



Usermanual

Version: December 9, 2022

IPBASIC

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To get more information about the commercial licence and the commercial use of IPBASIC and OCPBASIC please contact info@ocpbasic.com.

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2 Installing IPbasic and OCPbasic

2.1 Fast Installing Procedure

If you do not like to build IPBASIC and OCPBASIC yourself you can get precompiled packages for Ubuntu 22.04.

2.1.1 Installing via apt-get on Ubuntu 22.04

This section will install IPBASIC, OCPBASIC and LAPACK_WRAPPER using an apt-repository and the linear algebra packages libblas-dev and liblapack-dev.

First download and activate the ocpbasic gpg key and install it to your system.

Comment 2.1 *Only add software repositories from sources that you trust! Third-party software repositories are not checked for security or reliability and may contain software which is harmful to your computer. The use of these repositorys is at your own risk!*

```
$ wget https://archive.ocpbasic.com/dists/jammy/ocpbasic.gpg
$ sudo install -o root -g root -m 644 ocpbasic.gpg \
/etc/apt/trusted.gpg.d/
$ sudo rm ocpbasic.gpg
```

After that you can activate the ocpbasic apt-repository with the command:

```
$ sudo add-apt-repository \
'deb [arch=amd64] http://archive.ocpbasic.com/dists/jammy /'
```

If the apt-repository is installed correctly you can install OCPBASIC with the command:

```
$ sudo apt-get install ocpbasic
```

The software packages libblas-dev, liblapack-dev, LAPACK_WRAPPER and IPBASIC will then also be installed by apt.

Supported Ubuntu Versions:

Ubuntu version:	apt-repository
Ubuntu 22.04	'deb [arch=amd64] http://archive.ocpbasic.com/dists/jammy /'
Ubuntu 20.04	'deb [arch=amd64] http://archive.ocpbasic.com/dists/focal /'

Ubuntu 18.04	'deb [arch=amd64] http://archive.ocpbasic.com/dists/bionic /'
--------------	---

Table 1: Supported Ubuntu Versions

Comment 2.2 *The maintainer may remove or delete these apt-repositorys any time!*

2.2 Selection of linear algebra software packages

Package:	BLAS	LAPACK
BLAS / LAPACK:	✓	✓
OpenBLAS	✓	✓
Intel Math Kernel Library (MKL)	✓	✓
ATLAS	~	~
Accelerate - Apple Developer	~	~

Table 2: Dense linear algebra software packages used by LAPACK_WRAPPER

Table 2 shows all tested and supported dense linear algebra software packages from LAPACK_WRAPPER, that can be used by IPBASIC and OCPBASIC.

- ✓ Supported by LAPACK_WRAPPER and tested with IPBASIC and OCPBASIC.
- ~ Supported by LAPACK_WRAPPER but not tested with IPBASIC and OCPBASIC.

2.2.1 Selection of linear algebra software packages using qmake

Comment 2.3 *The user is responsible for linking against any thirdparty software and checking the licence compatibility.*

When using the default build method qmake for building LAPACK_WRAPPER, IPBASIC and OCPBASIC you will find a config.pri file in the src/ folder. In this file you will find a section SET DEFINES, where used packages can be selected. Fild the following lines:

```
#  
# Define used linear Algebra.  
#  
# Only supported by lapack_wrapper:
```



```
# LAPACK_WRAPPER_USE_ACCELERATE , LAPACK_WRAPPER_USE_ATLAS ,  
#  
# Currently supported values by IPbasic and OCPbasic:  
# LAPACK_WRAPPER_USE_LAPACK , LAPACK_WRAPPER_USE_OPENBLAS ,  
# LAPACK_WRAPPER_USE_MKL  
#  
DEFINES += LAPACK_WRAPPER_USE_LAPACK
```

You can change `LAPACK_WRAPPER_USE_LAPACK` into your preferred linear algebra software package define.

