

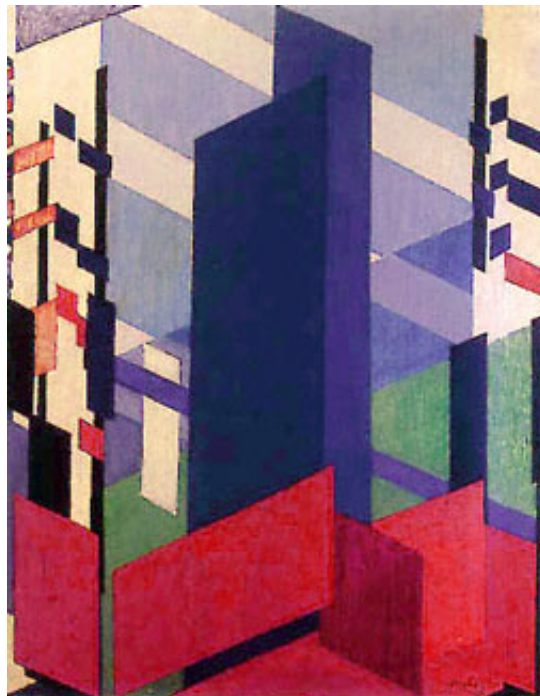
COURSE ON LMI OPTIMIZATION  
WITH APPLICATIONS IN CONTROL  
PART I.4

**SOLVING LMIs**

Didier HENRION

[www.laas.fr/~henrion](http://www.laas.fr/~henrion)

[henrion@laas.fr](mailto:henrion@laas.fr)



Architecture Philosophique (1913)  
František Kupka (1871-1957)

November 2003

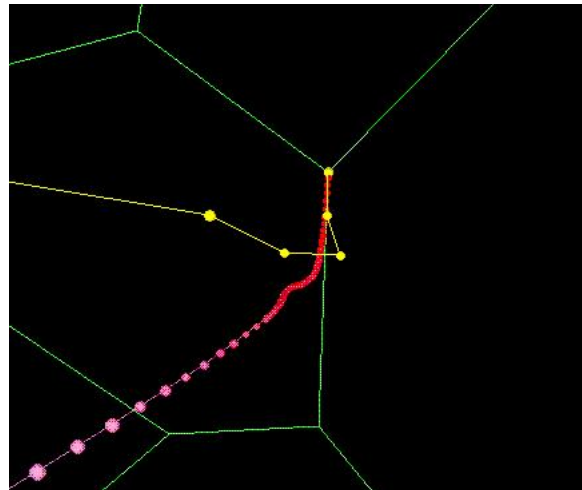
## History

### Convex programming

- logarithmic barrier function (Frisch 1955)
- method of centers (Huard 1967)

### Interior-point (IP) methods

- ellipsoid algorithm (Khachiyan 1979)  
polynomial bound on worst-case iteration count
- IP methods for LP (Karmarkar 1984)  
improved complexity bound and efficiency - now about 50% of commercial LP solvers
- self-concordant barrier functions (Nesterov, Nemirovski 1988, Alizadeh 1991) - IP methods for general convex programs, in particular SDP and LMI



Steve Wright's artistic view of interior point methods in action  
Iterates (yellow) LP edges (green) central path (red)

## Logarithmic barrier function

For the optimization problem

$$\begin{array}{ll} \min & f_0(x) \\ \text{s.t.} & f_i(x) \geq 0 \end{array}$$

where the  $f_i(x)$  are twice continuously differentiable convex functions, we define the **logarithmic barrier** function

$$\phi(x) = -\sum_i \log f_i(x) = \log \prod_i f_i(x)^{-1}$$

which is **convex** in the interior  $f_i(x) > 0$  of the feasible set

Then we solve the **unconstrained** convex problem

$$\min f_0(x) + \mu\phi(x)$$

where  $\mu > 0$  and the term  $\mu\phi(x)$  acts as a “repellent” of the boundary

The minimum is attained in the interior  
= **interior-point** method

## Descent methods

To solve an unconstrained optimization problem

$$\min f(x)$$

for  $x \in \mathbb{R}^n$  we produce a minimizing sequence

$$x_{k+1} = x_k + t_k \Delta x_k$$

where  $\Delta x_k \in \mathbb{R}^n$  is the **step** or **search direction** and  $t^{(k)} \geq 0$  is the **step size** or **step length**

