

# Improvement of the Measurement Update Step of EKF-SLAM

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**Abstract**—In this study, the measurement update step of the Extended Kalman Filter (EKF)-based Simultaneous Localization and Mapping (SLAM) is improved. The computational complexity of the measurement uncertainty matrix inversion operation in the measurement update step is reduced via using Jacobi iteration method. It is observed that, the calculation of the measurement uncertainty matrix inverse by using Jacobi iteration method generates numerically more stable results than naive single and batch update operations. Moreover, it produces more accurate results than the results of Cholesky decomposition with less complexity.

## I. INTRODUCTION

The simultaneous localization and map building (SLAM) techniques tries to solve the problem for an autonomous vehicle to start in an unknown location in an unknown environment to incrementally build a map of this environment. The robot uses this map to compute its own location simultaneously. It can only sense noisy, probabilistic observations of its surroundings without knowing its exact location. During the robot moves, uncertainty is added into an already uncertain pose.

Among the methods proposed in literature to solve the SLAM problem, the methods based on Bayesian estimation theory has been the most successful ones. Many applications in the literature use Extended Kalman Filter (EKF) to solve nonlinear estimation problems such as position tracking, localization and SLAM [1-10]. However, the quadratic computational complexity of the EKF makes it difficult to apply in real time. Unscented Kalman Filter (UKF) is a more reliable estimator than EKF while the system model is highly nonlinear. The past of the UKF is relatively short compared to EKF. By approximating the probability density function, instead of the nonlinear function itself, UKF SLAM [11, 12] received a considerable attention. Yet it did not make any improvement to the computational complexity of the EKF. FastSLAM [13-17] utilizes particle filters and improves the computational complexity considerably compared to EKF and UKF. However in case of having insufficient particle numbers in the filter; FastSLAM can not estimate the state successfully.

In this study, reducing the complexity of the measurement update step of the EKF-based SLAM applications is aimed. Inverse of the measurement uncertainty matrix is calculated in the measurement update step of EKF. Minimum complexity of the matrix inversion operation is  $O(n^{2.4})$ , achieved by Coppersmith–Winograd algorithm.

We present the Jacobi iteration method instead of matrix inversion process in EKF to reduce computational complexity. In EKF-based SLAM applications the innovation vector, the difference between the predicted and the actual observations, is taken into account for the state estimation via using batch or single update for each observation. The convergence of the measurement uncertainty matrix inversion by Jacobi iteration method produces numerically more stable and accurate results than the results of single and batch update approaches which use Cholesky decomposition with less complexity.

The measurement uncertainty matrix, whose inverse is calculated by Jacobi iteration method, is a symmetric and positive-definite matrix. The computational complexity of the inversion calculation of that special matrix by Jacobi iteration method is  $O(N)$ , while the size of the measurement uncertainty matrix is  $N \times N$ . However, the iterative solution of Jacobi method is converged to the inverse of the interested matrix in 8 of 10 experiments. Each of the EKF-based SLAM approaches with single and batch Cholesky updates, those use naive matrix inversion calculation; and batch update, that uses matrix inversion by Jacobi iteration, were run 10 times. The averages of the Root Mean Square Error (RMSE) results of the approaches are given in the experimental results.

In the second section of the paper, the EKF algorithm is described. In Third section the single and batch Cholesky updates and complexity of them are given. In forth section, the matrix inversion by Jacobi iteration method and the complexity analysis of the EKF algorithm which uses Jacobi iteration method for matrix inversion process are introduced. In the Fifth section the experimental results are illustrated and conclusion is given in the final section.

## II. EKF-BASED SLAM ALGORITHM

EKF-based SLAM algorithm bases on Bayes Filters. The state of the system, the state of the robot and the environment, at time  $t$  is expressed by random variables  $x_t$ . In the probabilistic SLAM methods, the state of the robot and the environment can only be expressed through the conditional probability distributions of the sensor data. The probability distribution representing the uncertainty at each point in the time is called *belief*,  $bel(x_t)$ . Bayes filters apply two update rules successively to estimate the system state [8].