Comment 2.4 *You must only use one single linear algebra software package define!*

After selecting your linear algebra software package you have to set the include and shared library path and the linked linear algebra package. Therefore you will find a section `SET PATH` in the `config.pri` file.

```
#  
# Define lib and include path for blas and lapack:  
#  
BLAS_INCLUDE_PATH = /usr/include/x86_64-linux-gnu/  
  
BLAS_PATH_LIBS = -L/usr/lib/x86_64-linux-gnu/ -lblas  
  
LAPACK_INCLUDE_PATH = /usr/include/x86_64-linux-gnu/  
  
LAPACK_PATH_LIBS = -L/usr/lib/x86_64-linux-gnu/ -llapack  
  
#  
# Define lib and include path for Openblas:  
#  
OPENBLAS_INCLUDE_PATH = \  
    /usr/include/x86_64-linux-gnu/openblas-pthread/  
  
OPENBLAS_PATH_LIBS = -L/usr/lib/x86_64-linux-gnu/ \  
    -lopenblas  
  
#  
# Define lib and include path for intel mkl:  
#
```

```
MKL_INCLUDE_PATH = /usr/include/mkl/

MKL_PATH_LIBS = -L/usr/lib/x86_64-linux-gnu/ \
               -lmkl_intel_lp64 \
               -lmkl_intel_thread \
               -lmkl_core \
               -lmkl_rt \
               -liomp5
```

Comment 2.5 *Do not delete or comment any unused include and shared library path. When not selected as define the path variables for other linear algebra software packages are ignored anyway.*

If you need to link additional software packages and libraries you can activate the define OCPBASIC_ADD_LINKED_LIBS by uncommenting the following line in the section SET DEFINES.

```
#
# Define use of additional linked libs.
#
DEFINES += OCPBASIC_ADD_LINKED_LIBS
```

After activating the define OCPBASIC_ADD_LINKED_LIBS you can link additional libraries in the variable ADD_LINKED_LIBS in section SET PATH of the config.pri file.

```
#
# Additional linked libs:
#
ADD_LINKED_LIBS = -lgfortran
```

2.2.2 Selection of linear algebra software packages using other buildsystem

Comment 2.6 *The user is responsible for linking against any thirdparty software and checking the licence compatibility.*

The linear algebra software packages can be selected in the configuration file of LAPACK_WRAPPER lapack_wrapper_config.hh. Uncomment the linear algebra software package define.

```
/*\
! values
! LAPACK_WRAPPER_USE_ACCELERATE, LAPACK_WRAPPER_USE_ATLAS,
```

```
! LAPACK_WRAPPER_USE_OPENBLAS , LAPACK_WRAPPER_USE_LAPACK ,
! LAPACK_WRAPPER_USE_MKL
\*/
#define LAPACK_WRAPPER_USE_LAPACK 1
```

Make sure you have your the corresponding linear algebra software package installed and linked with your buildsystem.

2.3 Selection supported of linear algebra solvers

Comment 2.7 *By default a LAPACK banded KKT solver can be used with OCPBASIC. When only using OCPBASIC with its LAPACK banded KKT solver you do not need to install any other linear solver, which is not included in LAPACK.*

Package:	IPBASIC	OCPBASIC
HSL	✓	✓
LAPACK banded KKT	x	✓
Linear solver interface	✓	x

Table 3: Linear algebra solver packages used by IPBASIC and OCPBASIC

IPBASIC provides a flexible linear solver interface, where users can define own linear solvers, which are then used by IPBASIC. OCPBASIC provides a powerfull structure exploiting linear solver for KKT systems using LAPACK.

For speedtesting also an interface for the sparse linear solvers of HSL is included in IPBASIC. If activated in IPBASIC, MA57 from HSL can also be used in OCPBASIC. Since IPBASIC and OCPBASIC do not provide any copy of HSL, by default the support of HSL and MA57 is not active. If you do have a copy of HSL and you are using qmake, you can activate the support in the SET DEFINES section of the file config.pri by uncommenting the IPBASIC_USE_HSL define.

Comment 2.8 *The user is responsible for linking against any thirdparty software and checking the licence compatibility.*

```
#
# Define use of hsl MA57 and MA48. Make sure a libhsl.so
# with all methods of MA57 and MA48 is installed to your
# system.
#
```

```
#DEFINES += IPBASIC_USE_HSL
```

Do not forget to also set a path to the hsl.so library. This can be done in the SET PATH section of the file config.pri.

```
#  
# Define lib path for hsl:  
#  
HSL_LIB_PATH = /usr/lib/
```

If you do have a copy of HSL and you are using an other buildsystem, you can activate the support in the file IPbasic_Config.h.

```
// Flag for use of hsl:  
#define IPBASIC_USE_HSL
```

