Model Predictive Control

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Contents

1	Mo	tion Control	2
2	Noı	a linear optimization under constraints	on under constraints 2
	2.1	Definition	2
	2.2	Example	3
	2.3	Solving with ipopt	3
3	Bas	ic Math utilities	6
	3.1	Translate into vehicle's coordinate system	6
	3.2	Fit a polynomial to a set of waypoints	6
	3.3	Compute the tangential angle of a polynomial curve	7
4	Vehicle Models		8
	4.1	Dynamic vs Kinematic Models	8
	4.2	Kinematic Model	8
		4.2.1 State	8
		4.2.2 Deriving the kinematic model	8
		4.2.3 Errors	11
		4.2.4 Actuator Constraints	13
		4.2.5 Constraints summary	13
	4.3	Dynamic Models	14
5	Model Predictive Control		14
	5.1	Minimization of a cost function	15
	5.2	Timsestep length and Elapsed duration	16
	5.3	Latency handling	16
	5.4	MPC Solver algorithm	17
	5.5	MPC Solver code	19

1 Motion Control

Motion Control deals with the last stage of an autonomous driving pipeline: the control module. The input to the control module will be provided by the output of the path planning module via a set of waypoints to follow as close as possible. The control module will have to provide the actuators commands (in our case steering angle and throttling; acceleration or deceleration) so that the automated driving comply with a set of rules:

- follow the planned waypoints as close as possible
- drives smoothly
- try to adjust the speed: as fast as a configurable reference when possible and driving more slowly during curves



Figure 1: Autonomous Driving pipeline

2 Non linear optimization under constraints

2.1 Definition

In its most generic form we are dealing with the following problem:

minimize
$$f_0(x)$$

subject to $lower_i \leq f_i(x) \leq upper_i, i = 1, ..., m.$

Note that by setting $lower_i = upper_i$ we can define constraints as equalities as well.

2.2 Example

```
minimize x_1 * x_4 * (x_1 + x_2 + x_3) + x_3

subject to x_1 * x_2 * x_3 * x_4 \ge 25

x_1^2 + x_2^2 + x_3^2 + x_4^2 = 40

1 \le x_1, x_2, x_3, x_4 \le 5
```

2.3 Solving with ipopt

ipopt and cppad are used to solve non-linear minimization problems. ipopt requires the computation of first order (Jacobians) and 2nd order derivatives (Hessians). These derivatives will be computed automatically thanks to cppad: providing automatic differentiation services.

The previous example is solved with ipopt and CppAD here: https://www.coin-or.org/CppAD/Doc/ipopt_solve_get_started.cpp.htm

