#### Computational Issues in Nonlinear Dynamics and Control

Arthur J. Krener

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Supported by AFOSR and NSF

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- Numerical Solution of Optimal Control Problems

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Hamilton Jacobi Bellman Equations

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#### Stabilization by Optimization

Problem: Find a feedback  $u = \kappa(x)$  so that the closed loop system is (locally) asymptotically stable around x = 0.

Solution: Choose a Lagrangian  $l(x, u) \ge 0$  and solve the infinite horizon optimal control problem. Under suitable conditions the optimal feedback  $u = \kappa(x)$  is stabilizing on some domain around x = 0 and this can be verified because the optimal cost  $\pi(x) \ge 0$  is a Lyapunov function,

$$rac{d}{dt}\pi(x(t))=rac{\partial\pi}{\partial x}(x(t))f(x(t),\kappa(x(t)))=-l(x(t),\kappa(x(t)))\leq 0$$

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$$0 = F'P + PF + Q - PGR^{-1}G'P, \quad K = -R^{-1}G'P$$

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Theorem: If  $Q \ge 0$ , R > 0, (F, G) stabilizable and  $(Q^{1/2}, F)$  detectable then there exist a unique nonnegative definite solution P to the Riccati equation and the feedback u = Kx is asymptotically stable, i.e., all the poles of F + GK are in the open left half plane.

$$\begin{split} \min_{u(0:T)} \int_{0}^{T} l(t,x,u) \ dt + \pi^{T}(x(T)) \\ \dot{x} &= f(t,x,u) \\ 0 &= g(x(0),X(T)) \\ u(t) &\in \mathcal{U}(t,x) \end{split}$$

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#### **Pontryagin Maximum Principle:**

If  $x^*(0:T), \ u^*(0:T)$  is optimal then there exists  $p:[0,T] \to I\!\!R^{1 imes n}$  such that

$$\begin{split} \dot{x}_i^* &= \ \frac{\partial \mathcal{H}}{\partial p_i}(t,p,x^*,u^*) \\ \dot{p}_i &= \ -\frac{\partial \mathcal{H}}{\partial x_i}(t,p,x^*,u^*) \\ u^* &= \ \operatorname{argmin}_{u \in \mathcal{U}(T,x^*)} \mathcal{H}(t,p,x^*,u^*) \\ t,p,x,u) &= \ pf(t,x,u) + l(t,x,u) \end{split}$$

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Commutative Diagrams?

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 $\begin{array}{cccc} \mbox{Finite Horizon OCP} & \rightarrow & \mbox{Nonlinear Program} \\ & \downarrow & & \downarrow \\ \mbox{PMP} & \rightarrow & \mbox{Karush Kuhn Tucker Conditions} \end{array}$ 

For simplicity of exposition assume n = 2, m = 1. Choose a rectangle around x = 0 and partition it with stepsize h. Let  $x^{i,j}$  denote the i, j node. Let  $\pi^{i,j}$  be the current computed approximation to the optimal cost at the  $x^{i,j}$ .

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For each i, j solve for the next approximation  $\kappa^{i,j}$  to the optimal feedback

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m argmin}_u \left\{ \left( \pi^{i+1,j} - \pi^{i-1,j}, \pi^{i,j+1} - \pi^{i,j-1} 
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The next approximation to the optimal cost  $\bar{\pi}^{i,j}$  is the solution to

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This is called **policy iteration** and it is not very efficient because it sweeps through the nodes many times. Discretization of the Optimal Contol Problem Instead we can discretize the optimal control problem. Asumme that the control takes on discrete values  $u_k$  and time is measured in steps of h. Discretization of the Optimal Contol Problem Instead we can discretize the optimal control problem. Asumme that the control takes on discrete values  $u_k$  and time is measured in steps of h.

Define  $ar{f}(x^{i,j},u^k)$  to be the state node closest to

$$x^{i,j} + f(x^{i,j}, u^k)h$$

Then on the state and control grids we have the discrete dynamics

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and we minimize by choice of control sequence

$$egin{array}{rll} \pi^{i,j}&=&\min_{u(0:\infty)}\sum\limits_{t=0:h:\infty}l(x,u)h\ \kappa^{i,j}&=&u^*(0) \end{array}$$

Discretization of the Optimal Contol Problem **Dynamic Programming Equation (DPE)** 

$$\begin{array}{lll} \pi^{i,j} & = & \min_{u^k} \left\{ l(x^{i,j}, u^k) h + \pi(\bar{f}(x^{i,j}, u^k)) \right\} \\ \kappa^{i,j} & = & \operatorname{argmin}_{u^k} \left\{ l(x^{i,j}, u^k) h + \pi(\bar{f}(x^{i,j}, u^k)) \right\} \end{array}$$

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This can be solved by policy iteration. Given the current approximation  $\pi(x^{i,j})$  to the optimal cost at grid points  $x^{i,j}$ , define the next approximation to the optimal feedback as

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This is again policy iteration but it is still slow.

