Robust Online Model Adaptation by Extended Kalman Filter with Exponential Moving Average and Dynamic Multi-Epoch Strategy

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Abstract
High fidelity behavior prediction of intelligent agents is critical in many applications. However, the prediction model trained on the training set may not generalize to the testing set due to domain shift and time variance. The challenge motivates the adoption of online adaptation algorithms to update prediction models in real-time to improve the prediction performance. Inspired by Extended Kalman Filter (EKF), this paper introduces a series of online adaptation methods, which are applicable to neural network-based models. A base adaptation algorithm Modified EKF with forgetting factor (MEKF$_\lambda$) is introduced first, followed by exponential moving average filtering techniques. Then this paper introduces a dynamic multi-epoch update strategy to effectively utilize samples received in real time. With all these extensions, we propose a robust online adaptation algorithm: MEKF with Exponential Moving Average and Dynamic Multi-Epoch strategy (MEKF$_{EMA-DME}$). The proposed algorithm outperforms existing methods as demonstrated in experiments. The source code is open-sourced in the following link https://github.com/intelligent-control-lab/MEKF_MAME.

Keywords: Online adaptation, extended Kalman filter, exponential moving average, optimization

1. Introduction
Supervised learning has been widely used to obtain models to predict the behaviors of intelligent agents Rudenko et al. (2019). Behavior prediction is a sub-topic of time series prediction Weigend (2018), which includes but is not limited to vehicle trajectory prediction during autonomous driving Lefe`vre et al. (2014) and human-motion prediction during human-robot collaboration Cheng et al. (2019). Although a trained model typically performs well on the training set, performance can drop significantly in a slightly different test domain or under a slightly different data distribution Si et al. (2019); Callison-Burch et al. (2010). For tasks without annotated corpora from the test domain, adaptation techniques are required to deal with the lack of domain-specific data. This paper studies robust online adaptation algorithms for behavior prediction.

In online adaptation, a prediction model observes instances sequentially over time. After every observation, the model outputs a prediction and receives the ground truth. Then the online adaptation algorithm updates the prediction model according to the error measured between the prediction and the ground truth. The goal of adaptation is to improve the prediction accuracy in subsequent rounds. An online adaptation algorithm is robust if it can efficiently adapt an existing model to a different (test) data distribution, without generating big transient errors.

For prediction models encoded in neural networks, most existing online adaptation approaches are based on stochastic gradients Kivinen et al. (2004). For example, the identification-based ap-
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Approach uses stochastic gradient descent (SGD) to adapt the model online Bhasin et al. (2012). However, these methods may be sub-optimal in minimizing the local prediction errors. Another solution is to use the recursive least square parameter adaptation algorithm (RLS-PAA) Ljung and Priouret (2010), which has been applied to adapt the last layer of a feedforward neural network Cheng et al. (2019) or the last layer of a recurrent neural network Si et al. (2019). RLS-PAA can only adapt the last layer of a neural network since it only applies to linear models. To adapt other layers, the adaptation problem becomes nonlinear, which requires the development of robust optimal nonlinear adaptation algorithms Jazwinski (2007); Cooper et al. (2014); Abuduweili et al. (2019).

Since a neural network parameterizes a nonlinear system with a layered structure, learning or adaptation of the neural network is equivalent to parameter estimation of the nonlinear system. The extended Kalman filter (EKF) is one promising method for nonlinear parameter estimation Jazwinski (2007), which is derived by linearizing the system equations at each time step and applying Kalman filter (an optimal filter that minimizes the tracking error) on the linearized system. The EKF approach has been demonstrated to be superior to the SGD-based algorithms in training feedforward neural networks Iiguni et al. (1992); Ruck et al. (1992). Nonetheless, in online adaptation, more recent data is more important Fink et al. (2001). Similar to adaptive EKF methods Yang et al. (2006); Ozbek and Efe (2004); Anderson and Moore (2012) that discount old measurements, this paper considers the Modified Extended Kalman Filter with forgetting factor, MEKF\(_\lambda\), as a base adaptation algorithm.

