Advanced Algorithm Implementation: Informed RRT* and D*

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1. INFORMED RRT*

a. Introduction and Methodology

The basic idea behind the INFORMED RRT* is additional constraints in the sampling process. Initially, the informed RRT* starts with the same process as the RRT* algorithm. Like RRT* the new node is connected to the nearest node. Then the neighbors of the new node are considered and the cost-to-come from start to the new node is checked through all the neighbors. The neighbor that provides the least cost-to-come to the new node is chosen as the parent for the new node. Then the connections are rewired such that the cost-to-come for the neighbors are less than the cost-to-come through the new node, if not the new node is assigned as the parent for the neighbor node.

After the path is found, then the sampling process is changed. Instead of sampling randomly from the entire space, the sampling is restricted to an ellipse (or ellipsoid in 3D) with the focal points being the start and the end nodes. Based on this restricted sampling, a new point is obtained. This new point is then used to rewire and replan the entire path.

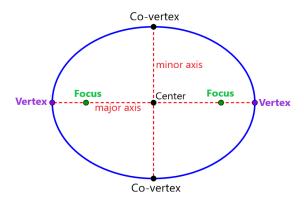
The following Pseudo code for Informed RRT* gives a brief flow of the process:

```
Algorithm 1: Informed RRT*(\mathbf{x}_{\text{start}}, \mathbf{x}_{\text{goal}})

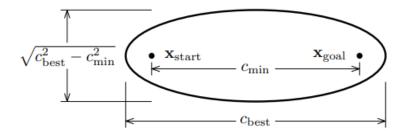
    V ← {x<sub>start</sub>};

  2 E ← ∅;
  X_{\mathrm{soln}} \leftarrow \emptyset;
  4 \mathcal{T} = (V, E);
  5 for iteration = 1 \dots N do
                  \begin{aligned} c_{\text{best}} &\leftarrow \min_{\mathbf{x}_{\text{soln}} \in X_{\text{soln}}} \left\{ \texttt{Cost} \left( \mathbf{x}_{\text{soln}} \right) \right\}; \\ \mathbf{x}_{\text{rand}} &\leftarrow \texttt{Sample} \left( \mathbf{x}_{\text{start}}, \mathbf{x}_{\text{goal}}, c_{\text{best}} \right); \end{aligned}
                   \mathbf{x}_{\text{nearest}} \leftarrow \text{Nearest} (\mathcal{T}, \mathbf{x}_{\text{rand}});
  8
                    \mathbf{x}_{\text{new}} \leftarrow \text{Steer}\left(\mathbf{x}_{\text{nearest}}, \mathbf{x}_{\text{rand}}\right);
                   if CollisionFree (\mathbf{x}_{\mathrm{nearest}}, \mathbf{x}_{\mathrm{new}}) then
10
11
                               V \leftarrow \cup \{\mathbf{x}_{\text{new}}\};
                              X_{\text{near}} \leftarrow \text{Near}\left(\mathcal{T}, \mathbf{x}_{\text{new}}, r_{\text{RRT}^*}\right);
12
                              \mathbf{x}_{\min} \leftarrow \mathbf{x}_{\mathrm{nearest}};
13
                              c_{\min} \leftarrow \text{Cost}(\mathbf{x}_{\min}) + c \cdot \text{Line}(\mathbf{x}_{\text{nearest}}, \mathbf{x}_{\text{new}});
14
                              for \forall \mathbf{x}_{\text{near}} \in X_{\text{near}} do
15
                                          c_{\text{new}} \leftarrow \text{Cost}(\mathbf{x}_{\text{near}}) + c \cdot \text{Line}(\mathbf{x}_{\text{near}}, \mathbf{x}_{\text{new}});
16
                                          if c_{\text{new}} < c_{\text{min}} then
17
                                                     if CollisionFree (\mathbf{x}_{\mathrm{near}}, \mathbf{x}_{\mathrm{new}}) then
18
                                                                 \mathbf{x}_{\min} \leftarrow \mathbf{x}_{\text{near}};
19
                                                                c_{\min} \leftarrow c_{\text{new}};
20
                               E \leftarrow E \cup \{(\mathbf{x}_{\min}, \mathbf{x}_{\text{new}})\};
21
                              for \forall \mathbf{x}_{\text{near}} \in X_{\text{near}} do
                                          c_{\text{near}} \leftarrow \text{Cost}(\mathbf{x}_{\text{near}});
23
                                          c_{\text{new}} \leftarrow \text{Cost}(\mathbf{x}_{\text{new}}) + c \cdot \text{Line}(\mathbf{x}_{\text{new}}, \mathbf{x}_{\text{near}});
24
                                          if c_{\text{new}} < c_{\text{near}} then
25
                                                     if CollisionFree (\mathbf{x}_{\mathrm{new}},\mathbf{x}_{\mathrm{near}}) then
26
27
                                                                \mathbf{x}_{\mathrm{parent}} \leftarrow \mathtt{Parent}\left(\mathbf{x}_{\mathrm{near}}\right);
                                                                 E \leftarrow E \setminus \{(\mathbf{x}_{\text{parent}}, \mathbf{x}_{\text{near}})\};
28
                                                                E \leftarrow E \cup \{(\mathbf{x}_{\text{new}}, \mathbf{x}_{\text{near}})\};
29
                              if InGoalRegion(x_{new}) then
30
                                        X_{\text{soln}} \leftarrow X_{\text{soln}} \cup \{\mathbf{x}_{\text{new}}\};
31
32 return T;
```

