

General Framework

- ① Initialization:  $G(V, E)$  as (undirected) search graph, where  $E$  is empty list of edges and  $V$  is a non-empty set of pts in  $C_{free}$ .

$V = \{q_I, q_G, \text{ and possibly some other points}\}$

Plays a role similar to priority queue in Section 2.2

- ② Vertex Selection Method (VSM):

Choose element of  $V$  for expansion;  $q_{cur} \in V$

- ③ Local Planning Method (LPM):

Main difference to search from Ch 2. Action is replaced by a path.

For some  $q_{new} \in C_{free}$  (possibly not in  $V$ ) attempt to construct a path  $\tau_s: [0, 1] \rightarrow C_{free}$  such that  $\tau(0) = q_{cur}$  and  $\tau(1) = q_{new}$ .

If connection fails (due to collision) go to step 2.

- ④ Insert edge in  $G$ : Insert  $\tau_s$  into  $G$  connecting  $q_{cur}$  to  $q_{new}$ . If  $q_{new} \notin V$ , insert it.

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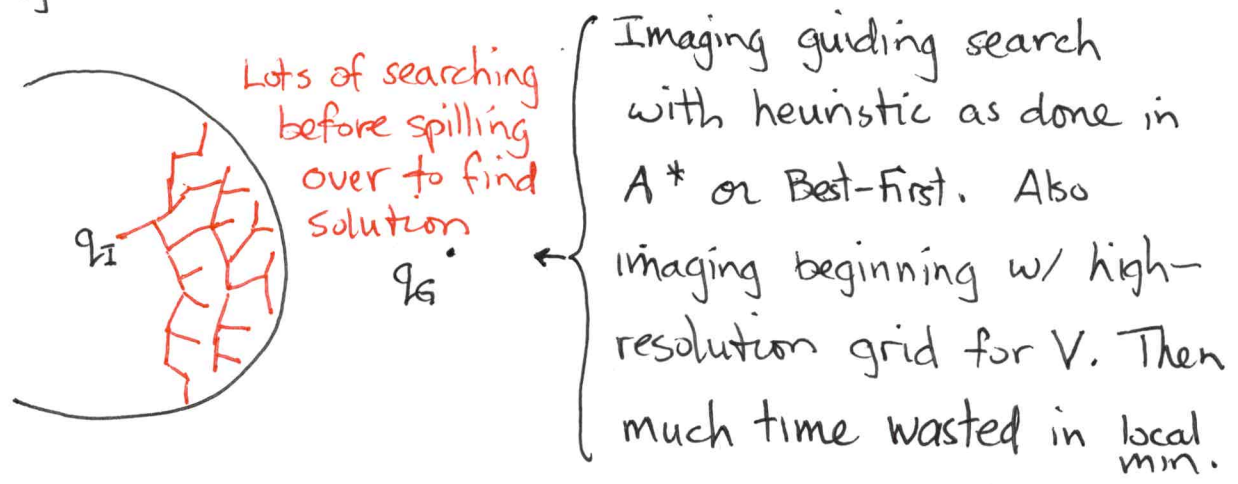
- ⑤ Check for Solution: Determine whether  $G$  encodes solution path. If  $G$  is a tree, this part is trivial. Otherwise not trivial. ②
- ⑥ Return to step 2: If ~~no~~ solution found in ⑤ or stopping criterion has been met, stop. Otherwise return to step ②.

Recall,  $G$  is a topological graph, i.e., every  $q \in V$  is in  $C_{free}$  and every  $e \in E$  is in  $C_{free}$ .

LPM is "local" planning method since it only operates close to  $q_{cur}$ .

Many possible planners can be created by pairing different methods for VSM and LPM

### Avoiding Local Minima



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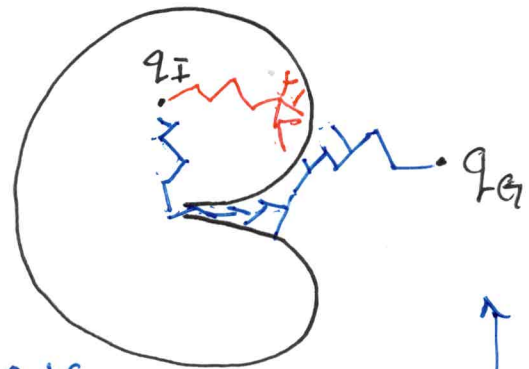
Classes of algorithms base on  
number of search trees

Unidirectional (single-tree) methods :

Very similar to discrete search methods in  
Chapter 2.