On top of the base adaptation algorithm, the following modifications are made. Generally, the step size of parameter update in EKF-based approaches may not be optimal, due to the error introduced during linearization. Inspired by exponential moving average (EMA) methods, this paper proposes EMA filtering to the base MEKF\(_\lambda\) in order to increase the convergence rate. The resulting algorithm is called MEKF\(_{EMA}\). Then in order to effectively utilize the samples in online adaptation, this paper proposes a dynamic multi-epoch update strategy to discriminate the “hard” samples from “easy” samples, and sets different weights for them. The dynamic multi-epoch update strategy can improve the effectiveness of online adaptation with any base optimizers, e.g., SGD or MEKF\(_{EMA}\). By incorporating MEKF\(_{EMA}\) with the dynamic multi-epoch update strategy, we propose the algorithm MEKF\(_{EMA-DME}\) (MEKF with Exponential Moving Average and Dynamic Multi-Epoch update strategy).

The remainder of the paper first formulates the online adaptation problem, then discusses the proposed algorithm, and finally validates the effectiveness and flexibility of the proposed algorithms.

2. Online adaptation framework

The behavior prediction problem is to make inference on the future behavior of the target agent given the past and current measurement of the target agent and its surrounding environment. The transition model for behavior prediction problem is formulated as

\[ Y_t = f(\theta, X_t), \]

where the input vector \( X_t = [x_t; x_{t-1}; \cdots; x_{t-n+1}] \) denotes the stack of \( n \)-step current and past measurements (e.g. trajectory of states or extracted features) at time steps \( t, t-1, \ldots, t-n+1 \). The output vector \( Y_t = [y_{t+1}; y_{t+2}; \cdots; y_{t+m}] \) denotes the stack of the \( m \)-step future behavior (e.g. future trajectory) at time steps \( t+1, t+2, \ldots, t+m \). The function \( f \) is the prediction model that maps the measurements to the future behavior. \( \theta \) denotes the (ground truth) parameter of the model. It is assumed that there are recurrent structures in \( f \) such that the prediction of \( Y_t \) is made
by rolling out the one-step predictions \(y_{t+1} = f_1(\theta, X_t)\). The function \(f_1\) is the one-step prediction function and is a recurrent part of the overall prediction model.

Online adaptation explores local overfitting to minimize the prediction error. At time step \(t\), the following prediction error is to be minimized

\[
\min_{\hat{\theta}} \| Y_t - f(\hat{\theta}_t, X_t) \|_p,
\]

where \(Y_t\) is the ground truth trajectory (to be observed in the future) and \(\hat{Y}_t := f(\hat{\theta}_t, X_t)\) is the predicted trajectory using the estimated model parameter \(\hat{\theta}_t\). The adaptation objective can be in any \(\ell_p\) norm. This paper considers \(\ell_2\) norm. Assume that the true model parameter changes slowly during adaptation, i.e., \(\dot{\theta} \approx 0\). Then the estimated model parameter that minimizes the prediction error in the future can be approximated by the estimated parameter that minimizes the fitting error in the past. Solving for the estimated parameter that minimizes the fitting error corresponds to a nonlinear least square (NLS) problem.

**Definition 1 (Problem NLS)** Given a dataset \(\{(X_i, Y_i), i = 1, 2, \cdots, T\}\), find \(\hat{\theta}_t \in \mathbb{R}^n\) that minimizes

\[
J_t(\hat{\theta}_t) = \frac{1}{2} \sum_{i=1}^{T} \| e_i \|_2^2,
\]

where error term is defined as \(e_i = Y_i - f(\hat{\theta}_t, X_i)\).

In online adaptation, the estimate of the model parameter is updated iteratively when new data is received. Then a new prediction is made using the new estimate. In the next time step, the estimate will be updated again given the new observation and the process repeats. It is worth noting that the observation we received at time \(t\) is \(y_t\). The other terms in \(Y_t\) remain unknown. This paper is focused on adaptation methods using only one-step observation. It is possible to adapt with multi-step observations, which will be studied in the future. The process for online adaptation is summarized in algorithm 1. \(\hat{\theta}_t\) is the estimate of the model parameter \(\theta\) at time \(t\).