In the above pseudo code, the best path length calculated in each iteration is considered. This path length (C_{best}) is assigned as the length of the major axis. The distance between the two focal points is taken as ' C_{min} '.



From the relation given below, using C_{best} and C_{min} , the minor axis is calculated.



Then a point is sampled normally from a unit circle (or sphere in 3D) and then it is transformed into the frame of the ellipse using rotation matrix as shown below in lines 8 and 9:

```
Algorithm 2: Sample (\mathbf{x}_{\text{start}}, \mathbf{x}_{\text{goal}}, c_{\text{max}})

1 if c_{\text{max}} < \infty then

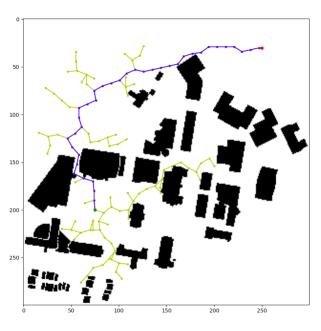
2 \begin{vmatrix} c_{\text{min}} \leftarrow ||\mathbf{x}_{\text{goal}} - \mathbf{x}_{\text{start}}||_2; \\ \mathbf{x}_{\text{centre}} \leftarrow (\mathbf{x}_{\text{start}} + \mathbf{x}_{\text{goal}})/2; \\ \mathbf{C} \leftarrow \text{RotationToWorldFrame} (\mathbf{x}_{\text{start}}, \mathbf{x}_{\text{goal}}); \\ \mathbf{r}_1 \leftarrow c_{\text{max}}/2; \\ \mathbf{6} & \{r_i\}_{i=2,\dots,n} \leftarrow \left(\sqrt{c_{\text{max}}^2 - c_{\text{min}}^2}\right)/2; \\ \mathbf{1} \leftarrow \text{diag} \{r_1, r_2, \dots, r_n\}; \\ \mathbf{1} \mathbf{x}_{\text{ball}} \leftarrow \text{SampleUnitNBall}; \\ \mathbf{x}_{\text{rand}} \leftarrow (\mathbf{CL}\mathbf{x}_{\text{ball}} + \mathbf{x}_{\text{centre}}) \cap X; \\ \mathbf{10} & \mathbf{else} \\ \mathbf{11} & \mathbf{x}_{\text{rand}} \sim \mathcal{U}(X); \\ \mathbf{12} & \mathbf{return} \ \mathbf{x}_{\text{rand}}; \\ \end{aligned}
```

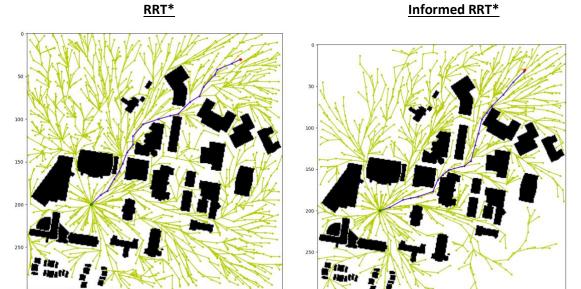
Using this new point, the path is replanned, and the new shortest path is obtained.

b. Results and Observations

The results shown below are for RRT, RRT* and informed-RRT*:







From the above plots it can be observed that the informed RRT* star has spread in a smaller region compared to the RRT* as well as both RRT* and informed RRT* have shorter path than the RRT algorithm.

Key difference between RRT* and informed RRT*:

RRT*	Informed-RRT*
New point is randomly sampled from the entire	New point is sampled from the restricted workspace
workspace.	(Ellipse/Ellipsoid region).
More number of nodes are explored because of large	Less number of nodes are explored to find a equivalent
workspace, to find an optimum path	optimum path because of confined sampling.
Path information such as path length is not utilized to	Path information such as path length is utilized to
help with the sampling process.	restrict the sampling process, as it is considered as
	major axis of ellipse/ellipsoid.