A **descent method** consists in finding a sequence  $\{x_k\}$  such that

$$f(x^*) \leq \dots f(x_{k+1}) < f(x_k)$$

where  $x^*$  is the optimum

### General descent method

0. given starting point  $x$
1. determine **descent direction**  $\Delta x$
2. line search: choose **step size**  $t > 0$
3. update:  $x = x + t \Delta x$
4. go to step 1 until a stopping criterion is satisfied

## Newton's method

A particular choice of search direction is the **Newton step**

$$\Delta x = -\nabla^2 f(x)^{-1} \nabla f(x)$$

where  $\nabla f(x)$  is the **gradient**  
and  $\nabla^2 f(x)$  is the **Hessian**

This step  $y = \Delta x$  minimizes the second-order Taylor approximation

$$\hat{f}(x + y) = f(x) + \nabla f(x)^T y + y^T \nabla^2 f(x) y / 2$$

and it is the steepest descent direction for the quadratic norm defined by the Hessian

**Quadratic convergence** near the optimum

## Self-concordance

Shortcomings of Newton's method:

- number of required Newton steps hardly estimated in practice
- analysis depends on used coordinate system

Theory of **self-concordant** functions:

- number of Newton steps easily estimated
- affine-invariant property

Smooth convex functions with 2nd derivatives Lipschitz continuous with respect to the metric induced by the Hessian:

$$|f'''(x)| \leq 2f''(x)^{3/2}$$

include many logarithmic barrier functions

For LP, QP or SDP **1st and 2nd derivatives** of standard self-concordant barriers can be found easily in closed form

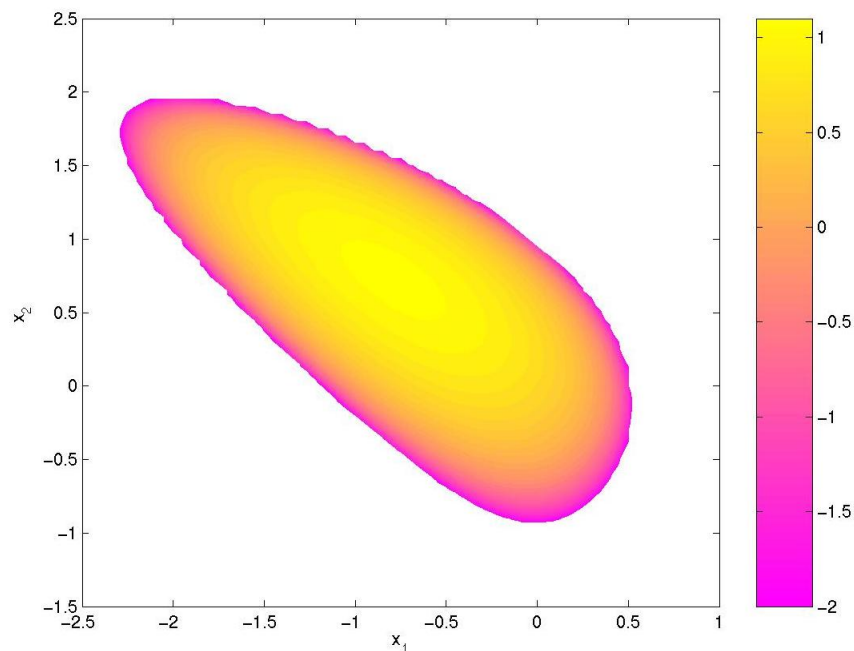
## Barrier function for an LMI

Given an LMI constraint  $F(x) \succeq 0$  we define its **logarithmic barrier** function

