# **OPTOOL** - Documentation

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#### Abstract

The OPTool package is an implementation of various state-of-the-art iterative optimization algorithms for differentiable cost functions along with algorithms to solve linear equations. Users can use the toolbox to solve optimization problems, although the code was written to researchers that want to compare their proposals with state-of-the-art implementation. New algorithms can be easily added and the software will be updated to have the most comprehensive list of solvers possible. It also comes with implemented functions to return optimal parameters for these algorithms based on a control-theoretical formulation of the algorithms.

Keywords: Optimization Problems, Control-theoretical Formalization, Gradient-descent-like Algorithms

## 1. Introduction and background

The package OPTOOL has various algorithms to solve optimization problems where x denote the variable and for a general cost function f written as:

$$\underset{x}{\text{minimize}} \quad f(x) \tag{1}$$

Function f is assumed differentiable, i.e., there exists  $\nabla f$ , so subgradient methods are currently not implemented. There exist multiple gradient-descent algorithms and we use two in this tutorial to help the user successfully run and understand its first example. The steepest gradient descent labeled as Gradient and the Nesterov method are:

GRADIENT: 
$$x^{(k+1)} = x^{(k)} - \beta \nabla f(x^{(k)})$$

NESTEROV: 
$$x^{(k+1)} = \xi^{(k)} - \beta \nabla f(\xi^{(k)})$$

$$\xi^{(k)} = (1+\gamma)x^{(k)} - \gamma x^{(k-1)}$$
(2)

with  $\beta$ ,  $\gamma$  selected as parameters.

The aforementioned methods are general to any function f and, given appropriate choice for parameters, they converge to the global minimum provided f is convex. Whenever this is not the case, convergence occurs for on of the points where  $\nabla f = \mathbf{0}$ , which can be a minimum, maximum or saddle point.

If we specialize function f to be quadratic, i.e.,  $f = \frac{1}{2} ||Mx - b||_2^2$ , the algorithms are linear as  $\nabla f$  is a linear function of x. Therefore, iterative algorithms to solve linear equations can also be applied to the equation Mx = b.

OPTOOL implements many methods that are present in the literature. See the README file for the complete list.

#### 2. Installation

To install simply download the zip file and extract it to the Matlab folder and add it to the path.

# 3. Usage

To use the package to solve a general optimization algorithm, one can use the function:

```
function [ stateVectors , errors ] = optSolver(
   algorithms , parameters , grad , errorFunction ,
   max_iterations , initialState , projectionFunction
   , tol)
```

With inputs:

**algorithms** a cell array containing pointers to functions implementing the next iteration of the algorithms;

**parameters** a cell array containing structures with the parameters for each algorithm in **algorithms**;

**grad** a pointer to the gradient function;

**errorFunction** a pointer to the function that computes the error for a particular x;

max\_iterations maximum allowed number of iterations to achieve the solution;

**initialState** initial guess for the minimum of f;

- **projectionFunction** optional input containing a pointer to a function that projects the state onto some constraining set;
- tol tolerance before the algorithm halts and sets all subsequent time steps equal to the current one.

and outputs:

- **stateVectors** a cell array containing the matrices  $[x(0) \cdots x(\text{max\_iterations})]$  for each of the selected algorithms;
- **errors** cell array containing the vectors of errors for each of the algorithms using **errorFunction**.

Similarly, the function:

```
function [ stateVectors , errors ] = linSolver(
   algorithms , parameters , A , b , errorFunction,
   max_iterations , initialState, projectionFunction
   , tol)
```

solves the linear equation Ax = b. Instead of providing the gradient function, the user must supply A and b and all the rest of the inputs are equal.

In case the optimization problem is quadratic, the user can call function:

```
function [ stateVectors , errors ] = quadSolver(
   algorithmNames , parameters , A , b ,
   errorFunction, max_iterations , initialState ,
   experimentName , projectionFunction , tol ,
   errorDescription)
```

giving the additional **experimentName** that will be used to save the resulting plots and variables in the folder *Stored Outputs*.

To facilitate the setup of the problems to be computed, inside the folder *Optimal Parameters* there is the function:

```
function [algorithms, parameters, names] =
  getParameters(A, b, r, methods)
```

With inputs:

**A** a matrix such that Q = A'A;

**b** a vector such that p = A'b;

**r** a value such that  $f(x) = 0.5x^{\mathsf{T}}Qx - p^{\mathsf{T}}x + r$ ;

**methods** a string array with all names of desired algorithms to solve the problem.

and outputs:

- algorithms a structure containing optimization that is the algorithms input to optSolver and linearEquations which should be used for linSolver:
- parameters a structure containing optimization that is the parameters input to optSolver and linearEquations which should be used for linSolver:
- names a sorted string array of the names of the selected methods present in the above data structures (badly specified methods will not appear in this output).