Could possibly  
be easy to  
solve with  
a single tree

Difficult to  
solve w/  
fwd search  
tree.



searching backward, if greedy search  
pushes search ~~to~~ toward gap.

Bidirection (two trees) :

Idea is that search frontiers of two trees will  
meet with less exploration than would occur  
when using one tree.

VSM alternates between the trees to choose  $q_{cur}$ .

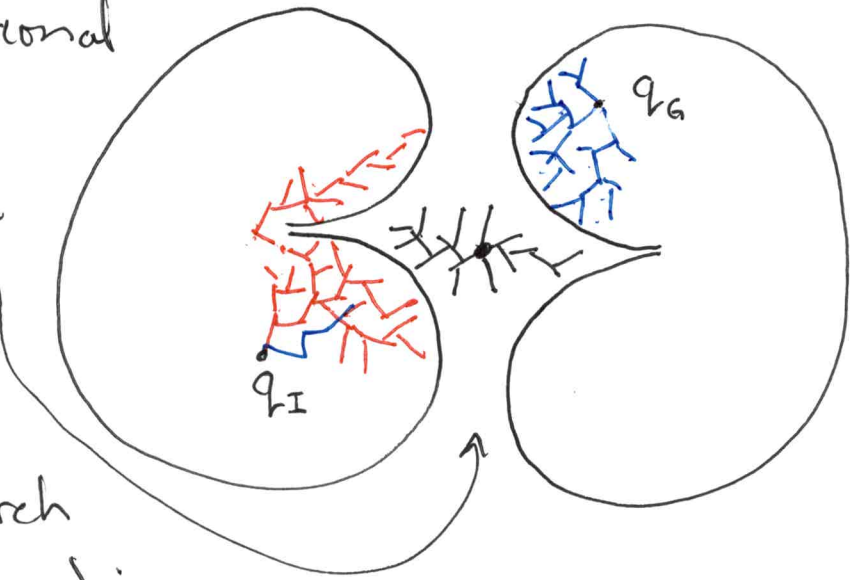
LPM " " connecting the trees and  
growing the trees by choosing  $q_{new} \in$  other  
tree or  $q_{new}$  not in a tree.

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Trouble for bi-directional search.

Maybe an extra tree can help.



Multi-directional search  
(more than two trees):

- More trees could help find paths into traps, but interleaving search and tree connection becomes more complicated.

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One has to accept that some problems just won't be solvable by sampling-based methods unless some structural info about C-space is available!

## 5.4.2 Adapting Discrete Search Algs

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Map all joint motions & other dof's onto  $[0,1]^n / \sim$

$$q_{i,\min} \leq q_i \leq q_{i,\max} \quad \forall i = 1, \dots, n$$

where  $n = \#$  of dof.

Discretization: grid on  $[0,1]^n / \sim$

$$\Delta q_i = [0 \dots 0 \ 1/k_i \ 0 \dots 0]$$

where  $k_i$  is # of grid cells in direction  $i$ .

Grid points are expressed as  $\sum_{i=1}^n j_i \Delta q_i = q$

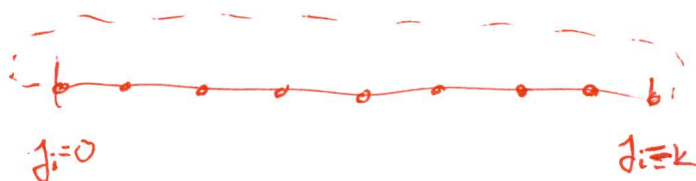
where  $j_i \in \{0, 1, \dots, k_i\}$  so that

$j_1, j_2, \dots, j_n$  are array indices

of the grid point,  $q$ .

For dimensions that are identified at  $0 \sim 1$ ,

then points with  $j_i = 0$  and  $j_i = k_i$  are equal.



identifications  
where needed  
revolute jnts  
that rotate  $2\pi$   
quaternions  
 $h = -h$

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(6)

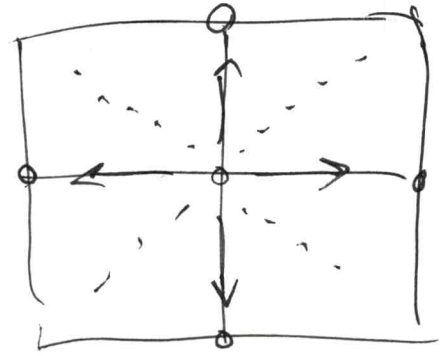
# Neighborhood

The 1-neighborhood is:

$$N_1(q) = \{q \pm \Delta q_1, \dots, q \pm \Delta q_n\}$$

There are at most  $2n$  valid ~~neighbors~~ <sup>1-neighbors</sup>.