$$\phi(x) = -\log \det F(x) = \log \det F(x)^{-1}$$

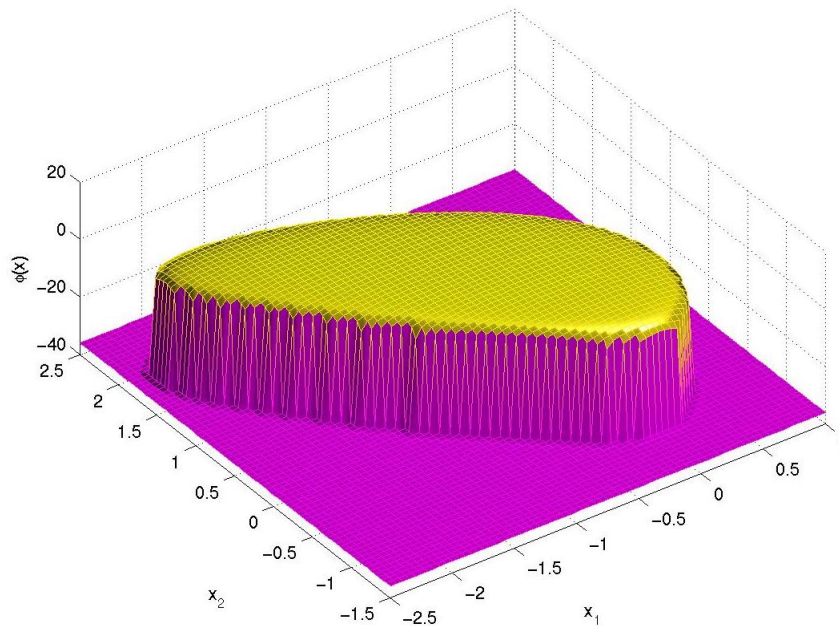
This function is analytic, convex and **self-concordant** on  $\{x : F(x) \succ 0\}$

The optimum of  $\min \phi(x)$  is called the **analytic center** of the LMI



## Gradient and Hessian for an LMI

The barrier function  $\phi(x)$  is flat in the interior of the feasible set and sharply increases toward the boundary



Closed-form expressions for gradient

$$\begin{aligned}(\nabla\phi(x))_i &= -\text{trace } F(x)^{-1}F_i \\ &= -\text{trace } F(x)^{-1/2}F_iF(x)^{-1/2}\end{aligned}$$

and Hessian

$$\begin{aligned}(\nabla^2\phi(x))_{ij} &= \text{trace } F(x)^{-1}F_iF(x)^{-1}F_j \\ &= \text{trace } (F(x)^{-1/2}F_iF(x)^{-1/2})(F(x)^{-1/2}F_jF(x)^{-1/2})\end{aligned}$$



## IP methods for SDP

Primal / dual SDP

$$\begin{array}{ll} \min_X & \text{trace } CX \\ \text{s.t.} & \text{trace } A_i X = b_i \\ & X \succeq 0 \end{array}$$

$$\begin{array}{ll} \max_y & b^T y \\ \text{s.t.} & Z = C - \sum_i y_i A_i \\ & Z \succeq 0 \end{array}$$

### Primal methods

$$\begin{array}{ll} \min_X & \text{trace } CX - \mu \log \det X \\ \text{s.t.} & \text{trace } A_i X = b_i \end{array}$$

where parameter  $\mu$  is sequentially decreased to zero and iterates  $X_k$  are always **primal feasible**

### Dual methods

$$\begin{array}{ll} \max_{y,Z} & b^T y + \mu \log \det Z \\ \text{s.t.} & Z = C - \sum_i y_i A_i \end{array}$$

where parameter  $\mu$  is sequentially decreased to zero and iterates  $y_k, Z_k$  are always **dual feasible**

$X_k \succeq 0$  or  $Z_k \succeq 0$  ensured via Newton process:

- large decreases of  $\mu$  require damped Newton steps
- small updates allow full (deep) Newton steps

## Primal-dual IP methods for SDP

### Primal-dual methods

$$\begin{aligned} \min_{x,y,Z} \quad & \text{trace } XZ - \mu \log \det XZ \\ \text{s.t.} \quad & \text{trace } A_i X = b_i \\ & Z = C - \sum_i y_i A_i \end{aligned}$$

Minimizers satisfy optimality conditions

$$\begin{aligned} \text{trace } A_i X &= b_i \\ \sum_i y_i A_i + Z &= C \\ XZ &= \mu I \\ X, Z &\succeq 0 \end{aligned}$$

