Numerical Methods for Optimal Control Problems. Part II: Local Single-Pass Methods for Stationary HJ Equations

Ph.D. course in OPTIMAL CONTROL



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- Hamilton-Jacobi equations for MTP
- Semi-Lagrangian discretization
- Classical iterative method
- Local single pass methods
- Fast marching method
- Can local single pass methods solve every HJ equation?

Hamilton-Jacobi equations

Hamilton-Jacobi equations arise in several applied contexts, e.g. *front propagation, control problems and differential games.*

Eikonal equation

$$\left\{egin{array}{c} |
abla v(x)| = 1 & x \in \mathbb{R}^d \setminus \mathcal{T} \ v(x) = 0 & x \in \partial \mathcal{T} \end{array}
ight.$$

The solution v represents the *distance* function from ∂T and it is well understood in the framework of viscosity solutions ¹.



Solution to the Eikonal equation in dimension d = 2 with $T = {\text{five random points}}.$

M.G. Crandall, P.-L. Lions, Viscosity solutions of Hamilton-Jacobi equations, Trans. Amer. Math. Soc., 277 (1983), 1–42.

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Semi-Lagrangian discretization of the HJB equation

Let G be a structured grid with nodes x_i , i = 1, ..., N and space step Δx .

SL discretization of the HJB equation

$$w(x_i) = \min_{a \in A} \left\{ w(\tilde{x}_{i,a}) + \frac{|x_i - \tilde{x}_{i,a}|}{|f(x_i, a)|} \right\}, \qquad x_i \in G$$

where $\tilde{x}_{i,a}$ is a *non-mesh* point, obtained by integrating, until a certain final time \hat{s} , the ODE

$$\left\{ egin{array}{ll} \dot{y}(s)=f(y,a), & s\in [0,\hat{s}] \ y(0)=x_i \end{array}
ight.$$

and then setting $\tilde{x}_{i,a} = y(\hat{s})$. To make the scheme fully discrete, the set of admissible controls A is discretized in N_c points.

We get different versions of the SL scheme varying \hat{s} , the method used to solve the ODE and the interpolation method used to compute $w(\tilde{x}_{i,a})$.

Semi-Lagrangian discretization of the HJB equation

Explicit forward Euler scheme for the ODE + linear interpolation



2pSL: $\tilde{x}_{i,a}$ intercepts the line connecting $x_{i,1}$ and $x_{i,2}$.

3pSL: $\tilde{x}_{i,a}$ is at distance Δx from x_i .

Stationary Hamilton-Jacobi equations

Equations we are interested in can be recast as *minimum time problems*. By choosing the set of admissible controls $A = B_1(0)$ we get the following

Reference equations HJ equation Name f(x, a) $|\nabla T(x)| = 1$ а homogeneous eikonal $c_1(x)a$ $c_1(x)|\nabla T(x)| = 1$ nonhomogeneous eikonal $c_2\left(\frac{\nabla T}{|\nabla T|}\right)$ $|\nabla T(x)| = 1$ $c_2(a)a$ hom. anisotropic eikonal $c_3\left(x, \frac{\nabla T}{|\nabla T|}\right) |\nabla T(x)| = 1$ $c_3(x, a)a$ nonhom. anisotropic eikonal

The functions c_1 , c_2 , c_3 are strictly positive and Lipschitz continuous.

Classical iterative method

How to solve the nonlinear system?

$$w(x_i) = S[w](x_i) := \min_{a \in A} \left\{ w(\tilde{x}_{i,a}) + \frac{|x_i - \tilde{x}_{i,a}|}{|f(x_i, a)|} \right\}, \qquad x_i \in G$$

Fixed point algorithm

Given an initial guess $w^{(0)}$ iterate on the grid G $w^{(k)} = S[w^{(k-1)}]$ k = 1, 2, 3, ...until $\max_{x_i \in G} |w^{(k)}(x_i) - w^{(k-1)}(x_i)| < \varepsilon$

Classical iterative method

Pros

- * Numerical approximation of *viscosity solution* in any dimension for any *f*.
- * Easy implementation.
- * Easy parallelization.
- * A priori error estimates in L^{∞} .
- * Structured or unstructured grids.

Cons

* "Curse of dimensionality" (exponentially increasingly nonlinear systems for high dimensional problems)

- \Rightarrow huge computational efforts
- \Rightarrow huge memory resources.

Causality as source of efficiency

At the continuous level, information emanates from the target set \mathcal{T} and propagates along characteristic lines.



By mimicking this behavior at the discrete level, one can produce a reordering of the grid nodes that decouples the nonlinear system.