The predictive belief at time  $t$  is calculated just before the observation ( $z_t$ ) is taken and uses the control data ( $u_{1:t}$ ) by the time  $t$  [8]. This step is also called control update, and expressed as in (1).

$$\overline{bel}(x_t) = P(x_t | z_{1:t-1}, u_{1:t}) \quad (1)$$

The state estimate given in (1) is corrected according to (2), using sensor measurements ( $z_{1:t}$ ) and the control data ( $u_{1:t}$ ) by the time  $t$ . This step is called as measurement update or the posterior belief of the system state and calculated whenever a sensor provides a new observation [8].

$$bel(x_t) = P(x_t | z_{1:t}, u_{1:t}) \quad (2)$$

Kalman filters are the most widely used variant of Bayes filters. Standard Kalman filter assumes that the state transitions and measurement transitions are linear with added Gaussian noise. EKF overcomes the linearity assumption by linearizing the nonlinear state and measurement transition functions via Taylor expansion.

The nonlinear state transition function  $g$  and measurement transition function  $h$  is expressed as in (3) and (4) respectively with added noises  $\varepsilon_t$  and  $\delta_t$ .

$$x_t = g(u_t, x_{t-1}) + \varepsilon_t \quad (3)$$

$$z_t = h(x_t) + \delta_t \quad (4)$$

EKF represents beliefs by the mean vector  $\mu_t$  and the covariance matrix  $\Sigma_t$  at time  $t$ .

The Taylor expansion of state transition function  $g$  is given as in (5) and (6). The Jacobian  $G_t$  is the value of first derivative of  $g$  at the point  $\mu_{t-1}$  [8].

$$g(\mathbf{u}_t, \mathbf{x}_{t-1}) \approx g(\mathbf{u}_t, \mu_{t-1}) + g'(\mathbf{u}_t, \mu_{t-1})(\mathbf{x}_{t-1} - \mu_{t-1}) \quad (5)$$

$$g(\mathbf{u}_t, \mathbf{x}_{t-1}) = g(\mathbf{u}_t, \mu_{t-1}) + G_t(\mathbf{x}_{t-1} - \mu_{t-1}) \quad (6)$$

Similarly, the Taylor expansion of measurement transition function  $h$  is given as in (7) and (8). The Jacobian  $H_t$  is the value of first derivative of  $h$  at the point  $\mu_t$  [8].

$$h(\mathbf{x}_t) \approx h(\overline{\mu}_t) + h'(\overline{\mu}_t)(\mathbf{x}_t - \overline{\mu}_t) \quad (7)$$

$$h(\mathbf{x}_t) = h(\overline{\mu}_t) + H_t(\mathbf{x}_t - \overline{\mu}_t) \quad (8)$$

The update rules corresponding to the prediction and correction steps of the Bayes filter and other details of the EKF algorithm can be found in [8-10].

In this study observation has two-dimensions, which are called the range and the bearing information of a laser scanner. Range is the distance between the robot and obstacles (landmarks); bearing is the angular difference between the robot and the obstacle. After taking the observation  $z_t$ , data association procedure is done to detect whether the measurements belong to an already existing landmark or to a new landmark. The measurements that belong to already existing landmarks cause update process in the filter. The single and batch update processes differ in terms of the calculating the innovation vectors and evaluating the effects of the associated landmarks on the update process. In the single update, the interested calculations and evaluations of the filter are handled one-by-one for each measurement value. In the batch update, they are handled together.

The pseudo code of the EKF-based SLAM algorithm [18] is given as follows:

1.  $[\mu_t, \Sigma_t] = \text{Predict}(\mu_{t-1}, \Sigma_{t-1})$
2.  $z_t = \text{Get observations}()$
3.  $[z_f, z_n] = \text{Data associate}(\overline{\mu}_t, \overline{\Sigma}_t, z_t, \mathbf{R})$
4.  $[\mu_t, \Sigma_t] = \text{Update Map}(\mu_t, \Sigma_t, z_f, \mathbf{R})$
5.  $[\mu_t, \Sigma_t] = \text{Augment Map}(\mu_t, \Sigma_t, z_n, \mathbf{R})$

where  $\mathbf{R}$  is the covariance matrix of the observation noise;  $z_f$  is the vector of measurements that belong to associated (already existing) landmarks;  $z_n$  is the vector of measurements that belong to unseen landmarks.