Duality gap

$$\text{trace } CX - b^T y = \text{trace } XZ \succeq 0$$

is minimized along the **central path** of solutions  $X(\mu)$ ,  $y(\mu)$ ,  $Z(\mu)$ , a smooth curve parametrized by scalar  $\mu$

For this reason, logarithmic barrier methods are also called **path-following methods**

## Newton step for LMI

For the SDP in LMI form

$$\begin{aligned} \min \quad & c^T x \\ \text{s.t.} \quad & F(x) = F_0 + \sum_i x_i F_i \succeq 0 \end{aligned}$$

the centering problem is

$$\min c^T x - \mu \log \det F(x)$$

and at each iteration Newton step  $\Delta x$  satisfies the [linear system of equations](#) (LSE)

$$H \Delta x = -g$$

where gradient  $g$  and Hessian  $H$  are given by

$$\begin{aligned} H_{ij} &= \text{trace } F(x)^{-1} F_i F(x)^{-1} F_j \\ g_i &= c_i / \mu + \text{trace } F(x)^{-1} F_i \end{aligned}$$

LSE typically solved via [Cholesky factorization](#) or [QR decomposition](#) (near the optimum)

## Newton step for primal-dual methods

For primal-dual IP methods, primal and dual directions  $\Delta X$ ,  $\Delta y$  and  $\Delta Z$  satisfy **non-linear** KKT optimality conditions

$$\begin{aligned}\text{trace } A_i \Delta X &= 0 \\ \sum_i \Delta y_i A_i + \Delta Z &= 0 \\ (X + \Delta X)(Z + \Delta Z) &= \mu I\end{aligned}$$

Key point is in **linearizing** and **symmetrizing** the latter equation



Long list of primal-dual **search directions**, the most known of which is Nesterov-Todd's

Dynamic updates of  $\mu$  result in **predictor-corrector** methods

## Complexity

For the  $n$ -by- $n$  LMI  $F(x) \succeq 0$  with  $m$  variables  
the **flops count** of IP methods for SDP is as follows:

For each iteration:

- (a)  $\mathcal{O}(n^2m)$  to form  $F(x)$
- (b)  $\mathcal{O}(n^3m)$  to form  $F(x)^{-1}F_iF(x)^{-1}F_j$
- (c)  $\mathcal{O}(n^2m^2)$  to form  $F(x)^{-1}F_i$
- (d)  $\mathcal{O}(m^3)$  to solve Newton LSE with Cholesky

Dominating terms are (b) and (c) so the complexity for solving one Newton step is:

$$\mathcal{O}(n^3m + n^2m^2)$$

..but **structure** can be exploited in these steps !

Number of iterations with Newton's method:

$$\mathcal{O}(\sqrt{n} \log \varepsilon^{-1})$$

where  $\varepsilon$  is the desired accuracy

In general, it is assumed that  $m = \mathcal{O}(n^2)$  otherwise redundant constraints can be removed, so the global **worst-case complexity** for a dense SDP is

$$\mathcal{O}(n^{6.5} \log \varepsilon^{-1})$$

Much less in practice !

## Primal-dual and barrier methods

In contrast with barrier methods,  
in primal-dual methods..

- At each step **both** primal and dual variables are updated simultaneously
- Search direction obtained from Newton's method applied to **modified KKT equations**
- Work when problem is **not strictly feasible**
- Iterates are primal and dual **infeasible** except when converging

Generally for LP, QP or SDP primal-dual methods outperform barrier methods

## IP methods in general

General characteristics of IP methods:

- **Efficiency**: about 5 to 50 iterations, almost independent of input data (problem), each iteration is a least-squares problem (well established linear algebra)
- **Theory**: worst-case analysis of IP methods yields polynomial computational time
- **Structure**: tailored SDP solvers can exploit problem structure

For more information see the Linear, Cone and SDP section at

[www.optimization-online.org](http://www.optimization-online.org)

and the Optimization and Control section at

[fr.arXiv.org/archive/math](http://fr.arXiv.org/archive/math)

## Penalty/augmented Lagrangian methods

Use similar ideas, but cannot be considered as an interior-point method

When applied to LMI problem

$$\min c^T x \text{ s.t. } F(x) = F_0 + \sum_i x_i F_i \succeq 0$$

- penalty method - **some** eigenvalues of  $F(x)$  can be negative
- barrier method - **no** eigenvalue of  $F(x)$  can be negative (use of logarithm)