Listing 1: Simple example with ipopt

```
1
 2
 3
   # include <cppad/ipopt/solve.hpp>
 4
 5
   namespace {
 6
         using CppAD::AD;
7
8
         class FG_eval {
9
         public:
              typedef CPPAD_TESTVECTOR( AD<double> ) ADvector;
10
11
              void operator () (ADvector& fg, const ADvector& x)
12
                     assert(fg.size() = 3);
                    assert(x.size() = 4);
13
14
15
                    // Fortran style indexing
                   AD < double > x1 = x[0];
16
17
                   AD < double > x2 = x[1];
                   AD < double > x3 = x[2];
18
                   AD < double > x4 = x[3];
19
20
                    // f(x)
                    fg[0] = x1 * x4 * (x1 + x2 + x3) + x3;
21
```

```
22
                   // g_{-1} (x)
23
                   fg[1] = x1 * x2 * x3 * x4;
24
                   // g_{-2} (x)
25
                   fg[2] = x1 * x1 + x2 * x2 + x3 * x3 + x4 * x4;
26
27
                   return;
28
              }
29
         };
30 }
31
32 bool get_started(void)
33
         bool ok = true;
34
         size_t i;
         typedef CPPAD_TESTVECTOR( double ) Dvector;
35
36
37
         // number of independent variables (domain dimension for f and g)
38
         size_t nx = 4;
39
         // number of constraints (range dimension for g)
40
         size_t ng = 2;
         // initial value of the independent variables
41
42
         Dvector xi(nx);
         xi[0] = 1.0;
43
44
         xi[1] = 5.0;
         xi[2] = 5.0;
45
46
         xi[3] = 1.0;
47
         // lower and upper limits for x
48
         Dvector xl(nx), xu(nx);
         for (i = 0; i < nx; i++)
49
50
               xl[i] = 1.0;
              xu[i] = 5.0;
51
52
53
         // lower and upper limits for g
54
         Dvector gl(ng), gu(ng);
55
         gl[0] = 25.0;
                            gu[0] = 1.0e19;
56
         gl[1] = 40.0;
                            gu[1] = 40.0;
57
58
         // object that computes objective and constraints
59
         FG_eval fg_eval;
60
         // options
61
```

```
62
        std::string options;
63
        // turn off any printing
        options += "Integer print_level
64
                                           0 \setminus n";
        options += "String sb
65
                                           yes \n";
66
        // maximum number of iterations
67
        options += "Integer max_iter
                                           10\n";
        // approximate accuracy in first order necessary conditions;
68
        // see Mathematical Programming, Volume 106, Number 1,
69
70
        // Pages 25-57, Equation (6)
71
        options += "Numeric tol
                                           1e-6 n;
72
        // derivative testing
73
        options += "String
                             derivative_test
                                                          second-order\n";
74
        // maximum amount of random pertubation; e.g.,
75
        // when evaluation finite diff
76
        options += "Numeric point_perturbation_radius 0.\n";
77
78
        // place to return solution
79
        CppAD::ipopt::solve_result < Dvector > solution;
80
81
        // solve the problem
82
        CppAD::ipopt::solve<Dvector, FG_eval>(
83
              options, xi, xl, xu, gl, gu, fg_eval, solution
84
        );
85
        // Check some of the solution values
86
87
88
        ok &= solution.status == CppAD::ipopt::solve_result < Dvector >::success
89
90
        double check_x[] = { 1.000000, 4.743000, 3.82115, 1.379408 };
91
        double check_zl[] = \{1.087871, 0.,
   0.
        double check_zu[] = \{ 0.,
92
                                          0.,
                                                     0.,
   0.
             };
                           = 1e-6; // relative tolerance
93
        double rel_tol
        double abs_tol = 1e-6; // absolute tolerance
94
95
        for (i = 0; i < nx; i++)
              ok &= CppAD:: NearEqual(
96
                   check_x[i], solution.x[i], rel_tol, abs_tol
97
98
99
              ok &= CppAD:: NearEqual(
```

```
check_zl[i], solution.zl[i], rel_tol, abs_tol
100
101
               );
               ok &= CppAD::NearEqual(
102
                     check_zu[i], solution.zu[i], rel_tol, abs_tol
103
104
               );
          }
105
106
107
          return ok;
108
```

3 Basic Math utilities

3.1 Translate into vehicle's coordinate system

To make the computations more simple we will work in vehicule's coordinate system at time t: so we have to operate $Rot(-\psi_t) \circ Translation(\begin{bmatrix} x_t \\ y_t \end{bmatrix})$ on current world's coordinates to translate them into vehicule's coordinates

As a reminder, for a rotation matrix of angle θ , we have $Rot(\theta) = \begin{bmatrix} \cos(\theta) & -\sin(\theta) \\ \sin(\theta) & \cos(\theta) \end{bmatrix}$. In our case we are dealing with $Rot(-\psi_t) = \begin{bmatrix} \cos(-\psi_t) & -\sin(-\psi_t) \\ \sin(-\psi_t) & \cos(-\psi_t) \end{bmatrix}$

So we have:

$$\begin{bmatrix} x' \\ y' \end{bmatrix} = \begin{bmatrix} \cos(-\psi_t) & -\sin(-\psi_t) \\ \sin(-\psi_t) & \cos(-\psi_t) \end{bmatrix} \begin{bmatrix} x - x_t \\ y - y_t \end{bmatrix}$$

3.2 Fit a polynomial to a set of waypoints

Let's assume the reference trajectory provided by the Path Planning module is a set of waypoints approximated or fitted by a (3rd order typically) polynomial $f(x_t) = a_0 + a_1x_t + a_2x_t^2 + a_3x_t^3$ and our real position on the y-axis is y_t

How can we frame the problem?

Our unknown vector corresponds to the unknown polynomial coefficients

$$X = [a_0, a_1, a_2, a_3]^T$$

Now let's assume we have 5 samples

$$A = \begin{bmatrix} 1 & x_t^{(1)} & x_t^{2(1)} & x_t^{3(1)} \\ 1 & x_t^{(2)} & x_t^{2(2)} & x_t^{3(2)} \\ 1 & x_t^{(3)} & x_t^{2(3)} & x_t^{3(3)} \\ 1 & x_t^{(4)} & x_t^{2(4)} & x_t^{3(4)} \\ 1 & x_t^{(5)} & x_t^{2(5)} & x_t^{3(5)} \end{bmatrix}$$

$$B = \begin{bmatrix} f(x_t^{(1)}) \\ f(x_t^{(2)}) \\ f(x_t^{(3)}) \\ f(x_t^{(4)}) \\ f(x_t^{(5)}) \end{bmatrix}$$

We have:

$$AX = B$$

A is not squarred, solve this problem with e.g. the pseudo-inverse.