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Partition the state space into a grid with spacing h and partition time with spacing k. Construct a Controlled Markov Chain with transition probability  $p(x^1|x^0, u)$  from gridpoint  $x^0$  to grid point  $x^1$  with control u. Choose a search radius r and define

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The cost is defined to be the expected value of

$$\sum_{t=0}^{\infty} l(x,u)$$

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But in higher dimensions it difficult to implement.

Suppose that the speed of propagation c(x) > 0 through a medium varies with location. Consider any path x(t) between the source  $x^0 = 0$  and  $x^1$ . Then the propagation time along this path is  $\int_0^t 1 d\tau$  so the Lagrangian l(x, u) = 1.

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Fast Marching Method for the Eikonal Equation This method is due to Tsitsiklis and it was refined by Sethian, Falcone and others. Fast Marching Method for the Eikonal Equation This method is due to Tsitsiklis and it was refined by Sethian, Falcone and others.

Partitition the nodes into three families called accepted, narrow band and far. Initially the only node in accepted family is the origin and  $\pi^{i_0,j_0} = 0$ .



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The FFM has been generalized to other optimal control problems but computing the minimum sums is more complicated because not every rectilinear path is feasible. Curse of Dimensionality
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Practical optimal control problems usually have state dimension larger than 2 or 3. For example, the attitude control problem for a spacecraft has state dimension n = 6 and control dimension typically m = 3. The position and attitude control problem for an airplane has state dimension n = 12 and control dimension at least m = 4.

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Consider trying to apply a grid based method.. For the solution to be reasonably accurate we would need a substantial number of grid points in each coordinate direction, e.g.,  $10^2$ . Then the total number of grid points is  $10^{12}$  for attitude control and  $10^{24}$  for position and attitude control. If we can process 100 nodes a second that works out to about 300 years for attitude control and  $3 \cdot 10^{14}$  years for position and attitude control.

HJB Equations and Conservation Laws Suppose we have a time varying problem of the form

$$egin{array}{rcl} \dot{x} &=& f(t,x) + g(t,x) u \ l(t,x,u) &=& q(t,x) + rac{1}{2} u' R(t,x) u \end{array}$$

Then the HJB PDEs reduce to the HJ PDE

$$0 \;\; = \;\; rac{\partial \pi}{\partial t} + rac{\partial \pi}{\partial x} f - rac{1}{2} rac{\partial \pi}{\partial x} g R g' \left(rac{\partial \pi}{\partial x}
ight)' + q$$

Let  $p = \frac{\partial \pi}{\partial x}$  and take the Jacobian of the HJ equation to obtain the conservation law

$$0 \;\; = \;\; rac{\partial p}{\partial t} + rac{\partial}{\partial x}F(t,x,p)$$

where the flux term is

$$F(t,x,p) = pf - rac{1}{2}pgRg'p' + q$$

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Another is that we are looking for a solution of the conservation law that is a closed one form,

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So we can compute  $\pi(x)$  by computing the stable manifold of the Hamiltonian dynamics.

### Hauser-Osinga Method

The Geometry of Optimal Control



Figure 1: Sketch of the balanced planar pendulum on a moving cart.

John Hauser & Hinke Osinga



Figure 2: Each point is colored according to how high the cost of getting to the origin using this point as initial condition. The cost increases as the color changes from blue to red.

10

### Min-Plus Methods

The min-plus semiring is defined as follows

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There is also a max-plus semiring where

$$\xi \oplus \zeta = \max\{\xi, \zeta\}$$

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$$egin{array}{rcl} \mathcal{S}_{T}(\phi)(x^{T}) &=& \min_{u(-T:0)} \left\{ \int_{-T}^{0} l(x(t),u(t)) \; dt + \phi(x(0) 
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But is linear in the min-plus sense and  $\pi(x)$  is an eigenvector corresponding to the eigenvalue 0 which is the  $\otimes$  identity.

$$0\otimes \pi(x) ~=~ \mathcal{S}_T(\pi(x))$$

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To make the calculation finite dimensional  $\pi(x)$  is chosen as a min-plus combination of basis functions.