### Augmented Lagrangian

$$L(x, Z, p) = c^T x + \text{trace } Z \Phi(x, p)$$

with dual variable  $Z$  and suitable **penalty** function, for example

$$\Phi(x, p) = p^2 (F(x) + pI)^{-1} - pI$$

with penalty parameter  $p$



## Penalty/augmented Lagrangian methods (2)

### General algorithm

1. find  $x_{k+1}$  such that  $\|\nabla_x L(x, Z_k, p_k)\| \leq \epsilon_k$
2. update dual variables:  $Z_{k+1} = f(x_{k+1}, Z_k)$
3. update penalty parameter:  $p_{k+1} < p_k$
4. go to step 1 until a stopping criterion is satisfied

Can be considered as a **primal-dual** method, but dual variables are obtained in **closed-form** at step 2

Complexity for the  $n$ -by- $n$  LMI  $F(x) \succeq 0$  with  $m$  variables depends mostly on Newton step 1 in  $O(n^3m + n^2m^2)$ , same as IP methods

Can be improved to  $O(m^2K^2)$  where  $K$  is the max number of non-zero terms in the  $F_i$

## SDP solvers

Available under the **Matlab** environment

Primal-dual path-following predictor-corrector algorithms:

- **SeDuMi** (Sturm)
- **SDPT3** (Toh, Tütüncü, Todd)
- **CSDP** (Borchers)
- **SDPA** (Kojima and colleagues)

parallel version available

Primal-dual potential reduction:

- **MAXDET** (Wu, Vandenberghe, Boyd)
- explicit max det terms in objective function

Dual-scaling path-following algorithms:

- **DSDP** (Benson, Ye, Zhang)

exploits structure for combinatorics

Barrier method and augmented Lagrangian:

- **PENSDP** (Kočvara, Stingl)

## Matrices as variables

Generally, in control problems we do not encounter the LMI in canonical or semidefinite form but rather with **matrix variables**

Lyapunov's inequality

$$A^T P + P A < 0 \quad P = P^T > 0$$

can be written in canonical form

$$F(\mathbf{x}) = F_0 + \sum_{i=1}^m F_i x_i > 0$$

with the notations

$$F_0 = 0 \quad F_i = -A^T B_i - B_i A$$

where  $B_i$ ,  $i = 1, \dots, n(n+1)/2$  are matrix bases for symmetric matrices of size  $n$

Most software packages for solving LMIs however work with canonical or semidefinite forms, so that a (sometimes time-consuming) **pre-processing step** is required

## LMI solvers

Available under the **Matlab** environment

Projective method: project iterate on ellipsoid within PSD cone = least squares problem

- **LMI Control Toolbox** (Gahinet, Nemirovski) exploits structure with rank-one linear algebra warm-start + generalized eigenvalues originally developed for INRIA's **Scilab**

LMI interface to SDP solvers

- **LMITOOL** (Nikoukah, Delebecque, El Ghaoui) for both Scilab and Matlab
- **SeDuMi Interface** (Peaucelle)
- **YALMIP** (Löfberg)

See Helmberg's page on SDP

[www-user.tu-chemnitz.de/~helmberg/semidef.html](http://www-user.tu-chemnitz.de/~helmberg/semidef.html)

and Mittelmann's page on optimization software with benchmarks

[plato.la.asu.edu/guide.html](http://plato.la.asu.edu/guide.html)

## LMI relaxation software

**GloptiPoly** is written as an open-source, general purpose and user-friendly **Matlab** software

Optionally, problem definition made easier with Matlab Symbolic Math Toolbox, gateway to **Maple** kernel

Gloptipoly solves small to medium **non-convex** global optimization problems with multivariate real-valued **polynomial** objective functions and constraints

Software and documentation available at

[www.laas.fr/~henrion/software/gloptipoly](http://www.laas.fr/~henrion/software/gloptipoly)

## Methodology

GloptiPoly builds and solves a **hierarchy** of successive **convex linear matrix inequality (LMI) relaxations** of increasing size, whose optima are **guaranteed** to converge asymptotically to the global optimum



