Probabilistic Robotics Course

Multi-Point Registration

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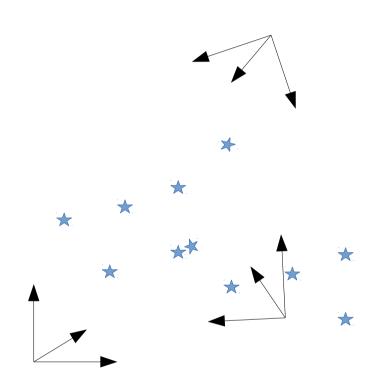
Problem Definition

The environment is populated with identifiable landmarks whose position is unknown

The landmarks are observed by a 3D sensor from multiple positions

We want to determine

- The landmark locations
- The robot positions



State

The state is a collection of transforms and landmark positions

$$\mathbf{X} \qquad : \quad \mathbf{X} = \{\mathbf{X}_r^{[1]}, \dots, \mathbf{X}_r^{[N]}, \mathbf{X}_l^1, \dots, \mathbf{X}_l^{[M]}\}$$
 robot $\mathbf{X}_r^{[n]} \in SE(3) : \qquad \mathbf{X}^{[n]} = \left(\mathbf{R}^{[n]} | \mathbf{t}^{[n]}\right)$ landmark $\mathbf{X}_l^{[m]} \in \Re^3 : \qquad \mathbf{X}_l^{[m]} = \left(x^{[m]} \; y^{[m]} \; z^{[m]}\right)^T$

The increments are represented by a large vector containing the minimal perturbation for each state variable

State (implementation)

In an Octave Implementation, we will represent the states as

- •an array of $\mathbf{4} \times \mathbf{4}$ transformation matrices, for the robot poses
- a matrix of 3 x num_landmarks elements, for the landmarks. Each column is a landmark

State (implementation)

With such an articulated representation of the state it might come in handy to have functions that translate

- a landmark or
- pose index

in the corresponding position in the perturbation vector

```
For a pose:
```

```
v_{idx}=1+(pose_{index}-1)*6;
```

For a landmark:

```
v_idx=1+(num_poses)*6 + (landmark_index-1)*3;
```

State (implementation)

```
function v_idx=poseMatrixIndex(pose_index, num_poses, num_landmarks)
  global pose dim; # 6 in our case
  global landmark dim; # 3 in our case
  if (pose index>num poses)
    v idx=-1:
    return:
  endif:
  v idx=1+(pose index-1)*pose dim;
endfunction:
function v idx=landmarkMatrixIndex(landmark index, num poses, num landmarks)
  global pose dim;
  global landmark dim;
  if (landmark index>num landmarks)
    v idx=-1;
    return:
  endif:
  v idx=1 + (num poses)*pose dim + (landmark index-1) * landmark dim;
endfunction:
```

Boxplus

The boxplus has to be adapted to apply the individual perturbations for each variable block

$$egin{array}{lll} \mathbf{X}' &=& \mathbf{X} oxplus \mathbf{\Delta} \mathbf{x} \ \mathbf{X}_r^{[n]'} &=& \mathbf{\Delta} \mathbf{x}_r^{[n]} oxplus \mathbf{X}_r^{[n]} \ &=& v2t(\mathbf{\Delta} \mathbf{x}_r^{[n]}) \mathbf{X}_r^{[n]} \ \mathbf{X}_l^{[m]'} &=& \mathbf{X}_l^{[m]} + \mathbf{\Delta} \mathbf{x}_l^{[n]} \end{array}$$

Boxplus (implementation)

```
function [XR, XL]=boxPlus(XR, XL, num_poses, num_landmarks, dx)
 global pose dim:
 global landmark dim;
  for(pose_index=1:num_poses)
    pose matrix index=poseMatrixIndex(pose index,
                                       num poses,
                                       num landmarks);
    dxr=dx(pose_matrix_index:pose_matrix_index+pose_dim-1);
    XR(:,:,pose index)=v2t(dxr)*XR(:,:,posea index);
 endfor:
  for(landmark_index=1:num_landmarks)
    landmark matrix index=landmarkMatrixIndex(landmark index,
                                               num poses,
                                               num landmarks);
    dxl=dx(landmark matrix index:landmark matrix index+landmark dim-1,:);
    XL(:,landmark index)+=dxl;
 endfor;
endfunction;
```