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**Algorithm 1** Generic Online Adaptation (Adaptable Prediction)

**Input:** Initial model parameters \(\hat{\theta}_0\)

**Output:** sequence of predictions \(\{\hat{Y}_t\}_{t=1}^T\)

1: for \(t = 1, 2, \cdots, T\) do
   2: obtain the ground truth observation value \(y_t\); construct the input features \(X_t\)
   3: adaptation step (supervised): \(\hat{\theta}_t = \text{Adapt}(\hat{\theta}_{t-1}, \hat{y}_t, y_t)\)
   4: prediction step: \(\hat{Y}_t = f(\hat{\theta}_t, X_t)\), where \(\hat{Y}_t = [\hat{y}_{t+1}; \hat{y}_{t+2}; \cdots; \hat{y}_{t+m}]\)
5: end for

6: return sequence of predictions \(\{\hat{Y}_t\}_{t=1}^T\)

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### 3. Robust nonlinear adaptation algorithms

#### 3.1. Modified EKF with forgetting factor

Our base adaptation algorithm is inspired by the recursive EKF method Moriyama et al. (2003); Alessandri et al. (2007). In EKF, the object being estimated is the state value of a dynamic system, while in adaptable prediction, the object to be adapted is the parameters that describe the system dynamics. Nonetheless, we can apply the EKF approach to adapt model parameters by regarding model parameters as system states. By assuming that the ground truth \(\theta\) changes very slowly, we can pose the parameter adaptation problem as a static state estimation problem Ruck et al. (1992); Nelson (2000) with the following dynamics,

\[
\dot{\hat{\theta}}_t = \hat{\theta}_{t-1} + \omega_t, \quad \omega_t \\
y_t = f_1(\hat{\theta}_{t-1}, X_{t-1}) + u_t
\]
where $\hat{\theta}_t$ is an estimate of the (ground truth) model parameter $\theta$; $y_t$ is the observation at time $t$; and $\hat{y}_t = f_1(\hat{\theta}_{t-1}, X_{t-1})$ is the prediction for time step $t$ made at time $t-1$. $f_1$ is the one-step prediction function. The injected process noise $\omega_t \sim \mathcal{N}(0, Q_t)$ and the injected measurement noise $u_t \sim \mathcal{N}(0, R_t)$ are assumed to be zero mean Gaussian white noise, and are identically and independently distributed. The symbol $\mathcal{N}$ represents Gaussian distribution. $Q_t$ and $R_t$ represent the covariance matrices for the process noise and the measurement noise respectively, which should be positive semidefinite. If there is no knowledge about the cross correlation of the noises, it is reasonable to assume that the entries in the noise vector are independent of each other, and set $Q_t$ and $R_t$ to be proportional to the identity matrix. For simplicity, we assume $Q_t = \sigma_q I$ and $R_t = \sigma_r I$ for $\sigma_q \geq 0$ and $\sigma_r > 0$.

In online adaptation, we assume that data in the distant past is no longer relevant for modeling the current dynamics, i.e. more recent data is more important. Hence, we consider a weighted nonlinear recursive least squares (NLS) problem:

$$\min_{\hat{\theta}_t} \sum_{i=1}^{t} \lambda^{t-i} ||y_i - f_1(\hat{\theta}_{i-1}, X_{i-1})||_2^2, \quad 0 < \lambda \leq 1, \quad (5)$$

where $\lambda$ is the “forgetting factor” which provides exponential decay to older samples. The forgetting factor prevents the EKF from saturation, and increases the algorithm’s ability to track a changing system. Algorithm 2 summarizes the modified extended Kalman filter algorithm with forgetting factor ($\text{MEKF}_\lambda$).

**Algorithm 2** Modified EKF algorithm with forgetting factor ($\text{MEKF}_\lambda$)

**Input:** Initial hyper-parameter for MEKF $\lambda$: $p_0 > 0$, $\lambda > 0$, $\sigma_r > 0$, $\sigma_q \geq 0$; $P_0 = p_0 I$

**Input:** previous parameter $\hat{\theta}_{t-1}$, previous prediction $\hat{y}_t = f_1(\hat{\theta}_{t-1}, X_{t-1})$, current observation $y_t$ at time step $t$

**Output:** Adapted parameter $\hat{\theta}_t$

1: $H_t = \frac{\partial f_1(\hat{\theta}, X)}{\partial \hat{\theta}}|_{\hat{\theta} = \hat{\theta}_{t-1}, X = X_{t-1}}$
2: $K_t = P_{t-1} \cdot H_t^T \cdot (H_t \cdot P_{t-1} \cdot H_t^T + \sigma_r I)^{-1}$
3: $\hat{\theta}_t = \hat{\theta}_{t-1} + K_t \cdot (y_t - f_1(\hat{\theta}_{t-1}, X_{t-1}))$
4: $P_t = \lambda^{-1}(P_{t-1} - K_t \cdot H_t \cdot P_{t-1} + \sigma_q I)$