Relaxations are build from LMI formulation of **sum-of-squares (SOS)** decomposition of multivariate polynomials (see last chapter)

In practice convergence is ensured **fast**, typically at 2nd or 3rd LMI relaxation

## Features

General features of GloptiPoly:

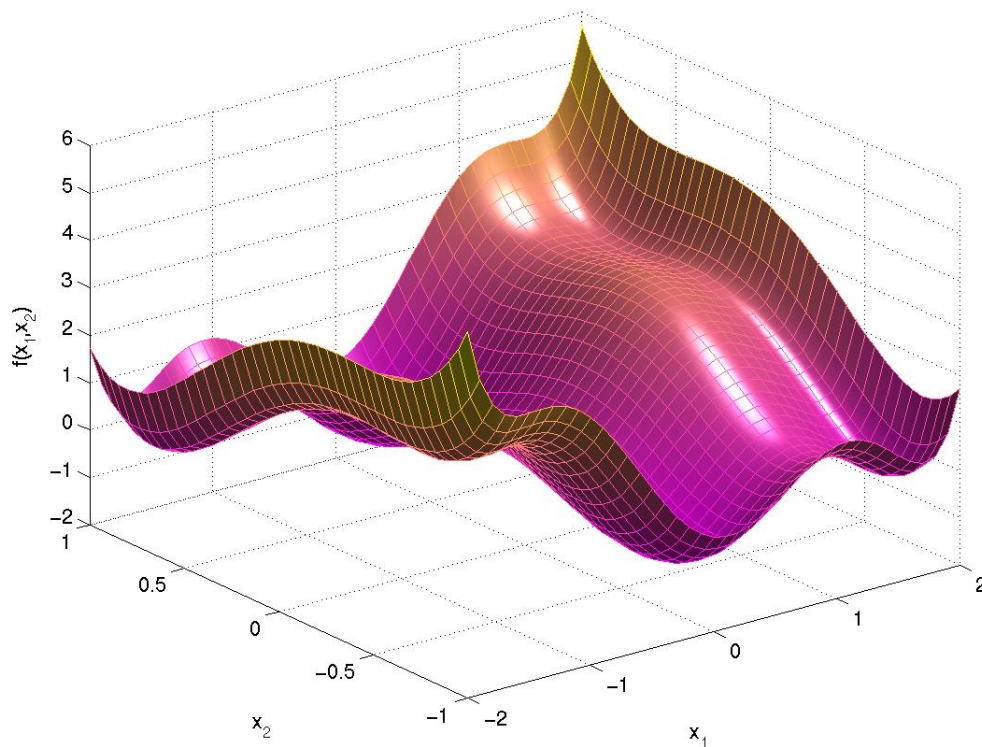
- Certificate of global optimality (rank checks)
- Automatic extraction of globally optimal solutions (multiple eigenvectors)
- 0-1 or  $\pm 1$  integer constraints on some of the decision variables (combinatorial optimization problems)
- Generation of input and output data in SeDuMi's format
- Generation of moment matrices associated with LMI relaxations (rank checks)
- User-defined scaling of decision variables (to improve numerical behavior)
- Exploits sparsity of polynomial data

Major update of Gloptipoly planned (hopefully !) for winter 2003

## Benchmark examples Continuous problems

Mostly from Floudas/Pardalos 1999 handbook

About 80 % of pbs solved with LMI relaxation of small order (typically 2 or 3) in less than 3 seconds on a PC Pentium IV at 1.6 MHz with 512 Mb RAM