Pseudo inverse quick recap:

AX = B with A a non squarred matrix

$$A^T A X = A^T B$$
 with $A^T A$ squarred

So we may inverse it now if A^TA is not degenerated (use samples that are different so that A^TA is not degenerated)

$$X = (A^T A)^{-1} A^T B$$

The pseudo-inverse of A a non squarred matrix is

$$A^{\dagger} = (A^T A)^{-1} A^T$$

3.3 Compute the tangential angle of a polynomial curve

https://en.wikipedia.org/wiki/Tangential_angle

If the curve is given by y = f(x), then we may take (x, f(x)) as the parametrization, and we may assume the tangential angle ψ is between $\pi/2$ and $\pi/2$. This produces the explicit expression:

$$\psi = arctan(f'(x))$$

4 Vehicle Models

4.1 Dynamic vs Kinematic Models

4.2 Kinematic Model

4.2.1 State

The state variables are the following:

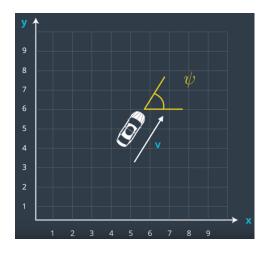


Figure 2: Model state

• x: x position

• y: y position

• ψ : angle between speed vector and x-axis

 \bullet v: speed vector

4.2.2 Deriving the kinematic model

Our state vector is

$$S_t = [x_t, y_t, \psi_t, v_t]$$

We derive an approximation model, kinematic, relating S_{t+1} and S_t . The smaller the dt the more accurate the model.

The model used id the kinematic bicycle model:

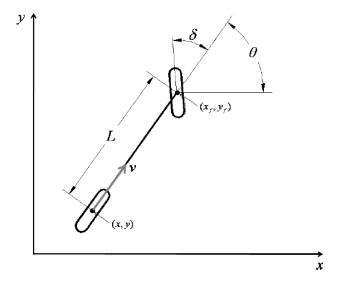


Figure 3: Kinematic bicycle model

Linear movement approximation: assuming during dt that v_t and ψ_t are constant:

$$x_{t+1} = x_t + v_t * \cos(\psi_t) * dt$$

$$y_{t+1} = y_t + v_t * \sin(\psi_t) * dt$$

Rotational movement approximation: assuming during dt that v_t and steering angle δ_t are constant we are creating a change from a starting ψ_t to end up with ψ_{t+1} :

Let first compute the radius of curvature R based on the steering angle δ and the distance L between rear and front wheels:

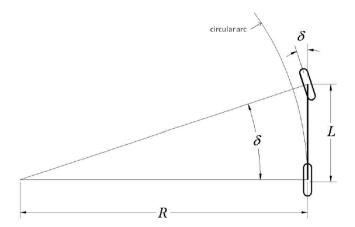


Figure 4: Radius of curvature R

We are moving on a circle from position M_t to position M_{t+1} at a constant speed v_t

$$M_{t+1}M_t = R * (\psi_{t+1} - \psi_t) \approx v_t * dt$$

We have

$$\tan(\delta) \approx L/R$$

Check that the 2 δ drawn are indeed the same. In the triangle we have $\pi/2 + \delta_{circle} + \alpha = \pi$ and along the line from the center of the circle of radius R going through the center of the front wheel we have $\pi/2 + \delta + \alpha = \pi$. So we have:

$$\delta_{circle} = \delta$$

And then:

$$\psi_{t+1} = \psi_t + (v_t/R) * dt$$

$$\psi_{t+1} = \psi_t + (v_t/L) * \tan(\delta_t) * dt$$

Note that for small δ_t we have $\tan(\delta_t) \approx \delta_t$

Speed update: assuming during dt that a_t is constant:

$$v_{t+1} = v_t + a_t * dt$$

So to summarize our kinematic bicycle model which is a **kinematic constraint model** is:

$$x_{t+1} = x_t + v_t * \cos(\psi_t) * dt$$

$$y_{t+1} = y_t + v_t * \sin(\psi_t) * dt$$

$$\psi_{t+1} = \psi_t + (v_t/L) * \tan(\delta_t) * dt$$

$$v_{t+1} = v_t + a_t * dt$$

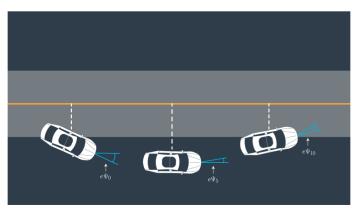
The state vector is $S_t = [x_t, y_t, \psi_t, v_t]$.

The actuator command $A_t = [a_t, \delta_t]$ defines a **constraint** between S_{t+1} and S_t .

