaling_factor} < +\text{inf}$ and its default value is 1.

nlp_scaling_method: Select the technique used for scaling the NLP

Selects the technique used for scaling the problem before it is solved. For user-scaling, the parameters come from the NLP. If you are using AMPL, they can be specified through suffixes (`scaling_factor`)

Possible values:

- `none`: no problem scaling will be performed
- `user_scaling`: scaling parameters will come from the user
- `gradient_based`: scale the problem so the maximum gradient at the starting point is `scaling_max_gradient`

nlp_scaling_max_gradient: Maximum gradient after scaling.

This is the gradient scaling cut-off. If the maximum gradient is above this value, then gradient based scaling will be performed. Scaling parameters are calculated to scale the maximum gradient back to this value. Note: This option is only used if "nlp_scaling_method" is chosen as "gradient_based". The valid range for this real option is $0 < \text{nlp_scaling_max_gradient} < +\text{inf}$ and its default value is 100.

bound_frac: Desired minimum relative distance from the initial point to bound

Determines how much the initial point might have to be modified in order to be sufficiently inside the bounds (together with "bound_push"). The valid range for this real option is $0 < \text{bound_frac} \leq 0.5$ and its default value is 0.01.

bound_mult_init_val: Initial value for the bound multipliers

All dual variables corresponding to bound constraints are initialized to this value. The valid range for this real option is $0 < \text{bound_mult_init_val} < +\text{inf}$ and its default value is 1.

bound_push: Desired minimum absolute distance from the initial point to bound

Determines how much the initial point might have to be modified in order to be sufficiently inside the bounds (together with "bound_frac"). The valid range for this real option is $0 < \text{bound_push} < +\text{inf}$ and its default value is 0.01.

bound_relax_factor: Factor for initial relaxation of the bounds.

Before start of the optimization, the bounds given by the user are relaxed. This option sets the factor for this relaxation. If it is set to zero, then then bounds relaxation is disabled. (See Eqn.(35) in implmentation paper.) The valid range for this real option is $0 \leq \text{bound_relax_factor} < +\text{inf}$ and its default value is $1 \cdot 10^{-08}$.

acceptable_compl_inf_tol: Acceptance threshold for the complementarity conditions.

Absolute tolerance on the complementarity. Acceptable termination requires that the (unscaled) complementarity is less than this threshold. The valid range for this real option is $0 < \text{acceptable_compl_inf_tol} < +\text{inf}$ and its default value is 0.01.

acceptable_constr_viol_tol: Acceptance threshold for the constraint violation.

Absolute tolerance on the constraint violation. Acceptable termination requires that the (unscaled) constraint violation is less than this threshold. The valid range for this real option is $0 < \text{acceptable_constr_viol_tol} < +\text{inf}$ and its default value is 0.01.

acceptable_dual_inf_tol: Acceptance threshold for the dual infeasibility.

Absolute tolerance on the dual infeasibility. Acceptable termination requires that the (unscaled) dual infeasibility is less than this threshold. The valid range for this real option is $0 < \text{acceptable_dual_inf_tol} < +\text{inf}$ and its default value is 0.01.

acceptable_tol: Acceptable convergence tolerance (relative).

Determines which (scaled) overall optimality error is considered to be acceptable. If the algorithm encounters "acceptable_iter" iterations in a row that are considered "acceptable", it will terminate before the desired convergence tolerance ("tol", "dual_inf_tol", etc) is met. The valid range for this real option is $0 < \text{acceptable_tol} < +\text{inf}$ and its default value is $1 \cdot 10^{-06}$.

alpha_for_y: Method to determine the step size for constraint multipliers.

This option determines how the step size (alpha_y) will be calculated when updating the constraint multipliers.

Possible values:

- primal: use primal step size
- bound_mult: use step size for the bound multipliers
- min: use the min of primal and bound multipliers
- max: use the max of primal and bound multipliers
- full: take a full step of size one
- min_dual_infeas: choose step size minimizing new dual infeasibility
- safe_min_dual_infeas: like "min_dual_infeas", but safeguarded by "min" and "max"

expect_infeasible_problem: Enable heuristics to quickly detect an infeasible problem.

This options is meant to activate heuristics that may speed up the infeasibility determination if you expect the problem to be infeasible. In the filter line search procedure, the restoration phase is called more quickly than usually, and more reduction in the constraint violation is enforced. If the problem is square, this is enabled automatically.

Possible values:

- no: the problem probably be feasible
- yes: the problem has a good chance to be infeasible

max_iter: Maximum number of iterations.

The algorithm terminates with an error message if the number of iterations exceeded this number. This option is also used in the restoration phase. The valid range for this integer option is $0 \leq \text{max_iter} < +\text{inf}$ and its default value is 3000.

max_refinement_steps: Maximum number of iterative refinement steps per linear system solve.

Iterative refinement (on the full unsymmetric system) is performed for each right hand side. This option determines the maximum number of iterative refinement steps. The valid range for this integer option is $0 \leq \text{max_refinement_steps} < +\text{inf}$ and its default value is 10.

max_soc: Maximum number of second order correction trial steps at each iteration.

Choosing 0 disables the second order corrections. (This is p_{max} of Step A-5.9 of Algorithm A in implementation paper.) The valid range for this integer option is $0 \leq \text{max_soc} < +\text{inf}$ and its default value is 4.

min_refinement_steps: Minimum number of iterative refinement steps per linear system solve.

Iterative refinement (on the full unsymmetric system) is performed for each right hand side. This option determines the minimum number of iterative refinements (i.e. at least "min_refinement_steps" iterative refinement steps are enforced per right hand side.) The valid range for this integer option is $0 \leq \text{min_refinement_steps} < +\text{inf}$ and its default value is 1.

output_file: File name of desired output file (leave unset for no file output)

NOTE: This option only works when read from PARAMS.DAT options file! An output file with this name will be written (leave unset for no file output). The verbosity level is by default, "print_level", or but can be overridden with "file_print_level".

Possible values:

- *: Any acceptable standard file name

file_print_level: Verbosity level for output file.

NOTE: This option only works when read from PARAMS.DAT options file! Determines the verbosity level for the file specified by "output_file". By default it is the same as "print_level". The valid range for this integer option is $0 \leq \text{file_print_level} \leq 10$ and its default value is 3.

D Detailed Installation Information