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The number of basis functions needed for a given accuracy is exponential in the state dimension n but it is probably grows slower than the number of grid points.

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It also suffers from a curse of complexity as it requires computing the pointwise maxima (or minima) of a large number of functions which can be expensive.

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If  $h_1 = 0.01$  then  $h_3 \approx 0.1$  so the number of grid points that is needed for a given accuracy is reduced by a factor of 10 in each dimension. If the state dimension is n then the reduction in grid points is by a factor of  $10^n$ .

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If the third order method takes  $k_3(n)$  times longer to compute for each node then the reduction in computational time is by the factor  $\frac{10^n}{k_3(n)}$ . Typically  $k_3(n)$  is polynomial in n.

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There are diminishing returns as we go to higher orders. Consider a fifth order method with step size  $h_5$ . Then for same level of accuracy

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 so  $h_5pprox h_1^{1/3}$   
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Multiply the first by 4/3 and the second by -1/3 and add,

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Szpiro and Dupuis have applied this technique to HJB equations.

#### A first order quasilinear PDE

$$0 \;\; = \;\; rac{\partial \phi}{\partial x}(x) a(x) + b(x,\phi(x))$$

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$$\begin{array}{rcl} a(x) &=& Ax + a^{[2]}(x) + a^{[3]}(x) + \dots \\ b(x,\phi(x)) &=& Cx + B\phi(x) + (b(x,\phi(x)))^{[2]} + (b(x,\phi(x)))^{[3]} + \dots \\ \phi(x) &=& Tx + \phi^{[2]}(x) + \phi^{[3]}(x) + \dots \end{array}$$

where  $(\cdot)^{[d]}$  denotes terms homogeneous of degree d.

Collect terms of first degree.

TA + BT = -C

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The higher degrees terms can be found in a similar fashion.

Many of the most important PDEs of nonlinear dynamics and control are essentially singular first order quasilinear including

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Discrete time and time varying problems can also be solved by similar power series methods.

In discrete time the degree two nonresonance conditions are

$$\alpha_{i_1}\alpha_{i_2} \neq \beta_j$$

Al'brecht developed the power series method for HJB equations for the optimal cost and optimal feedback,

$$\pi(x) = \frac{1}{2}x'Px + \pi^{[3]}(x) + \pi^{[4]}(x) + \dots$$
  

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This guarantees there are no resonances so the higher degree terms of  $\pi$ ,  $\kappa$  can be found by solving invertible linear equations.

This method has been implimented in the Nonlinear Systems Toolbox.

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- The HJB equations can be solved to degree 4 in  $\pi(x)$  and degree 3 in  $\kappa(x)$  for systems with state dimension n = 25 and control dimension m = 8 on a lap top.

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#### Patchy Methods



Figure : Optimal Cost of Inverting a Pendulum by a Torque at its Axis

# Sequence of Patches



Figure : Sequence of Patches

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Under suitable assumptions there is one positive root and one negative root. We take the positive root.

#### Invert a Pendulum



Figure : Periodicity of the Optimal Cost

The left axis is  $-15 \le \dot{\theta} \le 15$  and the right axis is  $-15 \le \theta \le 15$ . From points on the ridges there are two optimal trajectories, one going to the left well and the other going to the right well.

#### Adaptive Algorithm

The algorithm is adaptive. It splits a patch in two when the relative residue of the first HJB equation is too high at the lower corners of a patch. It also lowers the upper level of a ring of patches if the relative residue is too high on it.

Ring	1	2	3	4
Initial Patch Level	0.64	1.21	1.96	2.89
Final Patch Level	0.36	0.63	1.38	2.23
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The last ring (34) contains 78 patches.



Figure : Patchy Optimal Cost to Level Set 70



Figure : Patchy Optimal Cost to Level Set 140



Figure : Patchy Optimal Cost to Level Set 210



Figure : Patchy Optimal Cost to Level Set 280



Figure : Patchy Optimal Cost to Level Set 350



Figure : Patchy Optimal Cost to Level Set 420



Figure : Patchy Optimal Cost to Level Set 490



Figure : Patchy Optimal Cost to Level Set 566

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	Max Error	Max Rel Error	Error Factor
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This shows that the patchy method can be very accurate and it is parallelizable.

#### Three Dimensional Example Here is a level set of the patchy method applied to a three dimensional problem



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The complexity of keeping track of the patches probably makes the patchy method infeasible in higher dimesions.