In algorithm 2, $K_t$ is the Kalman gain. $P_t$ is a matrix representing the uncertainty in the estimates of the model parameter $\theta$. $H_t$ is the gradient matrix by linearizing the network. In online adaptation, $\theta_0$ is initialized by the offline trained parameter of the model. For $P_0$, due to the absence of a priori information, the $P_0$ matrix can be set to be proportional to the identity matrix, i.e. $P_0 = p_0 I$ for $p_0 > 0$.

**3.2. Extensions with exponential moving average filtering**

In the following discussion, for simplicity, an optimizer (e.g. MEKF$\lambda$) that solves the adaptation problem will be denoted as $A_P$ with internal state matrix $P$. The optimization process for adaptation can be compactly written as

$$\hat{\theta}_t = \hat{\theta}_{t-1} + V_t,$$

$$V_t = A_P(\hat{\theta}_{t-1}, \hat{y}_t, y_t). \quad (6)$$

where $V_t$ is the step size of the parameter update at time step $t$. 


To speed up the convergence of parameter estimation, we propose to apply exponentially-decayed moving average (EMA) for filtering in the MEKF$_{\lambda}$ optimization process. In SGD-based methods, numerous variants of EMA have been successfully used to speed up the convergence, including Polyak averaging Polyak (1964) and momentum Qian (1999). For MEKF$_{\lambda}$, we can either apply EMA on the step size $V$, to be discussed as EMA-V or momentum; or apply EMA on the optimizer’s inner state $P$, to be discussed as EMA-P.

**EMA-V**  EMA-V or momentum is widely used in SGD-based optimization algorithms Qian (1999), which helps accelerate gradient-based optimizers in relevant directions and dampen oscillations Qian (1999). Momentum can be regarded as an EMA filter on the step size of parameter update. It calculates the step size $V_t$ by decreasing exponentially the older step size with a factor $\mu_v \in [0, 1]$, i.e. $V_t = \mu_v V_{t-1} + (1 - \mu_v) A_P(\theta_{t-1}, \hat{y}_t, y_t)$.

**EMA-P**  As mentioned earlier in MEKF$_{\lambda}$, $P_t$ is a matrix representing the uncertainty in the parameter estimates. In order to attenuate instability during adaptation caused by anomaly data, we can smooth the inner state of the optimizer by pre-filtering $P_t$. The principle of pre-filtering the inner state (e.g., gradient, adaptive learning rate) before using them in optimization is applicable to many optimization algorithms. For example, in Adam Kingma and Ba (2014), the estimate of the first and second moment is filtered every step using EMA. Similarly, we can apply EMA on the inner state matrix $P_t$ of MEKF$_{\lambda}$.

By combining EMA-V and EMA-P, we propose the modified extended Kalman filter with exponential moving average (MEKF$_{\text{EMA}}$) algorithm as shown in algorithm 3, where $\mu_v$ is a momentum factor, and $\mu_p$ is a decay factor for the EMA filtering of $P_t$.

### Algorithm 3 Modified Extended Kalman Filter with Exponential Moving Average Filtering

**Input:** Initial base hyper-parameter: $p_0 > 0$, $\lambda > 0$, $\sigma_r > 0$, $\sigma_q \geq 0$; $P_0 = p_0 I$

**Input:** Initial EMA hyper-parameter: $0 \leq \mu_v < 1$, $0 \leq \mu_p < 1$

**Input:** previous parameter $\hat{\theta}_{t-1}$, previous prediction $y_{t-1} = f_1(\hat{\theta}_{t-1}, X_{t-1})$, current observation $y_t$ at time $t$

**Output:** Adapted parameter $\hat{\theta}_t$

1. $H_t = \frac{\partial f_1(\theta, x)}{\partial \theta} |_{\theta = \hat{\theta}_{t-1}, x = X_{t-1}}$
2. $K_t = P_{t-1} \cdot H_t^T \cdot (H_t \cdot P_{t-1} \cdot H_t^T + \sigma_r I)^{-1}$
3. $V_t = \mu_v V_{t-1} + (1 - \mu_v) K_t \cdot (y_t - f_1(\hat{\theta}_{t-1}, X_{t-1}))$
4. $\hat{\theta}_t = \hat{\theta}_{t-1} + V_t$
5. $P_t = \lambda^{-1}(P_{t-1} - K_t \cdot H_t \cdot P_{t-1} + \sigma_q I)$
6. $P_t = \mu_p P_{t-1} + (1 - \mu_p) P_t^*$