Advantages of using informed-RRT* over RRT*:

```
ajith@ajith-Inspiron-5558:~/Documents/Advanced Search Algorithms/informed_RRT$ python main.py
It took 118 nodes to find the current paths
The path length is 363.00
It took 1425 nodes to find the current path
The path length is 258.60
It took 1062 nodes to find the current path
The path length is 262.30
```

From the above output, it can be observed that informed-RRT* is able to obtain a significantly close optimum path like the RRT* but it utilizes a smaller number of nodes than the latter. Thus, informed-RRT* is both computationally as well as memory wise more efficient than the standard RRT* algorithm giving almost the same optimum path length.

From the above images of RRT* and informed-RRT*, it can be observed that the sampling region is restricted (within an ellipse) in the latter algorithm, thus reducing unnecessary sampling of nodes outside the required region, for a better path.

2. <u>D*</u>

a. Introduction and Methodology

In this section, the implementation of D* is discussed. Like Dijkstra or A*, D* algorithm also maintains an open list. Each node has a different state tag as follows:

- a. NEW
- b. OPEN
- c. CLOSED
- d. RAISE
- e. LOWER

Based on each tag, the nodes are processed using the Process_State function. In Process_State function, we initialize the goal's heuristic as zero and then insert into the open list. The function is called until robot's state is removed from the open list. During each process state, the Prepare_Repair function is called. This function checks if the neighbours of the current state is an obstacle or not. If the one of the neighbours is an obstacle.

If one of the neighbours is observed to be an obstacle in the Prepare_Repair function, then the Modify_cost function is called. This function modified the neighbours cost according to the new obstacles. After the neighbours cost are modified, then the Repair-Replan function is called which takes care of repairing the cost of the neighbours by replanning the path.

The following pseudo codes represent the operations that are described above:

D* Algorithm

```
1: for each X \in L do
2: t(X) = NEW
3: end for
4: h(G) = 0
5: INSERT(O, G, h(G))
6: X_c = S
7: P = INIT - PLAN(O, L, X_c, G)
8: if P = NULL then
9: Return (NULL)
10: end if
11: while X_c \neq G do
   PREPARE - REPAIR(O, L, X_c)
   P = REPAIR - REPLAN(O, L, X_c, G)
14: if P = NULL then
     Return (NULL)
15:
16: end if
   X_c = the second element of P {Move to the next state in P }.
18: end while
19: Return (X_c)
```

Process_State()

```
X = MIN-STATE()
                                                                           for each neighbor Y of X:
if X = NULL then return -1
                                                                                      if t(Y) = NEW or
k_{old} = GET-KMIN(); DELETE(X);
                                                                                      (b(Y) = X \text{ and } h(Y) \neq h(X) + c(X,Y)) then
if k_{old} < h(X) then
                                                                                            b(Y) = X; INSERT(Y, h(X)+c(X,Y))
   for each neighbor Y of X:
                                                                                      else
               if t(Y) \neq new and h(Y) \le k_{old} and h(X) > h(Y) +
                                                                                           if b(Y) \neq X and h(Y) > h(X)+c(X,Y) then
c(Y,X) then
                                                                                             INSERT(X, h(X))
                 b(X) = Y; h(X) = h(Y) + c(Y,X);
if k_{old} = h(X) then
                                                                                             if b(Y) \neq X and h(X) > h(Y) + c(X,Y) and
   for each neighbor Y of X:
                                                                                             t(Y) = CLOSED and h(Y) > k_{old} then
               if t(Y) = NEW or
                                                                                             INSERT(Y, h(Y))
               (b(Y) = X \text{ and } h(Y) \neq h(X) + c(X,Y)) \text{ or }
                                                                       Return GET-KMIN ()
               (b(Y) \neq X \text{ and } h(Y) > h(X)+c(X,Y)) then
```

b(Y) = X; INSERT(Y, h(X)+c(X,Y))

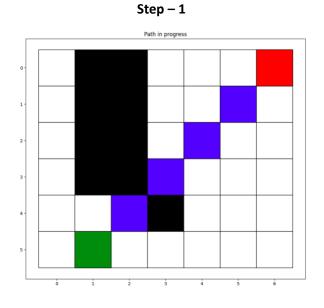
```
MODIFY - COST(O, X, Y, cval)
                                                         REPAIR - REPLAN(O, L, X_c, G)
1: c(X,Y) = cval
                                                         1: repeat
 2: if t(X) = CLOSED then
                                                         2: k_{min} = PROCESS - STATE(O, L)
      INSERT(O, X, h(X))
                                                         3: until (k_{min} \ge h(X_c)) or (k_{min} = -1)
                                                         4: P = GET - BACKPOINTER - LIST(L, X_c, G)
 4: end if
                                                         5: Return (P)
 5: Return GET - KMIN(O)
                       PREPARE - REPAIR(O, L, X_c)
                       1: for each state X \in L within sensor range of X_c and X_c do
                       2: for each neighbor Y of X do
                             if r(Y,X) \neq c(Y,X) then
                               MODIFY - COST(O, Y, X, r(Y, X))
                       4:
                             end if
                       6: end for
                       7: for each neighbor Y of X do
                             if r(X,Y) \neq c(X,Y) then
                       9:
                               MODIFY - COST(O, X, Y, r(X, Y))
```