2.4 Building and Installing IPbasic and OCPbasic on Ubuntu 20.04

2.4.1 Prepare buildsystem on Ubuntu 20.04

As default qmake is used for building IPBASIC and OCPBASIC. The installing procedure was tested on a fresh Ubuntu 20.04 LTS installation. The buildsystem can be installed by using the shell commands:

```
$ sudo apt-get install g++ -y  
$ sudo apt-get install build-essential -y  
$ sudo apt-get install qt5-default -y
```

Comment 2.9 *The user is responsible for installing and linking against any thirdparty software and checking the licence compatibility.*

2.4.2 Install preferred linear algebra package

Lapack and Blas

Installing lapack and blas by using the shell command:

```
$ sudo apt-get install libblas-dev liblapack-dev -y
```

OpenBLAS

Installing OpenBLAS by using the shell command:

```
$ sudo apt-get install libopenblas-dev -y
```

Intel MKL

Installing Intel MKL by using the shell command:

```
$ sudo apt-get install intel-mkl-full -y
```

Comment 2.10 *The user is responsible for installing and linking against any thirdparty software and checking the licence compatibility.*

Comment 2.11 *After installing the linear algebra packages you have to change to the correct linear algebra package define and enter the correct path variables in the config.pri file as described in section 2.2.1.*

2.4.3 Build and install lapack_wrapper, IPbasic and OCPbasic

Open a terminal in the OCPbasicMaster/ folder. First create a build folder and change your directory into the build folder.

```
$ mkdir build
$ cd build
```

Use qmake to generate a Makefile from the OCPbasicMaster/ folder.

```
$ qmake ../
```

Build LAPACK_WRAPPER, IPBASIC and OCPBASIC using make.

```
$ make
```

Install LAPACK_WRAPPER, IPBASIC and OCPBASIC system wide using make.

```
$ sudo make install
```

After installing LAPACK_WRAPPER, IPBASIC and OCPBASIC you can check the installation using make.

```
$ make check
```

Comment 2.12 *The command make check will execute some unittests. Some unittests return errors, if HSL is not used. Anyway by default OCPBASIC is compiled without HSL and you can use the LAPACK banded KKT solver instead.*

3 Using IPbasic

IPBASIC is an interior-point optimizer for nonlinear optimization problems. It is using a sparse matrix layout defined by the software package LAPACK_WRAPPER. IPBASIC is the nonlinear optimizer used by OCPBASIC. Therefore it is designed for the use of self defined linear solvers.

3.1 Nonlinear optimization problems in IPbasic

The class of nonlinear optimization problems, which can be solved by IPBASIC is defined by the following nlp.

Problem 3.1 (Nonlinear Optimization Problem)

Find the vector $x \in \mathbb{R}^n$ to minimize

$$f(x)$$

subject to the nonlinear constraints

$$h(x) = 0$$

$$g(x) \leq 0$$

with $f : \mathbb{R}^n \rightarrow \mathbb{R}$, $h : \mathbb{R}^n \rightarrow \mathbb{R}^p$ and $g : \mathbb{R}^n \rightarrow \mathbb{R}^q$.

The definition of the interior-point-method used by IPBASIC can be found in [1] and [2].

3.2 Parameter settings of IPbasic

μ_0 :

Notation:	mu_0
Description:	Startvalue of the barrier parameter μ .
Value range:	$[0, 1]$
Default value:	0.5

ϵ_{tol} :

Notation:	epsilon_tol
Description:	Tolerance of the interior-point-method.
Value range:	$[0, 1]$
Default value:	$1e - 8$

Θ^μ :

Notation:	Theta_mu
Description:	Correction parameter of the barrier parameter μ .
Value range:	(1, 2)
Default value:	$\frac{2}{3}$

 κ_ϵ :

Notation:	kappa_epsilon
Description:	Optimization constant.
Value range:	> 0
Default value:	10.0

 κ_μ :

Notation:	kappa_mu
Description:	Optimization constant.
Value range:	(0, 1)
Default value:	0.9

max_{iter}:

Notation:	max_iter
Description:	Max number of iterations.
Value range:	> 0
Default value:	1000

 α_{\max} :

Notation:	alpha_max
Description:	Max stepsize.
Value range:	> 0
Default value:	1

 α_{tol} :

Notation:	tol_alpha
Description:	Min stepsize.
Value range:	> 0

Default value:	$2e - 8$
-----------------------	----------

σ :