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Similar statements are true for the other PDEs of nonlinear control.

**Typical Problem** 

$$\min_{u(0:T)}\int_0^T l(x,u) \; dt$$

#### subject to

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$$\begin{split} \dot{x}_i &= \frac{\partial \mathcal{H}}{\partial p_i}(p, x, u^*) \\ \dot{p}_i &= -\frac{\partial \mathcal{H}}{\partial x_i}(p, x, u^*) \\ u^* &= \operatorname{argmin}_u \left\{ \mathcal{H}(p, x, u) : 0 \leq g(x, u) \right\} \end{split}$$

plus boundary and transversality conditions.

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If we discretize time with step size h then decision variables are

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Because of the development of excellent software for solving nonlinear programs, the direct approach has become more popular.

$$egin{aligned} \min_{u(0:T)} & \sum_{t=0:h:T-1} & ar{l}(x,u) \ x^+ &= ar{f}(x,u), & 0 \leq g(x,u) \ x(0) &= x^0, & x(T) = x^T \end{aligned}$$

where the discrete dynamics and discrete Lagrangian are defined by Lie differentiation

$$ar{f}(x,u) = x + f(x,u)h + L_{f(x,u)}f(x,u)rac{h^2}{2} + L_{f(x,u)}^2f(x,u)rac{h^3}{6} + ar{l}(x,u) = l(x,u)h + L_{f(x,u)}l(x,u)rac{h^2}{2} + L_{f(x,u)}^2l(x,u)rac{h^3}{6} + \dots$$

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If Lie differentiation is difficult use Runge-Kutta approximations.

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Discretization of the Optimal Trajectory Problem

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Recall that we have to find the minimum of a function of mT/h variables,  $u(0), u(h), u(2h), \ldots, u(T-1)$ .

The discretization of the continuous time problem is a form a quadrature so we could use any quadrature rule, e.g., Euler, Trapezoidal, Hermite-Simpson, etc. in either explicit or implicit form.

Perhaps the most efficient quadrature is Legendre-Gauss (LG). It uses only N nodes to exactly integrate any polynomial of degree 2N - 1 or less.

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On the standard interval [-1,1] it takes the form

$$\int_{-1}^{1} \phi(t) \; dt \;\; = \;\; \sum_{i=1}^{N} w_i \phi(t_i) \;\;$$

where the nodes  $t_i$  are the zeros of the  $N^{th}$  Legendre polynomial  $P_N(t)$ .

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But all the nodes  $t_i$  are in the open interval (-1, 1) so Legendre-Gauss quadrature is not suitable if there are boundary conditions.

legendre polynomials. 1 0.5 22 D -0.5 Prizz Pi/x Pairs Ps( P<sub>1</sub>/2  $\cdot 1$ Pi(x)  $\cdot 1$ -0.5D 0.5 1  $\mathbf{x}$ 

Figure : Legendre Polynomials to Degree 5

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The weights are  $\frac{2}{n(n-1)}$  at the endpoints and

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in between.

Gong, Kang and Ross have shown that the pseudospectral method converges for feedback linearizable systems. If m = 1 such a system can be transformed to

$$egin{array}{rcl} \dot{x}_1 &=& x_2 \ \dot{x}_2 &=& x_3 \ && dots \ \dot{x}_{n-1} &=& x_n \ \dot{x}_n &=& f_n(x) + g_n(x) u \end{array}$$

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Minimize

$$\int_{-1}^{1} l(x(t), u(t)) \ dt + \alpha(x(-1), x(1))$$

subject to

$$egin{array}{rcl} 0 &=& eta(x(-1),x(1)) \ 0 &\leq& \gamma(x(t),u(t)) \end{array}$$

Each  $x_i(t)$  is approximated by an  $N^{th}$  degree interpolating polynomial  $\bar{x}_i(t)$ . These polynomials are represented by their values at the N + 1 LGL nodes,

$$ar{x}_i = egin{bmatrix} ar{x}_i^0 & \dots & ar{x}_i^N \end{bmatrix}' \ ar{x}_i(t) = \sum_0^N ar{x}_i^j \phi_j(t)$$

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where the  $\phi_j(t)$  are the Lagrange polynomials at the LGL nodes. The dynamics is approximated by the equations

$$egin{array}{rll} ar{x}_{i+1} &=& Dar{x}_i, & i=1,\ldots,n-1 \ ar{u}^j &=& rac{(Dar{x}_n)^j-f(ar{x}^j)}{g(ar{x}^j)} \end{array}$$

so the N+1 decision variables are  $\bar{x}_1^0,\ldots,\bar{x}_1^N$ .