Note also that for a bicycle we have $\delta_t \in [-\pi/2, \pi/2]$ whereas for a car we have $\delta_t \in (-\delta_{max}, \delta_{max})$ with $\delta_{max} < \pi/2$.

4.2.3 Errors

The errors variables are the following:



The dashed white line is the cross track error.

Figure 5: Model errors

- cte: cross track error. It corresponds to distance of vehicle from the planned trajectory (as planned by path planning module)
- $e\Psi$: psie error is the angle difference of the vehicle trajectory with the planned trajectory (as planned by path planning module)

The new state vector is $[x_t, y_t, \psi_t, v_t, cte_t, e\Psi_t]$.

Let's assume the reference trajectory provided by the Path Planning module is a set of waypoints approximated or fitted by a (3rd order typically) polynomial $f(x_t) = a_0 + a_1 x_t + a_2 x_t^2 + a_3 x_t^3$ and our real position on the y-axis is y_t

Compute current Cross Track Error:

The current Cross Track Error is:

$$cte_t = y_t - f(x_t)$$

Note: we are dealing with a trajectory which is a set of (x, y, t) or (x_t, y_t) coordinates. We are not dealing with a path which a set of (x, y) coordinates.

Compute current Orientation Error:

The current Orientation Error is:

$$e\Psi_t = \psi_t - \text{tangential angle of curve of f at } x_t = \psi_t - \arctan(f'(x_t))$$

Update Cross Track Error:

The Cross Track Error will increase during dt proportionnaly to the speed and Orientation error (along local y-axis):

$$cte_{t+1} = cte_t + v_t * sin(e\Psi_t) * dt$$

Update Orientation Error:

The update rule for $e\Psi_t$ is the same as for ψ_t :

$$e\Psi_{t+1} = e\Psi_t + (v_t/L) * \tan(\delta_t) * dt$$

As we have:

$$\psi_{t+1} = \psi_t + (v_t/L) * \tan(\delta_t) * dt$$
$$e\Psi_t = \psi_t - \arctan(f'(x_t))$$
$$e\Psi_{t+1} = \psi_{t+1} - \arctan(f'(x_{t+1}))$$

We get:

$$e\Psi_{t+1} = \psi_t + (v_t/L) * \tan(\delta_t) * dt - \arctan(f'(x_{t+1}))$$

$$e\Psi_{t+1} = \psi_t - \arctan(f'(x_{t+1})) + (v_t/L) * \tan(\delta_t) * dt$$

And assuming our tangential angle remains constant during dt our orientation error will evolve such that:

$$e\Psi_{t+1} = \psi_t - \arctan(f'(x_t)) + (v_t/L) * \tan(\delta_t) * dt$$

$$e\Psi_{t+1} = e\Psi_t + (v_t/L) * \tan(\delta_t) * dt$$

So to summarize our kinematic bicycle error model which is a **kinematic** error model constraint is:

$$cte_{t+1} = cte_t + v_t * sin(e\Psi_t) * dt$$

$$e\Psi_{t+1} = e\Psi_t + (v_t/L) * tan(\delta_t) * dt$$

The underlying model is an approximation. Better models exist and are used in practice. Also the smaller the dt considered, the better the approximation. But in any cases no models are perfect (they are just usefull) and there will anyways be a delta (the smaller the better) between where we predict the vehicle will go and where we will end up. As an extreme case consider we are driving over some ice or oil that was not modelled. So the model will be reactive and corrective. Commands for the actuators will be reevaluated at every timestep even if we do predictions and plan actuator commands over multiple timesteps.

4.2.4 Actuator Constraints

In a real vehicle, actuators are limited by the design of the vehicle and fundamental physics. For example, a vehicle can't have a steering angle of 90 degrees. It is impossible. Thus it does not make sense for us to even consider these kinds of inputs. There is actually a vocabulary for describing this constraint. We call this model **nonholomonic** because the vehicle can't move in any arbitrary directions. It's limited by steering angle constraints. We can solve this by setting lower and upper bounds for the actuators. We have such **actuator constraints** as:

$$\begin{array}{|c|c|}
\hline
-30^o \le \delta_t \le +30^o \\
-1 \le a_t \le +1
\end{array}$$

where -1 is full brake and +1 is full acceleration.