If this function is used for a non-quadratic problem, matrix A should be such that its minimum and maximum eigenvalues are the smoothness and strong convexity values of the function. For a list of the available methods, just run "getParameters" with no inputs or to select all "getParameters(A, b, r, 'all ')".

## 4. Implementing additional algorithms

The OPTOOL was implemented such that adding other algorithms to the package is straightforward. If a NOVELALGORITHM needs to be added, the developer must:

1. Define the NOVELALGORITHM function that implements a single iteration of the proposed algorithm. The function should follow the definition:

```
function [ x , parameters] = novelAlgorithm( x ,
    previous_x , grad, parameters)
```

With inputs:

 $\mathbf{x}$  current vector estimation of the minimum of f;

**previous\_x** previous vector estimation of the minimum of f;

**grad** pointer to the gradient function;

**parameters** structure used to pass parameters between algorithm iterations.

and outputs:

 $\mathbf{x}$  new vector estimation of the minimum of f;

parameters updated structure of the parameters for this algorithm.

- 2. Store the above file in either folder *Optimization Algorithms* if it is a general gradient descent algorithm or in *Linear Equation Solver* if it solves Ax = b;
- 3. Add reference to the paper defining the algorithm in *References*;
- 4. In the file *getParameters.m* add the correspondence between name of the algorithm and the typical name of the parameters either in the variable **optAlgorithmNames** or **linEqAlgorithmNames** according to the type of algorithm;
- 5. Still in the file getParameters.m add an if clause similar to the one presented for the gradient descent:

```
% %==== Optimization Algorithms ====
% Gradient Descent
if strcmp(methods(i), "Gradient Descent") ||
    allMethods
optAlgorithms{optIndex} = @gradientDescent;
optParameters{optIndex} = struct('alpha',2/(L + m));
optIndex = optIndex + 1;
end
```

#### for example:

```
%==== Optimization Algorithms ====
  % Gradient Descent
  if strcmp(methods(i), "Gradient Descent") ||
3
      allMethods
  optAlgorithms{optIndex} = @gradientDescent;
4
   optParameters{optIndex} = struct('alpha',2/(L +
  optIndex = optIndex + 1;
6
  end
8
  % Novel Algorithm
  if strcmp(methods(i),"Novel Algorithm") ||
      allMethods
  optAlgorithms{optIndex} = @novelAlgorithm;
10
   optParameters{optIndex} = struct('myParameter',<</pre>
     myparameterValue>);
```

```
12 optIndex = optIndex + 1;
13 end
```

## 5. Illustrative Example

The following example comes in file PageRank.m. The PageRank problem consists of a ranking mechanism from Google, which was initially proposed in [1]. It corresponds to finding the eigenvector of the following matrix  $M \in \mathbb{R}^{n \times n}$ :

$$M := (1 - m)A + \frac{m}{n}S\tag{3}$$

where  $m \in (0,1)$  is a parameter defining the convex combination of the adjacency matrix A of the network with the matrix  $S := \mathbf{1}_n \mathbf{1}_n^{\mathsf{T}}$  ( $\mathbf{1}_n$  is the n-dimensional vector of ones). A typical choice is m = 0.15 [1]. The standard formulation can be efficiently computed through the power method:

$$x(k+1) = Mx(k) = (1-m)Ax(k) + \frac{m}{n}\mathbf{1}_n$$
 (4)

where  $x(k) \in \mathbb{R}^n$  and  $\forall k \geq 0 : \mathbf{1}_n^{\intercal} x(k) = 1$ .

The PageRank problem can also be formulated as an optimization problem or as the solution to a linear equation. In the former, the PageRank is the solution to the following optimization problem

minimize 
$$\frac{1}{2} \| ((1-m)A - I_n)x + \frac{m}{n} \mathbf{1}_n \|_2^2$$

If seen as the solution of a linear equation in matrix format we get:

$$(I_n - (1-m)A)x = \frac{m}{n}1_n.$$
 (5)

In [2], it is shown that the standard power method for the PageRank is equivalent to the Jacobi method applied to (5). The solution of the PageRank for a random Barabási–Albert generated network is included in the OPTOOL and the plot of the errors produced by the toolbox is given in Figure ??, where it is shown better alternatives to the PageRank if the optimal parameters are known by each node.

## 6. Plans for future releases

In future versions, we expect to add subgradient methods and also databased training such as the Stochastic Gradient Descent. It is also going to

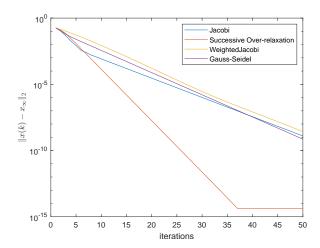


Figure 1: The error evolution for each of the tested algorithm in the PageRank case for a 20-node network.

be added work in progress on novel algorithms for optimization and other convergence rates calculations.

We also would like to make available additional examples of paper [3] and other works under development on power networks. Topics such as consenus, both deterministic and stochastic [4], [5].

#### References

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