Using 1-neighbors keeps # of neighbors to check in search alg. linear in dimension of prob.

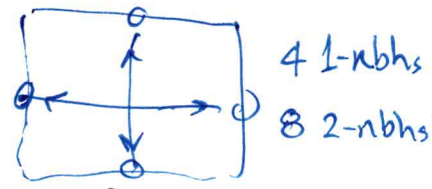


If neighbors included diagonal neighbors, then # of neighbors increases

~~which~~ ~~(exponentially)~~ ~~(quadratically)~~ quadratically with dimension.



The 2-neighborhood is:



$$N_2(q) = \{q \pm \Delta q_i \pm \Delta q_j \mid 1 \leq i, j \leq n, i \neq j\} \cup N_1(q)$$

LaValle says  $q$  has  $3^n - 1$  neighbors in an  $n$ -thbhd LaValle eq. (5.38)

→ How many 2-neighbors?

$q \pm \Delta q_i \pm \Delta q_j \rightarrow 4$ . How many ways to choose  $i \neq j$ ?  $n(n-1)$

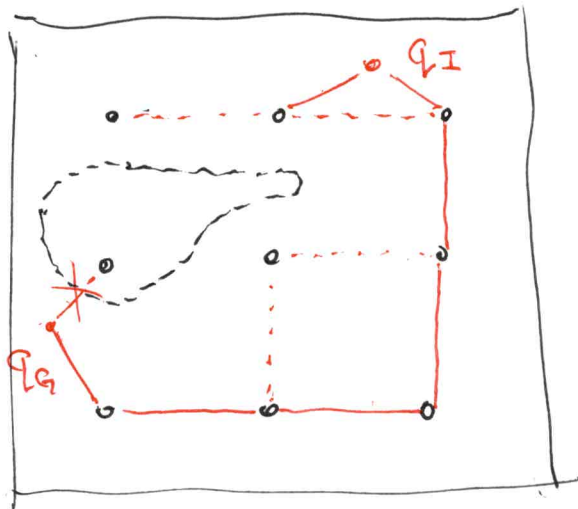
$$\text{Total \#2neighbors} = 4(n(n-1)) = \mathcal{O}(n^2).$$

# Obtaining a discrete planning problem

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- Grid defines  $V$
- Use LPM to connect  $q_I \neq q_G$  to nearest points on grid
- Not all edges need to be explored (necessarily)  $\rightarrow$  RDT or RRT algorithms (section 5.5)
- All vertices  $\neq$  edges could be explored in advance to yield a roadmap (section 5.6).



## Grid resolution issues

• Too much grid resolution slows planning

• Too little grid resolution misses solutions.

Two ways to deal w/ this problem:

- Interleave search  $\neq$  grid refinement

- Abandon grid/discrete search ~~with~~ using algs designed for the continuous problem

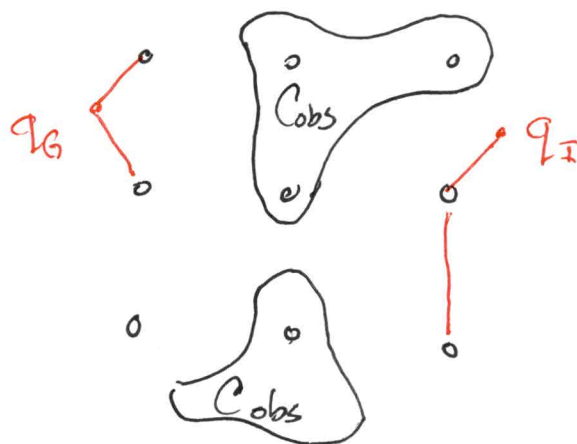
# grid Interleaved search and refinement

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- Start with 2 points per dimension

- search
- double resolution



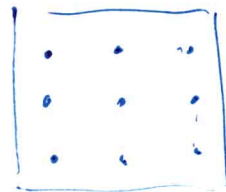
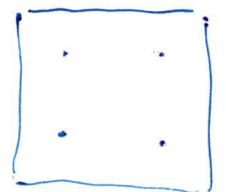
Problem is that # pts grows too rapidly if problem dimension is high,  $\mathcal{O}(2^n)$  where  $n$  is dimension of C-space.

resolution  
dimension  
 $\mathcal{O}(2^n)$

Better approach would be like iterative deepening -

use  $2^n$  pts, search, toss results  
then  $3^n$  pts, search, toss,  $4^n$  pts, ...

Amount of work ~~lost~~ is negligible for large  $n$ .