Measurements and Predictions

A measurement of the landmark m, performed by robot pose n is as follows

$$\mathbf{z}^{[n,m]} \in \mathbb{R}^3$$
 : $\mathbf{z}^{[m,n]} = (x^{[n,m]} y^{[n,m]} y^{[n,m]})^T$

The prediction and the error of this measurements are very similar to the ICP ones

$$\mathbf{h}^{[n,m]}(\mathbf{X}) = \mathbf{X}_r^{[n]} \mathbf{X}_l^{[m]}$$

$$\mathbf{e}^{[n,m]}(\mathbf{X}) = \mathbf{X}_r^{[n]} \mathbf{X}_l^{[m]} - \mathbf{z}^{[n,m]}$$

$$\mathbf{e}^{[n,m]}(\mathbf{X} \boxplus \mathbf{\Delta} \mathbf{x}) = v2t(\mathbf{\Delta} \mathbf{x}_r^{[n]}) \mathbf{X}_r^{[n]}(\mathbf{X}_l^{[m]} + \mathbf{\Delta} \mathbf{x}_l^{[m]}) - \mathbf{z}^{[n,m]}$$

Jacobians

The prediction depends only on the robot pose m and the landmark n, so it will be mostly 0

Non 0 only for

- pose block n
- landmark block m

Jacobians (implementation)

```
function [e,Jr,Jl]=errorAndJacobian(Xr,Xl,z)
   R=Xr(1:3,1:3);
   t=Xr(1:3,4);
   z hat=R*Xl+t; #prediction
   e=z hat-z;
   Jr=zeros(3,6);
   Jr(1:3,1:3)=eye(3);
   Jr(1:3,4:6) = -skew(z hat);
   Jl=R:
endfunction;
```

H Matrix and B vector

H and b for a measurement have only few non zero blocks

$$\mathbf{H}^{[n,m]} = \mathbf{J}^{[n,m]T} \mathbf{\Omega}^{[n,m]} \mathbf{J}^{[n,m]} \mathbf{J}^{[n,m]T} \mathbf{\Omega}^{[n,m]} \mathbf{J}^{[n,m]}_{r}$$
 $= \begin{pmatrix} \mathbf{J}^{[n,m]T}_{r} \mathbf{\Omega}^{[n,m]} \mathbf{J}^{[n,m]}_{r} \mathbf{J}^{[n,m]T}_{r} \mathbf{\Omega}^{[n,m]} \mathbf{J}^{[n,m]}_{l} \\ \mathbf{J}^{[n,m]T}_{l} \mathbf{\Omega}^{[n,m]} \mathbf{J}^{[n,m]}_{l} \mathbf{J}^{[n,m]T}_{l} \mathbf{\Omega}^{[n,m]} \mathbf{J}^{[n,m]}_{l} \end{pmatrix}$
 $= \begin{pmatrix} \mathbf{J}^{[n,m]T}_{r} \mathbf{\Omega}^{[n,m]} \mathbf{e}^{[n,m]} \\ \mathbf{J}^{[n,m]T}_{r} \mathbf{\Omega}^{[n,m]} \mathbf{e}^{[n,m]} \end{pmatrix}$

H and B Implementation

```
H=zeros(system size, system size);
b=zeros(system size,1);
chi stats(iteration)=0;
for (measurement_num=1:size(Z,2))
  pose index=associations(1,measurement num);
  landmark index=associations(2,measurement num);
  z=Z(:,measurement_num);
  Xr=XR(:,:,pose index);
  Xl=XL(:,landmark index);
  [e,Jr,Jl] = errorAndJacobian(Xr, Xl, z);
  Hrr=Jr'*Jr;
 Hrl=Jr'*Jl:
  Hll=Jl'*Jl;
  br=Jr'*e;
  bl=Jl'*e:
  pose matrix index=poseMatrixIndex(pose index, num poses, num landmarks);
  landmark matrix index=landmarkMatrixIndex(landmark index, num poses, num landmarks);
  H(pose matrix index:pose matrix index+pose dim-1,
    pose matrix index:pose matrix index+pose dim-1)+=Hrr;
  H(pose_matrix_index:pose_matrix_index+pose_dim-1,
    landmark matrix index:landmark matrix index+landmark dim-1)+=Hrl;
  H(landmark matrix index:landmark matrix index+landmark dim-1,
    landmark_matrix_index:landmark_matrix_index+landmark_dim-1)+=Hll;
  H(landmark matrix index:landmark matrix index+landmark dim-1,
    pose matrix index:pose matrix index+pose dim-1)+=Hrl';
  b(pose matrix index:pose matrix index+pose dim-1)+=br;
  b(landmark matrix index:landmark matrix index+landmark dim-1)+=bl;
endfor
```