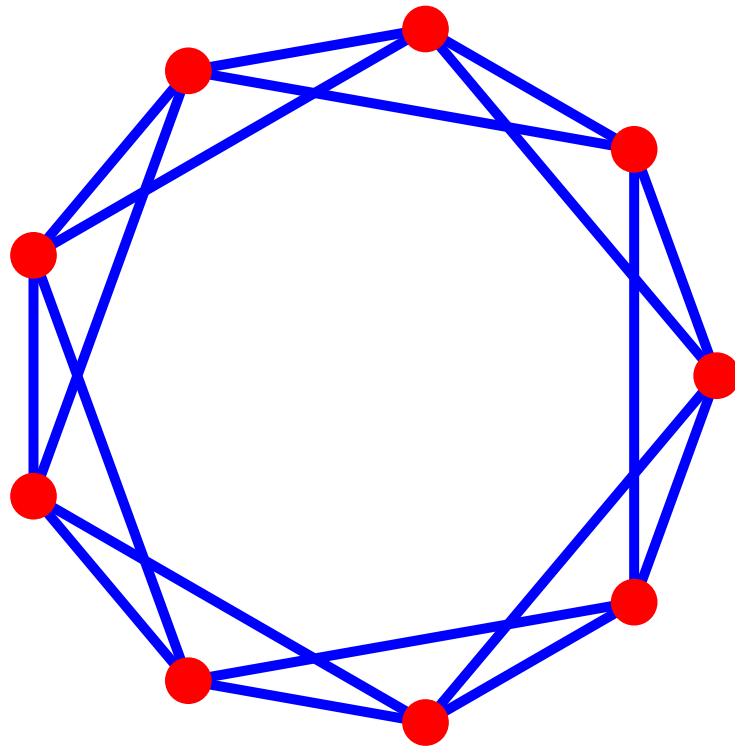
Six-hump camel back function



## Benchmark examples Discrete problems

From Floudas/Pardalos handbook and also  
Anjos' Ph.D (Univ Waterloo)

By perturbing criterion (destroys symmetry)  
global convergence ensured on **80 %** of pbs  
in **less than 4 seconds**



MAXCUT on antiweb  $AW_9^2$  graph

## Benchmark examples

### Polynomial systems of equations

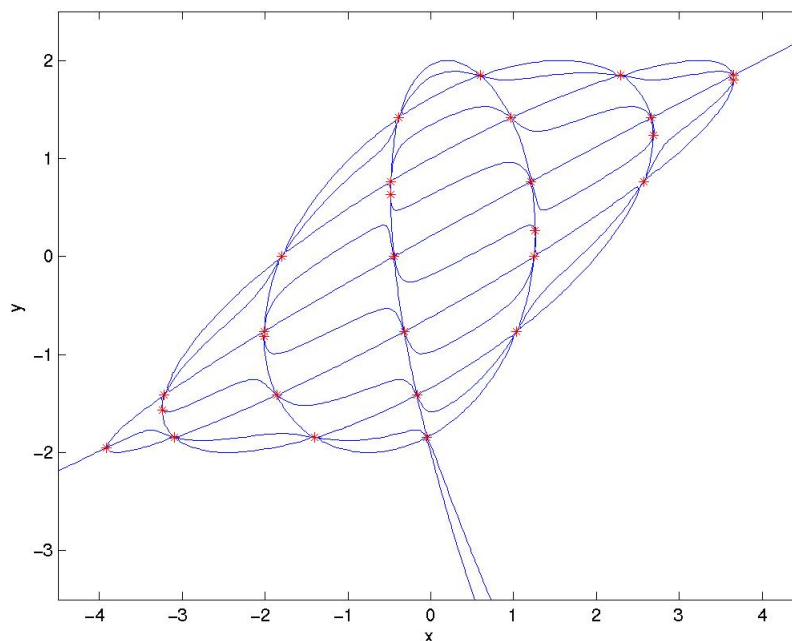
From Verschelde's and Posso database  
Real coefficients & coeffs only

Out of 59 systems:

- 61 % solved in  $t < 10$  secs
- 20 % solved in  $10 < t < 100$  secs
- 10 % solved in  $t \geq 100$  secs
- 9 % out of memory

No criterion optimized

No enumeration of all solutions



Intersections of seventh and eighth degree polynomial curves

## GloptiPoly: summary

GloptiPoly is a **general-purpose** software with a **user-friendly** interface

Pedagogical flavor, black-box approach, no expert tuning required to cope with **very distinct** applied maths and engineering pbs

**Not a competitor** to highly specialized codes for solving polynomial systems of equations or large combinatorial optimization pbs

**Numerical conditioning** (Chebyshev basis) deserves further study

See also the **SOSTOOLS** software

[www.cds.caltech.edu/sostools](http://www.cds.caltech.edu/sostools)