...

Possibly better. Increase resolution in one dimension at a time (set of nearest neighbors changes), then search

Maybe best. Keep track of connected components of  $\mathcal{G}$  w/ addition of every point. Search for path only when  $q_I$  &  $q_L$  in same component of  $\mathcal{G}$ .



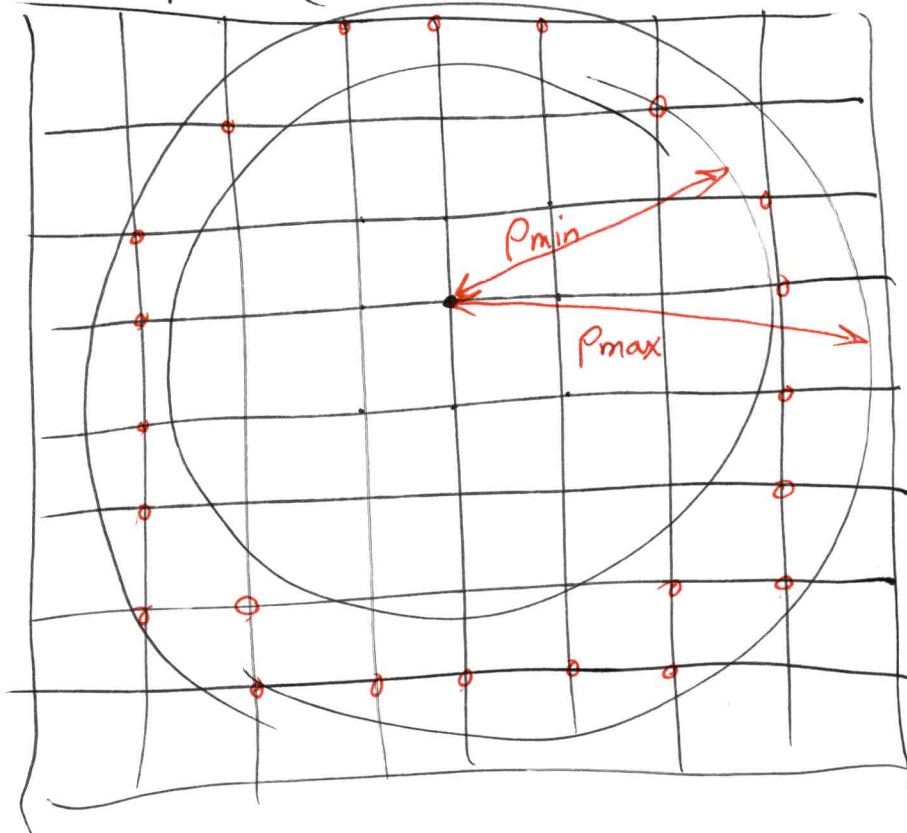
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⑨

Previous grid <sup>discussion</sup> approach tries to keep  $|V|$  small.

Opposite approach.

Declare outrageous resolution (100") at start & then  
hope search <sup>(eg, A\*, Best-First)</sup> has good heuristic to avoid visiting  
many points!

One possible approach, LPM selects "neighbor" far  
from  $q_{cur}$ . (I've not seen this done.)



Perhaps choose  
neighbors such  
that  
 $P_{min} < \rho(q_{cur}, q_{new})$   
 $\neq$   
 $\rho(q_{cur}, q_{new}) < P_{max}$

Could do something like simulated annealing.

$P_{min} \neq P_{max}$  large.

If solution progress slows, make  $P_{min} \neq P_{max}$  smaller.

Could also use large  $\Delta$ 's in one-neighbor selection.

Best-first  $\neq$   $A^*$  could solve w/o visiting many points  
and  $\therefore$  might not have to create a large  $G$ .

One way to guide BestFirst is by constructing  
and artificial potential field.

## Potential Fields

(show transparency)

Force field  
pushes robot away  
from obstacles

Designed to be  
global minimum

Attractive  
force field  
at  $q_g$

Local  
Minima



Very hard to design potential field w/o  
local minima and global minimum at goal.

"navigation functions" by Koditschek applies  
to special geometries.

It would be helpful to know shape of  $C_{obs}$  to

design potential field, but that won't be available

From Latombe,  
 "Motion Planning"  
 1992.