Pseudospectral Trajectory Optimization Multiplication of the interpolated values of a polynomial by the differentiation matrix D yields the interpolating values of its derivative. Pseudospectral Trajectory Optimization Multiplication of the interpolated values of a polynomial by the differentiation matrix D yields the interpolating values of its derivative.

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The cost is approximated by a LGL quadrature

$$\sum_{j=0}^N l(x^j,u^j)w_j+lpha(ar x^0,ar x^N)$$

The boundary conditions are approximated by a relaxed version of

$$0 = \beta(\bar{x}^0, \bar{x}^N)$$

and the constraints are approximated by a relaxed version of

$$0 ~\leq~ \gamma(ar{x}^j,ar{u}^j), ~~ j=0,\ldots,N$$

Pseudospectral Trajectory Optimization and the ISS

SIAM News, Volume 40, Number 7, September 2007

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#### Pseudospectral Optimal Control Theory Makes Debut Flight, Saves NASA \$1M in Under Three Hours

By Wei Kang and Naz Bedrossian



#### Suppose the problem of minimizing

$$\int_0^\infty l(x,u) \,\,dt$$

#### subject to

$$egin{array}{rcl} \dot{x}&=&f(x,u)\ x(0)&=&x^0\ 0&\leq&g(x,u) \end{array}$$

#### has been discretized into minimizing

$$\sum_{t=0:h:\infty}ar{l}(x(t),u(t))$$

subject to

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Minimization over the infinite horizon is too difficult so choose a time window T and a terminal cost  $\pi_T(x)$  defined on a terminal set  $\mathcal{X}_T$  which is a compact neighborhood of x = 0.

Consider the problem of minimizing

$$\sum_{t=0:h:T-h}ar{l}(x(t),u(t))+\pi_T(x(T))$$

subject to

$$egin{array}{rcl} x^+ &=& ar{f}(x,u) \ x(0) &=& x^0 \ 0 &\leq& g(x,u) \ x(T) &\in& \mathcal{X}_T \end{array}$$

The decision variables are  $u(0), \ldots, u(T-h)$ .

Then pass this nonlinear program to a fast solver to find the optimal  $u^0(0), \ldots, u^0(T-h)$ . This needs to be done in less than the time step h.

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Then between times h and 2h solve the problem of minimizing

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to obtain the optimal  $u^1(h),\ldots,u^1(T).$ 

Use the control  $u^1(h)$  to get the state to  $x^2 = x(2h)$ , etc.

The key issues are the following

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- The horizon T must be long enough and/or  $\mathcal{X}_T$  large enough so that  $x(t+T) \in \mathcal{X}_T$ .
- The initial guess of  $u^0(0), \ldots, u^0(T-1)$  that is fed to the solver must be close to optimal else the solver may fail to converge to the true solution.

• This is not as much a problem with later initial guesses because we can take  $u^0(h), \ldots, u^0(T-h)$  as the initial guess for  $u^1(h), \ldots, u^1(T-h)$ .

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- If the infinite horizon optimal control law  $\kappa_T(x)$  is known on the terminal set  $\mathcal{X}_T$  then the initial guess for  $u^1(T)$  should be  $\kappa_T(x^0(T))$  where  $\bar{x}^0(T)$  is the  $T^{th}$  state generated by the last control sequence.  $u^0(0), \ldots, u^0(T-1)$

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- It may be possible to use power series methods to compute the terminal cost  $\pi_T(x)$  and feedback  $\kappa_T(x)$  on a larger terminal set  $\mathcal{X}_T$ . This may allow us to lengthen the time step h and/or shorten the horizon T so that MPC can be used on faster processes.

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- Thank you! Questions?

- M. Bardi, I. Caputo-Dolcetta, Optimal Control and Viscosity Solutions of Hamilton-Jacobi-Bellman Equations, Birkhauser, 1997.
- M. Falcone, Numerical Solution of Dynamic Programming Equations, Appendix A, Optimal Control and Viscosity Solutions of Hamilton-Jacobi-Bellman Equations, Birkhauser, 1997.
- M. Boue and P. Dupuis, Markov Chain Approximations for Deterministic Control Problems with Affine Dynamics and Quadratic Cost in the Control, SIAM J. Numerical Analysis, v. 36, pp. 667-695, 1999.
- J. N. Tsitsiklis, Efficient Algorithms for Globally Optimal Trajectories, IEEE Transactions on Automatic Control, v. 40, pp. 1528-1538, 1995.