4.2.5 Constraints summary

We have now specified a set of constraints acounting for vehicle constraints. To summarize our kinematic model constraints are:

$$x_{t+1} = x_t + v_t * \cos(\psi_t) * dt$$

$$y_{t+1} = y_t + v_t * \sin(\psi_t) * dt$$

$$\psi_{t+1} = \psi_t + (v_t/L) * \tan(\delta_t) * dt$$

$$v_{t+1} = v_t + a_t * dt$$

$$cte_{t+1} = cte_t + v_t * \sin(e\Psi_t) * dt$$

$$e\Psi_{t+1} = e\Psi_t + (v_t/L) * \tan(\delta_t) * dt$$

And our actuator constraints are:

$$\begin{array}{|c|c|}
\hline
-30^o \le \delta_t \le +30^o \\
-1 \le a_t \le +1
\end{array}$$

So far these contraints are defined for a 1-step duration of dt milliseconds. But we can rollout these constraints over N-steps. Typically we may use N=10 and dt=20 ms.

4.3 Dynamic Models

Forces, Slip Angle, Slip ratio and Tire Models

5 Model Predictive Control

MPC reframes the task of following a trajectory as an optimization problem. The solution to the optimization problem is the optimal trajectory.

MPC involves simulating different actuator inputs, predicting the resulting trajectory and minimizing a set of constraints .

Once we found the lowest cost trajectory, we implement the very first set of actuation commands. Then we throw away the rest of the trajectory we calculated. Instead of using the old trajectory we predicted, we take our new state and use that to calculate a new optimal trajectory. In that sense, we are constantly calculating inputs over a future horizon. That's why this approach is also called Receding Horizon Control. We constantly reevaluate the trajectory because our vehicle model is not perfect and the

next predicted (or planned) state may (slightly...) differ with our prediction (in the sense of a consequence of a command sent).

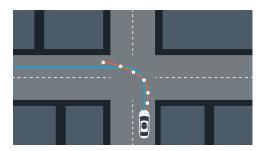


Figure 6: Minimization problem

5.1 Minimization of a cost function

We typically want to minimize things related to Tracking Error and comfort:

- Cross Track Error
- Orientation Error
- speed delta vs a target reference speed
- use of actuators (you prefer to use low values of a_t and δ_t as long as you can remain on track)
- limit the rate of change of actuator commands from one timestep to another

To account for these different objectives, we define different sub cost functions and weight them according to our priorities.

Listing 2: Cost function example

```
for (size_t t = 0; t < N; t++) {
    fg[0] += 4 * 2000 * CppAD::pow(vars[cte_start + t], 2);
    fg[0] += 4 * 2000 * CppAD::pow(vars[epsi_start + t], 2);
    fg[0] += CppAD::pow(vars[v_start + t] - ref_v, 2);
}

// Minimize the use of actuators.
for (size_t t = 0; t < N - 1; t++) {</pre>
```

```
fg[0] += 5 * CppAD::pow(vars[delta_start + t], 2);

fg[0] += 5 * CppAD::pow(vars[a_start + t], 2);

fg[0] += 5 * CppAD::pow(vars[a_start + t], 2);

// smooth
for (size_t t = 0; t < N - 2; t++) {
    fg[0] += 200 * CppAD::pow(vars[delta_start + t + 1] - vars[delta_start + t],

fg[0] += 10 * CppAD::pow(vars[a_start + t + 1] - vars[a_start + t],

fg[0] += 10 * CppAD::pow(vars[a_start + t],

fg[0] += 10 * CppAD::pow(vars
```

5.2 Timsestep length and Elapsed duration

N=10 and dt=100 ms are used so that we are working on 1 second of data. This is a trade-off: we need enough data visibility to ensure a good prediction, but we also have to limit the amount of computation. In general, smaller dt gives better accuracy, but that will require higher N for given horizon (N*dt). However, increasing N will result in longer computational time which increases the latency. The most common choice of values is N=10 and dt=0.1 but anything between N=20, dt=0.05 should work.

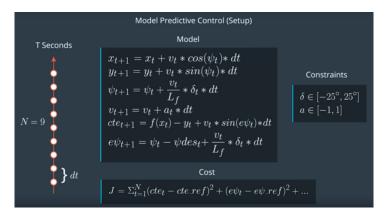


Figure 7: Solver setup with N*dt time horizon

5.3 Latency handling

A contributing factor to latency is actuator dynamics. For example the time elapsed between when you command a steering angle to when that angle is actually achieved. This could easily be modeled by a simple dynamic system and incorporated into the vehicle model. One approach would be running a simulation using the vehicle model starting from the current state for the duration of the latency. The resulting state from the simulation is the new initial state for MPC.

Thus, MPC can deal with latency much more effectively, by explicitly taking it into account, than a PID controller.