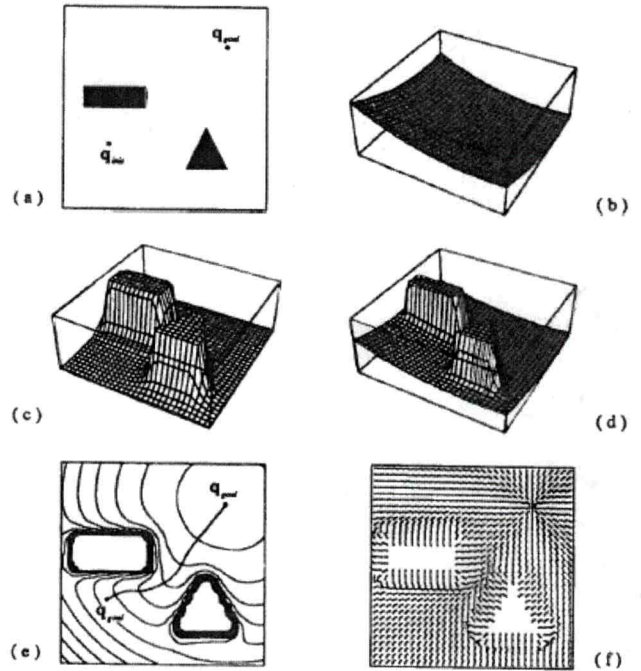
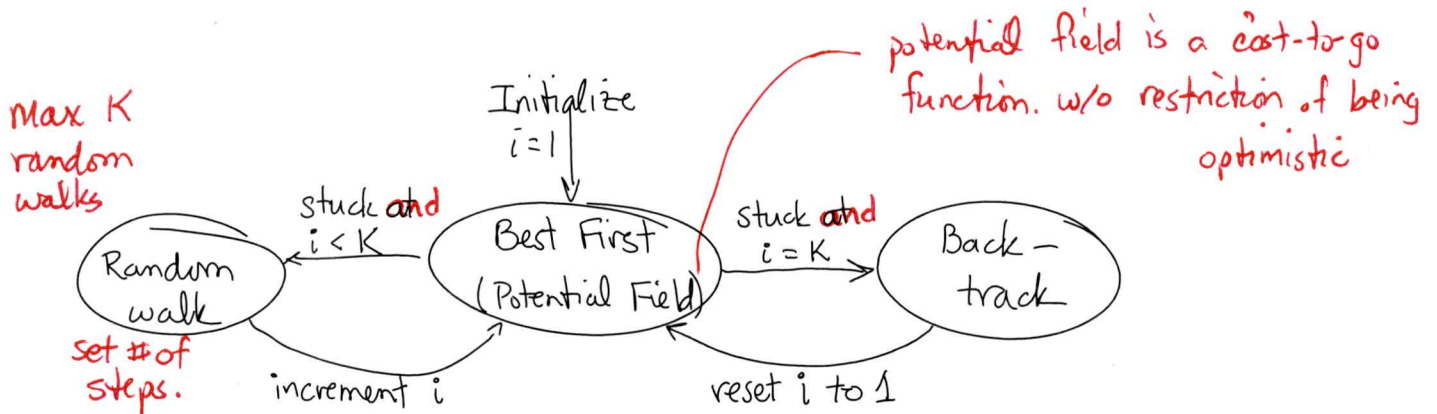


Figure 8. This series of diagrams illustrates the potential field approach. A simple two-dimensional configuration space with two polygonal C-obstacles depicted in Figure a. Figure b shows an attractive potential generated by goal configuration  $q_{goal}$ . Figure c shows a repulsive potential generated by the C-obstacles. Figure d shows the sum of the two potentials. Figure e displays equipotential contours of the total potential and a path obtained by following its negated gradient. Figure f shows orientations of the negated gradient of the

### Randomized Potential Field Method



Follow best first search

random walk  $i \geq 1$

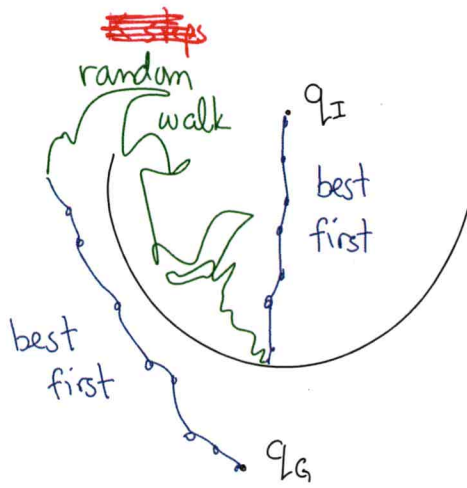
Follow best first search until all neighbors make negative progress.

Switch to random walk for a "while."

Key

You must know how to find neighboring configurations!

This is a vote in favor of grids.