Notation:	sigma
Description:	Parameter for the armijo-rule.
Value range:	(0, 1)
Default value:	0.01

β :

Notation:	beta
Description:	Parameter for the armijo-rule.
Value range:	(0, 1)
Default value:	0.9

τ_{\min} :

Notation:	tau_min
Description:	Distance parameter for the bounds.
Value range:	(0, 1)
Default value:	$1e - 4$

Printlevel:

Notation:	PrintLev
Description:	Printlevel of IPbasic.
Value range:	PRINT_NONE = No printing. PRINT_LESS = Printing only solution. PRINT_MORE = Printing iterations and solution. PRINT_MARKED = Printing iterations and solution with colors.
Default value:	PRINT_MORE

λ_{tol} :

Notation:	lambda_tol
Description:	Tolerance of the lagrange multipliers.
Value range:	(0, 1)
Default value:	$1e - 10$

RegDiagMatrixEqu:

Notation:	RegDiagMatrixEqu
Description:	Regularisation parameter of the diagonalmatrix of the equations.
Value range:	
Default value:	0

RegDiagMatrixInitialGuessLambda:

Notation:	RegDiagMatrixInitialGuessLambda
Description:	Regularisation parameter of the diagonalmatrix lagrange multipliers.
Value range:	
Default value:	$1e - 9$

CheckMatricesFor_INF_NAN:

Notation:	CheckMatricesFor_INF_NAN
Description:	Tells if matrices are checked due to inf and nan.
Value range:	bool
Default value:	true

LineSearchCouples:

Notation:	MyLineSearchCouples
Description:	Used LineSearchCouples.
Value range:	COUPLES_NONE = No couples. COUPLE_STATE_SLACK = Couple linesearch of slack and state variables. COUPLE_STATE_SLACK_LAMBDA = Couple linesearch of slack variables, state variables and lagrange multipliers.
Default value:	COUPLE_STATE_SLACK

BreakpointMaxTime:

Notation:	BreakpointMaxTime
Description:	Tells, if there is a breakpoint with a maximal time.
Value range:	bool
Default value:	false

maximalTime:

Notation:	maximalTime
Description:	If BreakpointMaxTime, maximal time for IPbasic to solve. [in s]
Value range:	> 0
Default value:	1

3.2.1 Optimizer flags for IPbasic

LineSearchMethod:

Notation:	UsedLineSearchMethod
Description:	Used linesearch method in IPBASIC.
Value range:	LINESEARCH_NONE = No linesearch used. LINESEARCH_ARMIJO_L1_PENALTY = Armijo linesearch with ℓ_1 -penalty-function.
Default value:	LINESEARCH_ARMIJO_L1_PENALTY

4 Examples

Some examples for the use of IPBASIC can be found in the examples/examples_IPbasic/ folder, which is included in OCPbasicMaster.

4.1 IPbasic using dense linear solver

For using IPBASIC without OCPBASIC or HSL you will need a linear solver for solving nlps. For small nlps you might want to use a dense linear solver from LAPACK. IPBASIC is desinged for overloading the linear solver. This example will show how a dense linear solver can be defined with IPBASIC and LAPACK_WRAPPER.

Dense linear solver

IPBASIC is using the SparseMatrixBase class from LAPACK_WRAPPER. To create a dense linear solver you need to overload the LinearSolver_interface class from IPBASIC. A header only definition of a dense linear solver, using linear algebra methods of LAPACK_WRAPPER, can be created by the following code example.

```
using lapack_wrapper;

class DenseSolver : public
IPbasic::LinearSolver_interface<double>
{
lapack_wrapper::LU<double> lu;

public:

DenseSolver()
: LinearSolver_interface<double>()
{
}

virtual ~DenseSolver()
{
}

bool factorize(SparseMatrixBase<double> const & Sparse,
               bool is_stored_symmetric) IPBASIC_OVERRIDE
{
    int const *    pRows;
    int const *    pCols;
    double const * pValues;
    int            numRows, numCols, nnz;
    Sparse.get_info(numRows, numCols, nnz);
    Sparse.get_data(pRows, pCols, pValues);
    lu.allocate(numRows, numCols);
    if (is_stored_symmetric)
    {
        lu.load_sparse_sym(nnz, pValues, pRows, pCols);
    }
    else
    {
        lu.load_sparse(nnz, pValues, pRows, pCols);
    }
    lu.factorize("factorize");
    return true;
}
```

```
}  
  
bool solve(real_type RHS[]) IPBASIC_OVERRIDE  
{  
    lu.solve(RHS);  
    return true;  
}  
};
```