5.4 MPC Solver algorithm

To summarize we have defined a cost function we want to minimize under a set of constraints. The cost function accounts for what we want to produce ideally: a comfortable trajectory that follows a reference path as close as possible. Whereas the constraints account for the constraints we are dealing with: typically at this stage of the Autonomous Driving pipeline, we are dealing with vehicle model constraints. So our non-linear optimization problem is defined and can be summarized visually as:

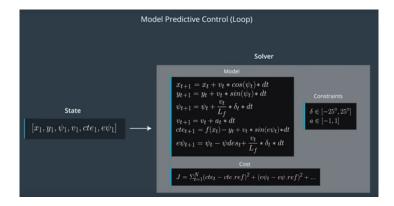


Figure 8: Solver input

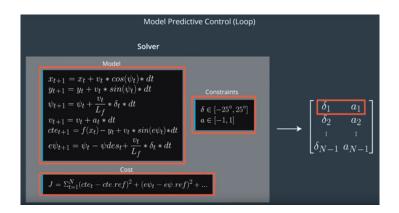


Figure 9: Solver output

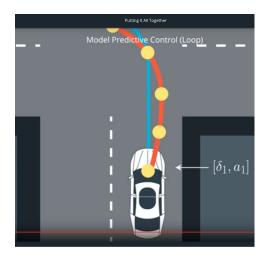


Figure 10: Solver actuator commands

The next step is to solve it in software. The model, a cost function to minimize and a set of constraints to comply with, will be described in software and a mathematical library will be used to solve this non-linear optimization problem.