### III. SINGLE AND BATCH UPDATE IN EKF ALGORITHM

If the effects of the measurements of the associated landmarks on the update process of the filter are taken into account one-by-one for each measurement, it is called single update process; if they are handled together, it is called batch update. Pseudo codes and complexity analysis of two updating approaches are given in the following subsections.

#### A. Single Update Approach

The associated measurements are handled one-by-one in the single update process. The difference between the actual and the predicted measurements is called innovation vector. The size of the innovation vector is two for only one measurement, because each measurement has two-dimensions, which are called range and bearing information. The measurement estimation uncertainty matrix is a  $2 \times 2$  matrix for only one measurement.

While  $N$  is the number of associated measurements, pseudo code of the single update approach:

1. For  $i=1$  to  $N$ 
  - a.  $[z_p, H_t] = h(\overline{\mu}_t) // h$  is the observ. model
  - b.  $v = z_p - z_f // v$  is the innovation vector

- c.  $\mathbf{S}_t = \mathbf{H}_t \bar{\Sigma}_t \mathbf{H}_t^T + \mathbf{R}$  // meas. uncert. matrix
- d.  $\mathbf{K}_t = \bar{\Sigma}_t \mathbf{H}_t^T \mathbf{S}_t^{-1}$  // Kalman gain
- e.  $\boldsymbol{\mu}_t = \boldsymbol{\mu}_t + \mathbf{K}_t \mathbf{v}$  // update of the mean vec.
- f.  $\bar{\Sigma}_t = (\mathbf{I} - \mathbf{K}_t \mathbf{H}_t) \bar{\Sigma}_t$  // update of the cov.

2. End for

The Kalman gain,  $\mathbf{K}_t$ , is calculated by using the inverse of the measurement uncertainty matrix,  $\mathbf{S}_t$ . To calculate the inverse of  $\mathbf{S}_t$ ; Cholesky decomposition can be used as in [18].

The complexity of the inversion of a  $2 \times 2$  matrix is  $O(2^{2.4})$ ; therefore the computational complexity of the  $\mathbf{S}_t$ 's inversion operation is  $O(Nx2^{2.4})$  in the single update for all observations, while  $N$  is the number of the associated measurements.

### B. Batch Update Approach

The associated measurements are handled together within a joint matrix in the batch update process. The innovation vector is a  $2Nx1$  vector when the number of the associated measurements is  $N$ , because each measurement has two-dimensions (features). And also the measurement estimation uncertainty matrix is a  $2Nx2N$  matrix.

Pseudo code of the batch update approach:

1. For  $i=1$  to  $N$ 
  - a.  $[\mathbf{z}_p, \mathbf{H}_t(2i-1:2i, 2i-1:2i)] = h(\bar{\boldsymbol{\mu}}_t)$
  - b.  $\mathbf{v}(2i-1:2i) = \mathbf{z}_p - \mathbf{z}_f$  //  $\mathbf{v}$  is innovation vec.
  - c.  $\mathbf{RR}(2i-1:2i, 2i-1:2i) = \mathbf{R}$
2. End for
3.  $\mathbf{S}_t = \mathbf{H}_t \bar{\Sigma}_t \mathbf{H}_t^T + \mathbf{RR}$  // meas. uncertainty matrix
4.  $\mathbf{K}_t = \bar{\Sigma}_t \mathbf{H}_t^T \mathbf{S}_t^{-1}$  // Kalman gain
5.  $\boldsymbol{\mu}_t = \boldsymbol{\mu}_t + \mathbf{K}_t \mathbf{v}$  // update of the mean vec.
6.  $\bar{\Sigma}_t = (\mathbf{I} - \mathbf{K}_t \mathbf{H}_t) \bar{\Sigma}_t$  // update of the cov.

To calculate the inverse of the measurement uncertainty matrix,  $\mathbf{S}_t$ ; Cholesky decomposition can be used as in [18].