Random walk until  
1) cost is improved  
OR 2) Nw steps performed

If ~~K~~ walks fails  
Start Best First from random  $q \in V$

Questions:

How do we design good pot. fields?

Value of K?

what resolution?

From which node should walk begin?

More detail

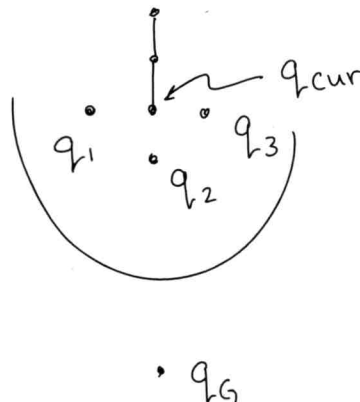
e.g.  $g_r = \frac{a}{\psi_n(q)} = a \frac{1}{\text{dist}}$

$g_a = b \|q - q_G\|$

Let  $g = g_r + g_a$   
 $\uparrow$  repulsive       $\uparrow$  attractive

Let  $g(q)$  denote the potential of  $q$ .

Currently at  $q_{curr}$   
Move next to  $q_i \ni$

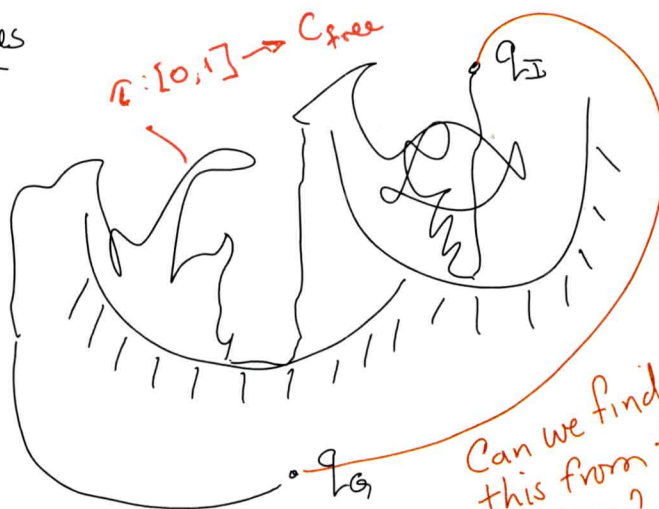


$$g(q_i) < g(q_{cur})$$

If  $g(q_i) \geq g(q_{cur}) \forall i$ , then best first is stuck.

Random walks make very bad paths!

Smoothing issues



One way:

Recursively

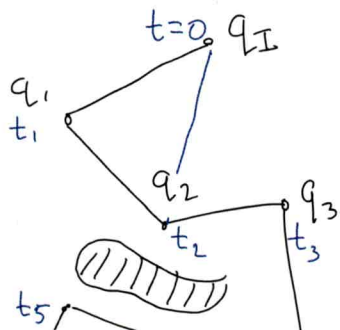
smooth the path,  $\tau$ ,

is to pick pairs of points

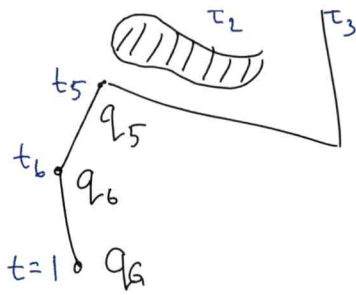
along the path & connect them with a

straight line in  $C$ . Then the new path is:

$$\tau' = \begin{cases} \tau(t) & 0 \leq t \leq t_1 \\ a\tau(t_1) + (1-a)\tau(t_2) & t_1 \leq t \leq t_2 \\ \tau(t) & t_2 \leq t \leq 1 \end{cases}$$



Try  $q_I \rightarrow q_2$ . If successful, cut out  $q_1$  (and cost-to-come lower)



What if we tried  $q_1 \rightarrow q_6$ ?

we could potentially cut out  $q_2, q_3, q_4, q_5$

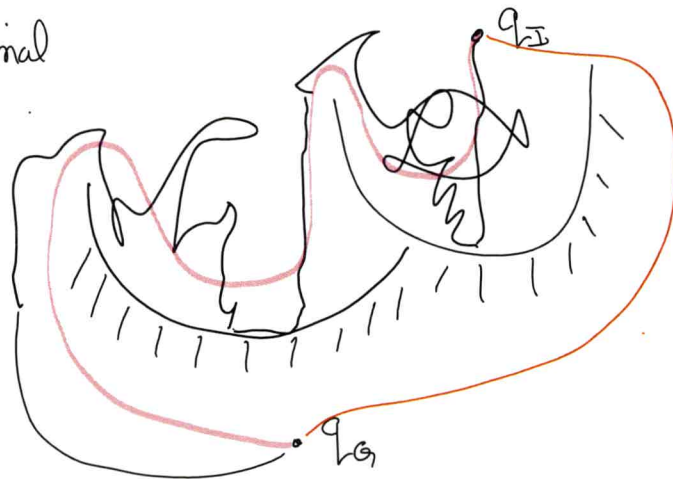
The order of pairs to connect is not cut and dry.