5.5 MPC Solver code

Listing 3: MPC solver with ipopt

```
1 #include "MPC.h"
 2 #include <cppad/cppad.hpp>
 3 #include <cppad/ipopt/solve.hpp>
 4 #include "Eigen -3.3/Eigen/Core"
 5 #include "Eigen -3.3/Eigen/QR"
 7 using CppAD::AD;
9 // TODO: Set the timestep length and duration
10 size_t N = 10;
11 double dt = 0.1;
12
13 // This value assumes the model presented in the classroom is used.
14 //
15 // It was obtained by measuring the radius formed by running the vehicle in
16 // simulator around in a circle with a constant steering angle and velocity
17 // flat terrain.
18 //
19 // Lf was tuned until the the radius formed by the simulating the model
20 // presented in the classroom matched the previous radius.
21 //
22 // This is the length from front to CoG that has a similar radius.
23 const double Lf = 2.67;
24
25 // NOTE: feel free to play around with this
26
  // or do something completely different
27
28 \quad double \quad ref_v = 120;
29
30 // The solver takes all the state variables and actuator
31 // variables in a singular vector. Thus, we should to establish
```

```
32 // when one variable starts and another ends to make our lifes easier.
33 \text{ size_t } x_{\text{start}} = 0;
34 \text{ size_t y\_start} = \text{x\_start} + \text{N};
35 \text{ size\_t psi\_start} = \text{y\_start} + \text{N};
36 \text{ size_t } \text{v_start} = \text{psi_start} + \text{N};
37 size_t cte_start = v_start + N;
38 size_t epsi_start = cte_start + N;
39 size_t delta_start = epsi_start + N;
40 \text{ size\_t a\_start} = \text{delta\_start} + N - 1;
41
42 class FG_eval {
43
    public:
     // Fitted polynomial coefficients
44
45
     Eigen::VectorXd coeffs;
     FG_eval(Eigen::VectorXd coeffs) { this->coeffs = coeffs; }
46
47
48
      typedef CPPAD_TESTVECTOR(AD<double>) ADvector;
49
      void operator()(ADvector& fg, const ADvector& vars) {
50
        // TODO: implement MPC
51
        // 'fg' a vector of the cost constraints, 'vars' is a vector of variab
52
        // NOTE: You'll probably go back and forth between this function and
53
        // the Solver function below.
54
        // The cost is stored is the first element of 'fg'.
55
56
        // Any additions to the cost should be added to 'fg[0]'.
57
        fg[0] = 0;
58
59
        // Reference State Cost
        // TODO: Define the cost related the reference state and
60
        // any anything you think may be beneficial.
61
62
        for (size_t t = 0; t < N; t++) {
63
          fg[0] += 4 * 2000 * CppAD::pow(vars[cte_start + t], 2);
64
          fg[0] += 4 * 2000 * CppAD::pow(vars[epsi_start + t], 2);
65
          fg[0] \leftarrow CppAD::pow(vars[v_start + t] - ref_v, 2);
66
        }
67
68
        // Minimize the use of actuators.
        for (size_t t = 0; t < N - 1; t++) 
69
70
          fg[0] += 5 * CppAD::pow(vars[delta_start + t], 2);
71
          fg[0] += 5 * CppAD::pow(vars[a_start + t], 2);
```

```
72
        }
73
74
        // smooth
75
        for (size_t t = 0; t < N - 2; t++) 
76
          fg[0] += 200 * CppAD::pow(vars[delta_start + t + 1] - vars[delta_sta
77
          fg[0] += 10 * CppAD::pow(vars[a_start + t + 1] - vars[a_start + t],
78
79
80
81
        // Setup Constraints
82
83
        // NOTE: In this section you'll setup the model constraints.
84
85
        // Initial constraints
86
        // We add 1 to each of the starting indices due to cost being located
87
88
        // index 0 of 'fg'.
89
        // This bumps up the position of all the other values.
90
        fg[1 + x_start] = vars[x_start];
91
        fg[1 + y_start] = vars[y_start];
92
        fg[1 + psi_start] = vars[psi_start];
        fg[1 + v_start] = vars[v_start];
93
94
        fg[1 + cte_start] = vars[cte_start];
        fg[1 + epsi_start] = vars[epsi_start];
95
96
97
        // The rest of the constraints
        for (size_t t = 1; t < N; t++) {
98
99
          // at time t+1
          AD < double > x1 = vars[x_start + t];
100
          AD < double > y1 = vars[y_start + t];
101
102
          AD<double> psi1 = vars[psi_start + t];
103
          AD < double > v1 = vars[v_start + t];
104
          AD<double> cte1 = vars[cte_start + t];
105
          AD<double> epsi1 = vars[epsi_start + t];
106
107
          // at time t
108
          AD < double > x0 = vars[x_start + t - 1];
          AD < double > y0 = vars[y_start + t - 1];
109
110
          AD < double > psi0 = vars[psi_start + t - 1];
111
          AD < double > v0 = vars[v_start + t - 1];
```

```
112
          AD < double > cte0 = vars[cte\_start + t - 1];
113
          AD < double > epsi0 = vars[epsi_start + t - 1];
114
          AD < double > delta0 = vars[delta_start + t - 1];
115
116
          AD < double > a0 = vars[a_start + t - 1];
117
118
          // XXX: to be updated
119
          AD < double > f0 = coeffs[0] + coeffs[1] * x0 + coeffs[2] * x0 * x0 + c
120
          AD < double > psides0 = CppAD :: atan(coeffs[1] + 2 * coeffs[2] * x0 + 3
121
          // Here's 'x' to get you started.
122
123
          // The idea here is to constraint this value to be 0.
124
125
          // NOTE: The use of 'AD<double>' and use of 'CppAD'!
126
          // This is also CppAD can compute derivatives and pass
127
          // these to the solver.
128
129
          // TODO: Setup the rest of the model constraints
130
          fg[1 + x_start + t] = x1 - (x0 + v0 * CppAD:: cos(psi0) * dt);
          fg[1 + y_start + t] = y1 - (y0 + v0 * CppAD:: sin(psi0) * dt);
131
132
          fg[1 + psi_start + t] = psi1 - (psi0 - v0 * delta0 / Lf * dt); // XX
133
          fg[1 + v_start + t] = v1 - (v0 + a0 * dt);
134
135
          // BUG fg[1 + cte_start + t] = cte_1 - (cte_0 + v_0 * CppAD::sin(epsi_0))
136
          // BUG fg[1 + epsi_start + t] = epsi1 - (epsi0 + v0 * delta0 / Lf *
137
          fg[1 + cte_start + t] = cte_1 - ((f0 - y0) + (v0 * CppAD::sin(epsi0))