### 3.3. Dynamic multi-epoch update strategy

In generic online adaptation, all data are equally considered. We run the adaptation algorithm chronologically from the first data $X_1$ to the last data $X_T$. Every data sample is used only once, as shown in algorithm 1. We call the adaptation method that uses every data sample only once as **single-epoch online update strategy**.

Inspired by curriculum learning Bengio et al. (2009) in offline training, we introduce a more effective way to determine the adaptation epochs for every data sample during online adaptation. A curriculum can be viewed as a sequence of training criteria. Each training criterion in the sequence is associated with a different set of weights on the training examples. That said, it is practically
Definition 2 (Dynamic multi-epoch online update strategy) In online adaptation, the predicted output $\hat{y}_t$ generated by the estimated parameter $\theta^*$ is $\hat{y}_t = f_1(\theta^*, X_{t-1})$. Define a criterion $C$ to determine the number of epochs $\kappa_t \ (\kappa_t \in \mathbb{N})$ to adapt the parameter with the current sample, i.e., $\kappa_t = C(x_{t-1}, y_t, \hat{y}_t, \theta^*)$. In other words, we reuse the input-output pair $(X_{t-1}, y_t)$ $\kappa_t$ times to adapt the parameter $\theta^*$. This approach is called the dynamic multi-epoch online update strategy or dynamic multi-epoch update.

We propose a very straightforward criterion $C$ to determine the number of epochs $\kappa_t$ for every sample, as shown in algorithm 4. Two thresholds $\xi_1$ and $\xi_2$ are used to discriminate “easy”, “hard”, and “anomaly” samples. Before updating the parameter, we calculate the prediction error $j_t = \|y_t - \hat{y}_t\|_2$ at the current step. If the error satisfies $j_t < \xi_1$, the sample is considered as an “easy” sample. Then we run single-epoch update for this sample. If the error satisfies $\xi_1 \leq j_t < \xi_2$, the sample is considered as a “hard” sample. Then we reuse this sample and run the adaptation twice. The rationale is that for a “hard” sample, an adaptation optimizer may not learn enough under single-epoch update. If the error satisfies $j_t \geq \xi_2$, the sample is considered as an “anomaly” sample. Then we skip the update of $\hat{\theta}_t$. The rationale is that if the cost is too high, the sample is likely to be an anomaly point in the data distribution, which may destabilize the model adaptation process if learned. It is crucial to identify and learn more from those “hard” samples without sacrificing the generalizability of the model.

Algorithm 4 Dynamic Multi-epoch Update Strategy

**Input:** threshold $0 \leq \xi_1 \leq \xi_2$, optimizer $A_P$ (e.g. MEKFEMA), model prediction function $f_1$

**Input:** previous parameter $\hat{\theta}_{t-1}$, previous input $X_{t-1}$, previous prediction $\hat{y}_t = f_1(\hat{\theta}_{t-1}, X_{t-1})$, current observation $y_t$

**Output:** Adapted parameter $\hat{\theta}_t$

1. $j_t = \|y_t - \hat{y}_t\|_2$
   - $1$, if $j_t < \xi_1$
   - $2$, if $\xi_1 \leq j_t < \xi_2$
   - $0$, if $j_t \geq \xi_2$
2. $\kappa_t = 1$
3. $\hat{\theta}_t, 0 = \hat{\theta}_{t-1}$
4. for $i = 1, \ldots, \kappa_t$ do
5.   $\hat{\theta}_t, i = f_1(\theta_t, i-1, X_{t-1})$
6.   $\hat{\theta}_t, i = \hat{\theta}_t, i-1 + A_P(\theta_t, i-1, \hat{\theta}_t, i, y_t)$
7. end for
8. $\hat{\theta}_t = \hat{\theta}_t, \kappa_t$

The thresholds $\xi_1$ and $\xi_2$ can be determined by the validation set empirically. If the dataset is noise-free, there is no need to identify “anomaly” samples and we set $\xi_2 \to +\infty$. In general, we recommend the following method to find the desired $\xi_1$ and $\xi_2$. First, we need to run the single-epoch adaptation on the validation set and record each sample’s prediction error $\{j_1, j_2, \cdots\}$. Second, we set $\xi_1$ as the $50\% \sim 95\%$ quantile value of the errors, and set $\xi_2$ as the $99.9\% \sim 100\%$ quantile value.