A path metric would be helpful. What's important?

# of turns, energy expended, path length, ...

Suppose we suspect that our smoothed path is from a sub-optimal equivalence class (homotopy class)?

Maybe the red one is better.



You could continue searching until you believe you have a path from every homotopy class, but this would be very hard to determine definitively. Determining this is likely to be as hard as computing Cobs exactly.

In general much parameter tuning is need for good performance. And one tuning may not work well for all problems.

## Other methods

Ariadne's Clew - ~~Use 2 graphs  $G_I$  &  $G_G$~~

Key idea: interleave search & exploration

Algorithm:

- ① VSM - choose vertex  $q_e$  in  $G$  at random
- ② LPM - find new point  $q_{new}$  maximally far from  ~~$q_e$~~  all vertices in  $G$ .  
that can be "easily" connected.
- ③ Try to connect  $q_{new}$  to the other components of  $G$

Drawback - finding  $q_{new}$  is a difficult optimization problem. Optimization method have to be tuned to each motion planning problem.

~~Init w/  $q_I \neq q_G$~~

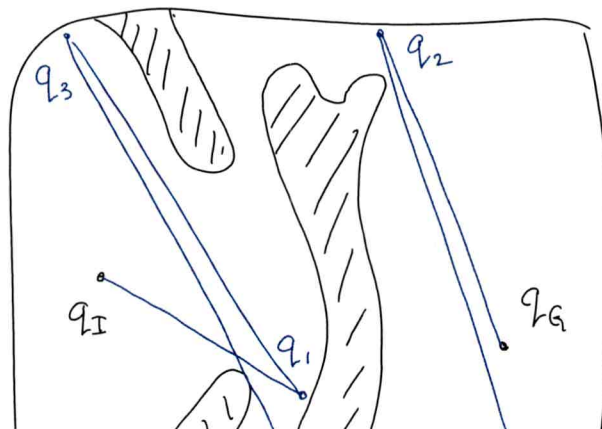
~~Attempt connect ( $q_I, q_G$ )  
fail.~~

~~Choose  $q_e = q_I$~~

~~find  $q_1$~~

~~Connect ( $q_1, q_G$ ) = fail~~

~~...~~



Connect ( $q_1, q_G$ ) = fail

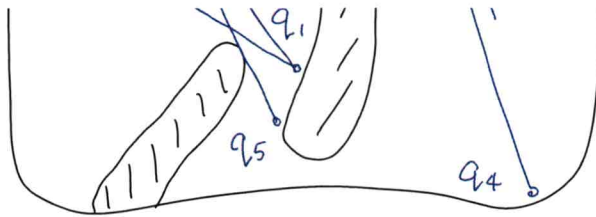
Choose  $q_e = q_G$

Find  $q_2$

Connect ( $q_2, q_I$ ) = fail

Choose  $q_e = q_1$ , Find  $q_3$ , Connect ( $q_3, q_G$ ) = fail.

Alg fails in this example, since only 5 new  $q$ 's will ever be found. That is choosing any  $q_e \in S$  will find  $q_{new}$  already in  $S$ .



### Expansive space planner

Goal: generate samples in unexplored regions of  $C_{free}$

① VSM - select  $q_e$  from  $S$  w/ probability inv. proportional to its # of <sup>connected</sup> neighbors

② LPM - expand  $q_e$  to get  $q_{new}$  w/in a nbhd of  $q_e$

③ Insert  $q_{new}$  into  $S$  w/ probability inv. prop. to # of neighbors of  $q_{new}$

Don't let the graph get high degree.

Three parameters to tune.

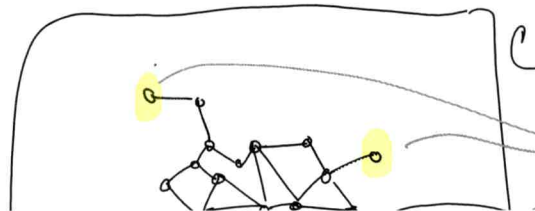
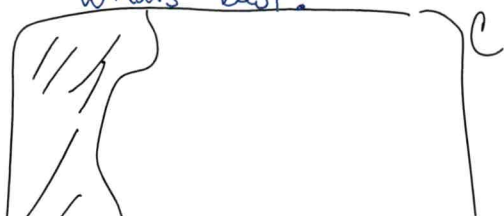
Sizes of nbhds important:

Too small yields uniform prob. over  $S$

Too large " " " "  $S$ .