          fg[1 + epsi\_start + t] = epsi1 - ((psi0 - psides0) - v0 * delta0 / L
138
139
140
141 };
142
143 //
144 // MPC class definition implementation.
145 //
146 MPC::MPC() {}
147 MPC::~MPC() {}
148
149 vector<double> MPC::Solve(Eigen::VectorXd state, Eigen::VectorXd coeffs) {
150
      bool ok = true;
151
      size_t i;
```

```
152
      typedef CPPAD_TESTVECTOR(double) Dvector;
153
154
      double x = state[0];
155
      double y = state[1];
156
      double psi = state[2];
157
      double v = state[3];
      double cte = state [4];
158
159
      double epsi = state [5];
160
161
      // TODO: Set the number of model variables (includes both states and inp
      // For example: If the state is a 4 element vector, the actuators is a 2
162
      // element vector and there are 10 timesteps. The number of variables is
163
164
      // 4 * 10 + 2 * 9
165
      size_t n_vars = N * 6 + (N - 1) * 2;
166
      // TODO: Set the number of constraints
167
      size_t n_constraints = N * 6;
168
169
      // Initial value of the independent variables.
170
171
      // Should be 0 except for the initial values.
      Dvector vars (n_vars);
172
173
      for (i = 0; i < n_vars; i++)
174
        vars[i] = 0.0;
175
176
      // Set the initial variable values
177
      vars[x_start] = x;
178
      vars[y_start] = y;
      vars[psi_start] = psi;
179
180
      vars[v_start] = v;
      vars[cte_start] = cte;
181
182
      vars[epsi_start] = epsi;
183
184
      // Lower and upper limits for x
      Dvector vars_lowerbound(n_vars);
185
      Dvector vars_upperbound(n_vars);
186
187
      // TODO: Set lower and upper limits for variables.
      // Set all non-actuators upper and lowerlimits
188
      // to the max negative and positive values.
189
190
      for (i = 0; i < delta_start; i++)
191
        vars_lowerbound[i] = -1.0e19;
```

```
192
        vars_upperbound[i] = 1.0e19;
193
194
      // The upper and lower limits of delta are set to -25 and 25
195
196
      // degrees (values in radians).
      // NOTE: Feel free to change this to something else.
197
      for (i = delta\_start; i < a\_start; i++) {
198
        vars_lowerbound[i] = -0.436332 * Lf; // *Lf ? XXX
199
        vars\_upperbound[i] = 0.436332 * Lf; // *Lf?
200
201
      }
202
      // Acceleration/decceleration upper and lower limits.
203
204
      // NOTE: Feel free to change this to something else.
      for (i = a_start; i < n_vars; i++)
205
        vars_lowerbound[i] = -1.0;
206
207
        vars\_upperbound[i] = 1.0;
      }
208
209
210
211
      // Lower and upper limits for the constraints
      // Should be 0 besides initial state.
212
213
      Dvector constraints_lowerbound(n_constraints);
214
      Dvector constraints_upperbound(n_constraints);
215
      for (i = 0; i < n_{constraints}; i++)
216
        constraints\_lowerbound[i] = 0;
217
        constraints_upperbound[i] = 0;
218
      constraints_lowerbound [x_start] = x;
219
220
      constraints_lowerbound[y_start] = y;
      constraints_lowerbound[psi_start] = psi;
221
222
      constraints_lowerbound[v_start] = v;
      constraints_lowerbound[cte_start] = cte;
223
224
      constraints_lowerbound[epsi_start] = epsi;
225
226
      constraints_upperbound [x_start] = x;
227
      constraints_upperbound[v_start] = y;
      constraints_upperbound[psi_start] = psi;
228
      constraints_upperbound[v_start] = v;
229
230
      constraints_upperbound[cte_start] = cte;
      constraints_upperbound[epsi_start] = epsi;
231
```

```
232
233
      // Object that computes objective and constraints
      FG_eval fg_eval(coeffs);
234
235
236
      // NOTE: You don't have to worry about these options
237
238
      // options for IPOPT solver
239
      std::string options;
240
241
      // Uncomment this if you'd like more print information
      options += "Integer print_level 0 \ n";
242
      // NOTE: Setting sparse to true allows the solver to take advantage
243
244
      // of sparse routines, this makes the computation MUCH FASTER. If you
      // can uncomment 1 of these and see if it makes a difference or not but
245
      // if you uncomment both the computation time should go up in orders of
246
      // magnitude.
247
      options += "Sparse
248
                          {
m true}
                                       forward\n";
249
      options += "Sparse true
                                      reverse\n";
      // NOTE: Currently the solver has a maximum time limit of 0.5 seconds.
250
251
      // Change this as you see fit.
      options += "Numeric max_cpu_time
                                             0.5 \setminus n";
252
253
254
      // place to return solution
      CppAD::ipopt::solve_result < Dvector > solution;
255
256
257
      // solve the problem
258
      CppAD::ipopt::solve<Dvector, FG_eval>(
          options, vars, vars_lowerbound, vars_upperbound, constraints_lowerbound
259
260
          constraints_upperbound , fg_eval , solution );
261
262
      // Check some of the solution values
263
264
      ok &= solution.status == CppAD::ipopt::solve_result < Dvector >::success;
265
      // Cost
266
267
      auto cost = solution.obj_value;
268
      //std::cout << "Cost " << cost << std::endl;
269
270
      // TODO: Return the first actuator values. The variables can be accessed
      // 'solution.x[i]'.
271
```

```
272
      // {...} is shorthand for creating a vector, so auto x1 = \{1.0, 2.0\}
273
      // creates a 2 element double vector.
274
      //return {solution.x[delta_start], solution.x[a_start]};
275
276
277
      vector<double> result;
278
279
      result.push_back(solution.x[delta_start]);
      result.push_back(solution.x[a_start]);
280
281
282
      for (size_t i = 0; i < N - 1; i++)
        result.push_back(solution.x[x_start + i + 1]);
283
        result.push_back(solution.x[y_start + i + 1]);
284
      }
285
286
287
      return result;
288 }
```