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Local Single Pass algorithms

Causality

Exploit physical/geometric properties of the HJ equations to find an ordering of the grid nodes that avoid useless computations.

Locality

The computation is dynamically localized on the grid nodes carrying relevant information (few, compared to the entire grid). Each node is computed using only neighboring nodes.

Single Pass property

Each node is re-computed at most r times, where r only depends on the equation and the grid structure, not on the number of grid nodes.

Inspired by Dijkstra's algorithm³ for the *shortest path problem* on a graph, FMM (by Tsitsiklis⁴ and Sethian⁵) is a *local single pass* method for the *Eikonal equation*.



●Accepted ●Considered ○Far

FMM algorithm

Set T = 0 in ACC and $T = +\infty$ in FAR Compute T in CONS While(CONS $\neq \emptyset$) Find $\bar{x} = \underset{x \in CONS}{\operatorname{arg min}} T(x)$ Move \bar{x} from CONS to ACC Move !ACC neighbors of \bar{x} in CONS (if yet in) and (re)compute T on them

- ³E. W. Dijkstra, A note on two problems in connexion with graphs, 1959.
- ⁴ J. N. Tsitsiklis, Efficient algorithms for globally optimal trajectories, 1995.
- ⁵ J. A. Sethian, A fast marching level set method for monotonically advancing fronts, PNAS USA, 1996.

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Why FMM works?

FMM computes each node in CONS by means of nodes with smaller values (practical implementations enforce the use of nodes in ACC only!).

The solution is computed in *ascending order*, so that the node in CONS with *minimal value* is the only *not influenced* by other nodes in CONS.

The minimal value rule corresponds to get information from the simplex containing $-\nabla T$ (and implies that CONS approximately expands as a *level set* of T).

For the Eikonal equation, *characteristic lines coincide with gradient lines* of the solution itself, hence FMM computes the *correct* solution.

For general HJ equations this is not true!

FMM's FAILURE: Anisotropic Eikonal equation



Wide divergence between characteristic and gradient lines!

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Beyond FMM

Several directions of research: high order accuracy, smart implementations, different schemes (FD, SL, DG, FV), other competitive approaches (FS, FI, MaxPlus), hybrid methods, more general HJ equations.

Some references

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Let us classify HJ equations in two classes:

(EIK) Eikonal-like equations, whose characteristic lines coincide or lie in the same simplex of the gradient lines of their solutions.

 $(\neg EIK)$ Non Eikonal-like equations, for which there exists at least a grid node where the characteristic line and the gradient of the solution do not lie in the same simplex.

By construction FMM works for equations of type EIK and fails for equations of type \neg EIK (e.g. the Anisotropic Eikonal equation).

Is the minimal value rule really needed?

In order to solve \neg EIK equations, CONS *cannot be* at any time an approximation of a *level set*, i.e. we have to *drop* the minimal value rule.

We consider another classification:

(DIFF) Equations with smooth characteristics. Information spreads from the target \mathcal{T} to the rest of the space along smooth lines, without shocks. The solution \mathcal{T} is differentiable.

(\neg DIFF) Equations with non smooth characteristics. Information starts from the target \mathcal{T} and then crashes, creating shocks. The solution \mathcal{T} is Lipschitz continuous.



$$T(x_i) = \min_{a \in A} \left\{ T(\tilde{x}_{i,a}) + \frac{|x_i - \tilde{x}_{i,a}|}{|f(x_i, a)|} \right\}, \qquad x_i \in G$$

Safeness

A node $x_i \in \text{CONS}$ is said to be *safe* if $T(x_i)$ is computed using values at neighboring interpolation points which are in ACC only.



Warning! Safeness makes sense if nodes in CONS can be computed using nodes both in ACC and CONS.

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Safe "Method" (SM)

At each step, every *safe* node in CONS enters ACC.

SM *can* solve DIFF equations (both EIK and \neg EIK), it is much faster than FMM (multiple node acceptance, no search of min value in CONS).

SM fails for equations of type \neg DIFF.



EIK&¬DIFF: FMM works

M works EIK&¬DIFF: SM fails

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How to handle the shocks?

As in the continuous case, a grid node Δx -close to a shock has to be approached by the ACC region approximately *at the same time* from the directions corresponding to the characteristic lines.

This property is satisfied by FMM in the case EIK, since CONS is approximately a *level set*.



Requiem for Local Single Pass methods?

¬EIK&¬DIFF equations are very hard (if not impossible) to solve

 \neg EIK requires CONS not to be a level set, whereas CONS \sim level set seems the only way to handle shocks in \neg DIFF.



A shock crossing a region with strong anisotropy. What to do?

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