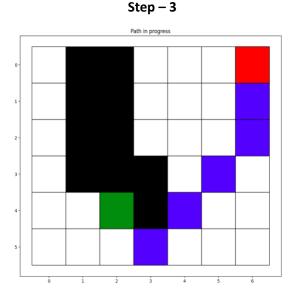
b. Results and Observations

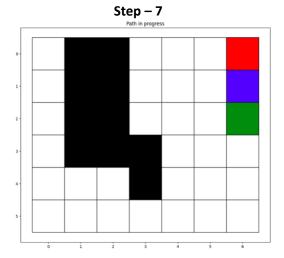
10:

11: end for 12: end for

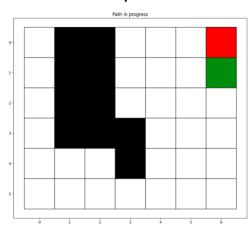
SCENARIO -1



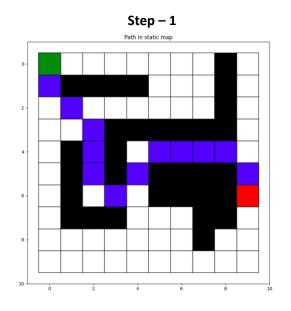


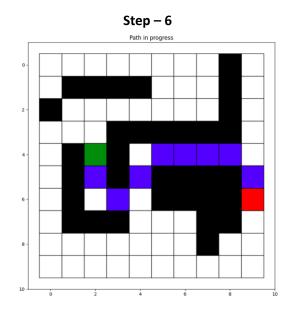


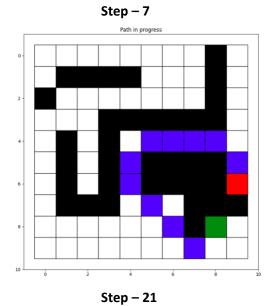
Step – 8

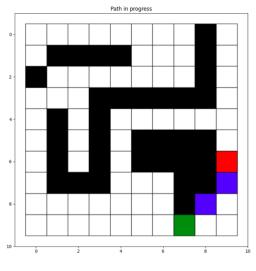


SCENARIO – 2

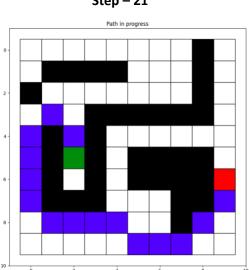


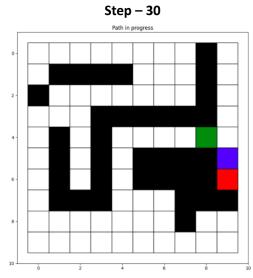




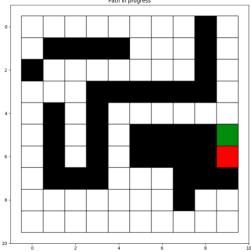


Step - 20

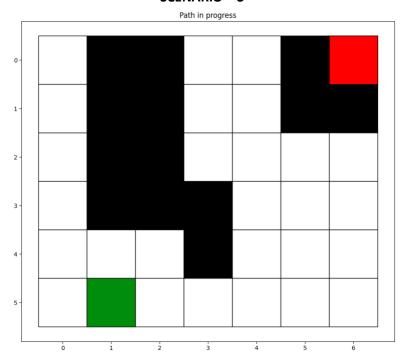








SCENARIO - 3



Outputs

For three different scenarios/maps, the following results were obtained:

```
ajith@ajith-Inspiron-5558:~/Documents/Advanced Search Algorithms/D_star$ python3 main.py ajith@ajith-Inspiron-5558:~/Documents/Advanced Search Algorithms/D_star$ python3 main.py ajith@ajith-Inspiron-5558:~/Documents/Advanced Search Algorithms/D_star$ python3 main.py No path is found
```

D* vs A* or Dijkstra

- 1. A* and Dijkstra are used for static environment whereas D* is used for dynamic environment.
- 2. In case we use A* and Dijkstra in the dynamic environment, then the entire path must be replanned from the current position to the goal position. During this, the entire environment is again reevaluated. Whereas in the case of the D* algorithm, the changes are made locally when the obstacle is faced and based on the change, a different path is followed which is optimum from the changed direction. Thus, it can be observed that A* and Dijkstra are computationally more expensive than D*. Therefore, D* is comparatively faster than A* and Dijkstra algorithm.