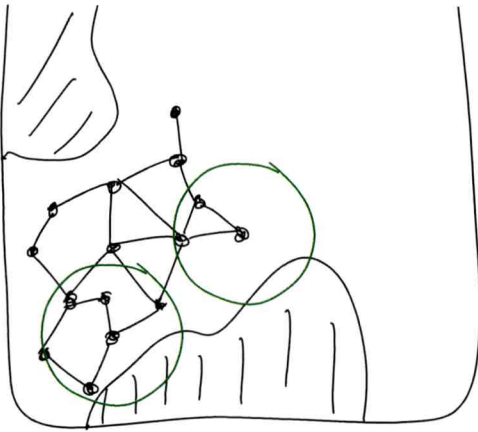
What's best?

Size of nbhd should be tuned to feature resolution



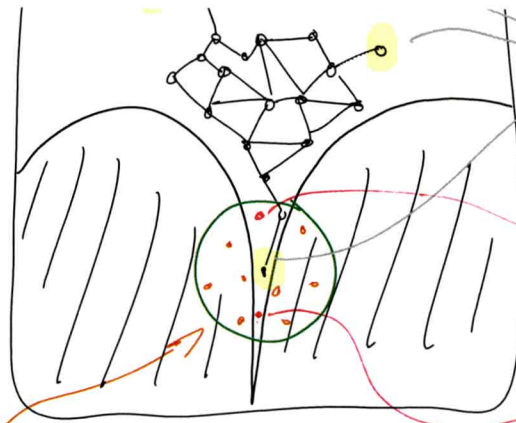
most likely to be chosen





Bias search to explore from nodes on bndry of graph.

Best performance if size of nbhd was tuned to sizes of feature of Cobs, but this is unknown.



most likely to be chosen

don't pick this point. it would have 3 neighbors

pick this point - only one neighbor

Bias search from node with little hope of success.

Could keep failed points, but that would be bad if there was a narrow passage.

### Random Walk Planner

Choose new points from a pdf that adapts to Cobs.

For example new points could be drawn from a multivariate Gaussian with changing mean and covariance matrix



As search progresses in the example, the pdf narrows & aligns in passages and becomes less

narrows & aligns in passages and becomes less biased in wide open areas.

Recall the definition of a multivariate gaussian

$$f(q) = \frac{\exp(-\frac{1}{2}(q-\mu)^T \Sigma^{-1} (q-\mu))}{2\pi^{N/2} |\Sigma|^{1/2}}$$

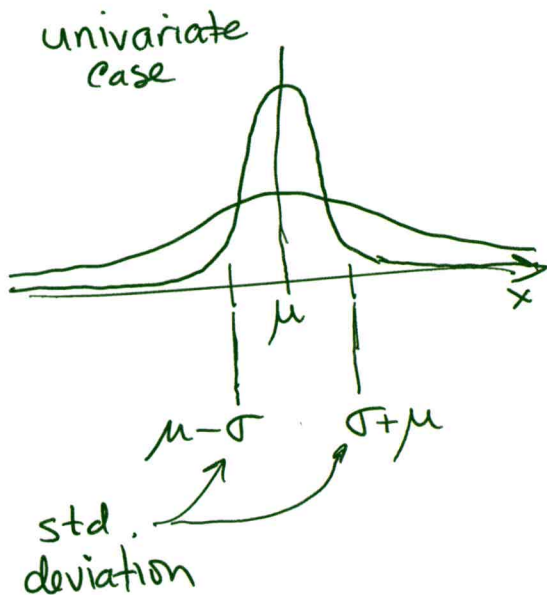
where  $\mu$  is mean of p.d.f.

$\Sigma$  is covariance matrix

$N$  is length of vector  $q$

$|\Sigma|$  is determinant of  $\Sigma$

Use moving average to create changing mean over time.



$$\mu = \sum_{i=1}^k x_i / \frac{1}{k}$$

$$\Sigma_{ij} = \sum_{i=1}^k \sum_{j=1}^k (x_i - \mu)^T (x_j - \mu)$$

summation

not summation, unfortunate choice of variable name

This method also has difficulty w/ winding narrow passages.

So you could use a sum of gaussians, but would be more complicated to compute the distribution