- J. A. Sethian, A fast marching level set method for monotonically advancing fronts, Proc.Natl. Acad. Sci. USA, 93 (1996), 1591-1595.
- J. A. Sethian and A. Vladimirsky, Ordered upwind methods for static Hamilton-Jacobi equations: theory and algorithms, SIAM J. Numer. Anal., 41 (2003), 325-363.
  - E. Carlini, M. Falcone, N. Forcadel, and R. Monneau, Conver- gence of a Generalized Fast Marching Method for an Eikonal equation with a velocity changing sign, SIAM J. Numer. Anal. v. 46, pp. 2920-2952, 2008.
- **J**. Hauser and H. Osinga On the Geometry of Optimal Control: the Inverted Pendulum Example, in Proceedings of the American Control Conference, Arlington VA, USA, Volume 2, pp. 17211726, 2001.

- M. Akian, S. Gaubert and A. Lakhoua, A max-plus finite element method for solving finite horizon determinsitic optimal control problems, Proc. 16th International Symposium on Mathematical Theory of Networks and Systems (MTNS), Leuven, 2004.
- W. McEneaney, A Curse of Dimensionality Free Method for the Solution of Certain HJB PDEs, SIAM J. Control and Opt. (2007).
- A. Szpiro and P. Dupuis, A Second Order Numerical Method for First Order Hamilton-Jacobi Equations, SIAM J. Numer. Anal., v. 40, pp.1136-1183, 2002.

- **E. G. Al'brekht, On the Optimal Stabilization of Nonlinear** Systems, J. Appl. Math. Mech., v. 25, pp. 1254-1266, 1961.
- C. Navasca and A. J. Krener, Patchy Solutions of Hamilton-Jacobi-Bellman Partial Differential Equations, in A. Chiuso et al. (eds.), Modeling, Estimation and Control, Lecture Notesin Control and Information Sciences, 364 (2007), 251-270.
- **T**. Hunt, A proof of the higher order accuracy of the patchy method for solving the Hamilton-Jacobi-Bellman equation, Ph.D. Thesis, University of California, Davis, 2011.
- S. Cacace, E. Cristiani, M. Falcone and A. Picarelli, A patchy dynamic pro- gramming scheme for a class of Hamilton-Jacobi-Bellman equations, accepted by SIAM Journal on Scientic Computing.

- C. O. Aguilar, C. O., T. W. Hunt and A. J. Krener, An Adaptive Patchy Method for Solving Hamilton Jacobi Bellman Equations, Proceedings of the International Conference of Numerical Analysis and Applied Mathematics, American Institute of Physics Conference Proceedings no. 1479, 2012.
- C. O. Aguilar and A. J. Krener, Patchy Solution of a Francis-Byrnes-Isidori Partial Diffential Equation, International Journal of Robust and Nonlinear Control,
- B. Krauskopf, H.M. Osinga, E.J. Doedel, M.E. Henderson, J. Guckenheimer, A. Vladimirsky, M. Dellnitz and O. Junge, A survey of methods for computing (un)stable manifolds of vector fields, Int. J. Bifurcation and Chaos, v. 15, pp. 763-791, 2005.

- M. Gerdts, Numerical Optimal Control, OMPC 2013 -SADCO Summer School and Workshop on Optimal and Model Predictive Control. Slides available at http://num.math.uni-bayreuth.de/en/conferences/
- **F.** Fahroo and I.M. Ross. Direct Trajectory Optimization by a Chebyshev Pseudospectral Method. Journal of Guidance Control and Dynamics, 25, 2002.
- Q. Gong, W. Kang and I. M. Ross, A Pseudospectral Method for the Optimal Control of Constrained Feedback Linearizable Systems, IEEE Trans. on Automatic Control, v. 51, pp.1115-1129, 2006
- D. Q. Mayne, J. B. Rawlings, C. V. Rao and P. O. M. Scokaert Constrained model predictive control: Stability and optimality. Automatica, v. 36, pp.789-814, 2000.

- J. B. Rawlings and D. Q. Mayne. Model Predictive Control: Theory and Design. Nob Hill Publishing, Madison, WI, 2009.
- J. B. Rawlings, Model Predictve Control, OMPC 2013 -SADCO Summer School and Workshop on Optimal and Model Predictive Control. Slides available at http://num.math.uni-bayreuth.de/en/conferences/