The optimal complexity of the inverse calculation of a  $2Nx2N$  matrix is  $O((2N)^{2.4})$ ; hence the computational complexity of the  $\mathbf{S}_t$ 's inversion operation is  $O(N^{2.4}x2^{2.4})$  in the batch update for whole observations.

## IV. MATRIX INVERSION BY JACOBI ITERATION METHOD

The linear equation system, which has  $N$  variables and  $N$  equations, is given in (9) [19]:

$$\mathbf{Ax} = \mathbf{b} \quad (9)$$

The solution of (9) is obtained as,

$$\mathbf{x} = \mathbf{A}^{-1}\mathbf{b}. \quad (10)$$

If  $\mathbf{b}$  is chosen as an unit matrix  $\mathbf{I}$ , the solution of (9) is the inverse of  $\mathbf{A}$ . Due to the computational load involved in taking the inverse of  $\mathbf{A}$ , rearrange (9) as

$$(\mathbf{D} + \mathbf{L} + \mathbf{U})\mathbf{x} = \mathbf{I}, \quad (11)$$

where  $\mathbf{D}$ ,  $\mathbf{L}$  and  $\mathbf{U}$  are the diagonal matrix, lower and upper triangular matrices respectively, yields

$$\mathbf{x} = \mathbf{D}^{-1}[(\mathbf{D} - \mathbf{A})\mathbf{x} + \mathbf{I}]. \quad (12)$$

Now assign initial matrix  $\mathbf{x}_0$  to the  $\mathbf{x}$  on the right-hand side and calculate  $\mathbf{x}_1$  on the left-hand side. in general case, we can write

$$\mathbf{x}_{k+1} = \mathbf{D}^{-1}[(\mathbf{D} - \mathbf{A})\mathbf{x}_k + \mathbf{I}]. \quad (13)$$

If  $\mathbf{A}$  is a dominant matrix, the absolute values of the eigenvalues of the matrix  $\mathbf{D}^{-1}(\mathbf{D} - \mathbf{A})$  are less than 1. When this condition exists,  $\mathbf{x}_k$  approaches the inverse of  $\mathbf{A}$ . A dominant matrix has the absolute value of the diagonal entry in a row which is larger than or equal to the sum of the absolute values of all the other entries in that row. If that condition is satisfied, inverse of the matrices, which have even large size; can be calculated in a few iterations by using Jacobi iteration method [19].

The measurement uncertainty matrices, those are generated in this study, are symmetric and positive-definite matrices. They also provide the above conditions generally. The sufficient iteration number is  $M$  for an  $M \times M$  matrix to converge the solution by Jacobi iteration method in this study.

If Jacobi iteration method is used for the matrix inversion in the single update, the complexity is  $O(2N)$  where  $N$  is the number of the associated measurements.  $O(2)$  is the complexity of calculating the inverse of a  $2 \times 2$  matrix by Jacobi iteration method. This is a bit smaller than the complexity of the naive single update, it is not significant. However in the naive batch update for all of the measurements, calculating the inverse of the measurement uncertainty matrix costs  $O(N^{2.4}x2^{2.4})$ , while it costs  $O(2N)$  by using Jacobi iteration method for matrix inversion in the batch update. It is significantly less than the complexity of naive batch update.

All of the operations, except the inverse matrix calculation, are the same in the naive update approaches and the approaches which use Jacobi iteration method for inverse matrix calculation. The complexity results for both single and batch update are the same,  $O(2N)$ , when Jacobi iteration method is used for matrix inversion. The difference between the single updates, which uses and does not use Jacobi iteration method, is not significant. Hence the results of the single update that uses Jacobi iteration method are not given in the experimental results.

TABLE I  
COMPLEXITY ANALYSIS OF THE APPROACHES

Approach	Single Cholesky update	Batch Cholesky update	Batch update that uses Jacobi iteration method for matrix inverse
Complexity	$O(2^{2.4}xN)$	$O(M^{2.4})=O(2^{2.4}xN^{2.4})$	$O(M) = O(2xN)$



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