Dense nlp problem

Since the default matrixtype of IPBASIC is using sparse matrices, we should use the `lapack_wrapper::SparseCCOOR` type for the jacobian and hessian. Still we can use a dense matrix storage layout in `LAPACK_WRAPPER`. This storage layout is created by the method `setup_as_full_column_major(...)`.

```
class Problem : public IPbasic::NLP<real_type>  
{  
    lapack_wrapper::SparseCCOOR<real_type> Constraint_Jacobian;  
    lapack_wrapper::SparseCCOOR<real_type> Lagrangian_Hessian;  
    lapack_wrapper::SparseCCOOR<real_type> MyIPMatrix;  
    bool Is_equality[2];  
  
public:  
    Problem()  
    :NLP<real_type>(&this->Constraint_Jacobian,  
    &this->Lagrangian_Hessian, &this->MyIPMatrix,  
    this->Is_equality, false, false)  
    {  
        this->Constraint_Jacobian.setup_as_full_column_major(2, 2);  
        this->Lagrangian_Hessian.setup_as_full_column_major(2, 2);  
        this->MyIPMatrix.setup_as_full_column_major(4, 4);  
        this->is_equality_vec[0] = false;  
        this->is_equality_vec[1] = false;  
    }  
  
    int_type state_dimension() const IPBASIC_OVERRIDE  
    {  
        return 2;  
    }  
};
```

```
}

virtual void initial_guess(real_type x[]) const
IPBASIC_OVERRIDE
{
    x[0] = 0.6;
    x[1] = 2.0;
}

virtual real_type objective(real_type const x[])
IPBASIC_OVERRIDE
{
    return x[0] * x[0] + x[1] * x[1];
}

virtual void objective_gradient(real_type const x[],
real_type grad[])
IPBASIC_OVERRIDE
{
    grad[0] = 2 * x[0];
    grad[1] = 2 * x[1];
}

virtual int_type constraint_dimension()
const IPBASIC_OVERRIDE
{
    return 2;
}

virtual void constraint(real_type const x[],
real_type c[])
IPBASIC_OVERRIDE
{
    c[0] = -x[1] + 1.5;
    c[1] = -x[1] + exp(x[0]);
}

void constraint_jacobian_update(real_type const x[])
```

```
IPBASIC_OVERRIDE
{
    lapack_wrapper::MatrixWrapper<real_type> M;
    this->Constraint_Jacobian.get_full_view(M);
    M(0, 0) = 0.0;
    M(0, 1) = -1.0;
    M(1, 0) = exp(x[0]);
    M(1, 1) = -1.0;
}

void lagrangian_hessian_update(real_type sigma,
real_type const x[],
real_type const lambda[])
IPBASIC_OVERRIDE
{
    lapack_wrapper::MatrixWrapper<real_type> M;
    this->Lagrangian_Hessian.get_full_view(M);
    M(0, 0) = 2 * sigma + lambda[1] * exp(x[0]);
    M(0, 1) = 0.0;
    M(1, 0) = 0.0;
    M(1, 1) = 2 * sigma;
}
};
```

Using a dense linear solver

The created dense linear solver can directly be used by IPBASIC. All linear systems are now solved by a dense lu factiozation.

```
int main(int, char *[])
{
    Problem      prob;
    DenseSolver  MyDenseSolver;

    IPbasic::Solver<double> *      MySolver =
        new IPbasic::Solver<double>(&prob, &MyDenseSolver);
    IPbasic::IPparameter<double> * Param    =
        MySolver->getIPparameter();
    Param->mu_0      = 1e-3;
```

```
if (MySolver->optimize())
{
    std::cout << "Job_done!" << std::endl;
}
else
{
    std::cout << "Job_not_done!" << std::endl;
}

delete MySolver;
return 0;
}
```

References

- [1] A. Huber, M. Gerds, and E. Bertolazzi. Structure exploitation in an interior-point method for fully discretized, state constrained optimal control problems. *Vietnam Journal of Mathematics*, Oct 2018.
- [2] A. K.-P. Huber. *Methoden zur Berechnung optimaler Rennlinien im dynamischen Grenzbereich*. PhD thesis, Universität der Bundeswehr München, 2020.