Degrees of Freedom

A state configuration X, has the same error of all states obtained by applying a rigid transform to all variables.

The system will be under-determined

6 degrees of freedom (dof of a rigid transform)

To solve the system we need to either

- eliminate the redundant variables
- adding a prior

Eliminating Variables

We can impose that one robot pose stays fixed To this extent we can

- reduce the linear system by suppressing the rows and the columns of the variables that stay fixed.
- solve the reduced system
- propagate the computed perturbation to all variables that have not been eliminated

Eliminating Variables (impl)

```
dx=zeros(system_size,1);
```

```
% we solve the linear system, blocking the first pose
% this corresponds to "remove" from H and b the locks
% of the 1st pose, while solving the system
```

```
dx(pose_dim+1:end)=
    -(H(pose_dim+1:end,pose_dim+1:end)\
    b(pose_dim+1:end,1));
[XR, XL]=boxPlus(XR,XL,num_poses, num_landmarks, dx);
```

Adding a Prior

Alternatively we can add a "flexible" constraint of the type $\Delta \mathbf{x}_r^{[1]} = \mathbf{0}$ to the least squares problem.

This is done by adding an error term

$$egin{array}{lll} \mathbf{e}_{\mathrm{prior}} &=& \mathbf{\Delta}\mathbf{x}_r^{[1]} \ \mathbf{J}_{\mathrm{prior}} &=& \mathbf{I} \ \mathbf{H}_{\mathrm{prior}} &=& \mathbf{\Omega}_{\mathrm{prior}} \ \mathbf{b}_{\mathrm{prior}} &=& \mathbf{0} \end{array}$$

Can be implemented by adding to the final H a 6x6 (large) information matrix in correspondence of the first variable

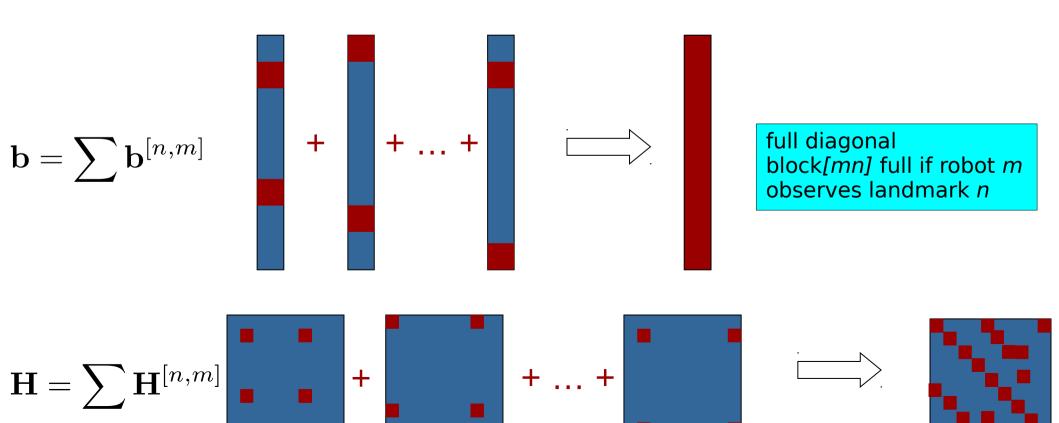
The larger the Omega, the more stiff the constraint will be

Adding a Prior (impl)

Structure

Structure of **H**: location of the non zero elements

The structure of H depends only on the structure of the measurements



Conclusions

We approached a complex multi-robot multilandmark problem

- The measurement independence leads to a sparse structure of H
- The structure of H does not change during the iterations
- It can be efficiently solved by using sparse methods (we will see these later in this course)