1. The criterion can be application-specific. This criterion is proposed since it is observed in our experiment that a large number of epochs will lead to overfitting to the historical data, which is harmful to generalization. We will investigate more reasonable and effective criterion for dynamic multi-epoch update in the future.
of the errors. That means, we regard $50\% \sim 95\%$ of the samples as “easy” samples, $5\% \sim 50\%$ of the samples as “hard” samples, and $0\% \sim 0.1\%$ of the samples as “anomaly” samples.

We use $\text{MEKF}_{\text{EMA-DME}}$ to denote $\text{MEKF}_{\text{EMA}}$ with the dynamic multi-epoch update strategy.

4. Experiments

Experimental design In the experiments, we consider multi-task prediction problems for simultaneous intention and trajectory prediction of either humans or vehicles. We construct Recurrent Neural Network Salehinejad et al. (2017) (RNN) based architectures to conduct experiments on Mocap dataset (human) and NGSIM dataset (vehicle) Colyar and Halkias (2007). We leave the details of the experiment in the Appendix A.1. Before online adaptation, the prediction models are trained offline. In the following discussion, we studied the performance of various adaptation algorithms on these offline-trained models (with online adaptation on the test set). In particular, we evaluate the accuracy (0-1) for intention prediction, and the mean squared error (MSE) for trajectory prediction.

Comparison among different optimizers The proposed algorithm $\text{MEKF}_{\text{EMA-DME}}$ is compared with the based algorithm $\text{MEKF}_\lambda$ and other commonly used optimizers such as SGD, Adam, and Amsgrad. For fair comparison, we apply the dynamic multi-epoch update strategy on SGD, Adam, and Amsgrad. Here is a table showing the comparison of different optimizers.

| Dataset | Metrics       | w/o adapt | SGD         | Adam        | Amsgrad     | MEKF_\lambda | MEKF_{EMA-DME} |
|---------|---------------|-----------|-------------|-------------|-------------|---------------|----------------|
| Mocap   | accuracy      | 0.984     | 0.984       | 0.984       | 0.984       | 0.985         | 0.985          |
|         | MSE(dm^2)     | 3.271     | 3.185       | 3.149       | 3.156       | 2.788         | 2.746          |
| NGSIM   | accuracy      | 0.951     | 0.954 (.0427) | 0.951 (.0426) | 0.951 (.0427) | 0.956 (.0430) | 0.956 (.0430) |
|         | MSE(m^2)      | 2.559     | 2.367 (1.981) | 2.402 (1.975) | 2.407 (1.993) | 2.157 (1.902) | 2.092 (1.893) |

Table 1 shows the prediction performance of online adapted models using different optimizers on the Mocap dataset and the NGSIM dataset. Compared to the stochastic gradient-based algorithms, the EKF-based methods $\text{MEKF}_\lambda$ and $\text{MEKF}_{\text{EMA-DME}}$ perform better. In addition, $\text{MEKF}_{\text{EMA-DME}}$ has the best performance among all, due to the extensions inspired by EMA and dynamic multi-epoch update. On the CMU Mocap dataset, Adam reduces the trajectory MSE by 3.73%. $\text{MEKF}_\lambda$ reduces the trajectory MSE by 14.77%. $\text{MEKF}_{\text{EMA-DME}}$ reduces the trajectory MSE by 16.05%. These improvements are important to ensure safe and efficient operation of behavior prediction Zhao et al. (2020). The variance of performance using different optimizers on the NGSIM dataset is shown in the parenthesis. In particular, the standard deviation (Std.) of the prediction accuracy as well the Std. of MSE are shown. For intention prediction, $\text{MEKF}_\lambda$ and $\text{MEKF}_{\text{EMA-DME}}$ have slightly higher Std. than other SGD based optimizers. However, for trajectory prediction, $\text{MEKF}_\lambda$ and $\text{MEKF}_{\text{EMA-DME}}$ have lower Std. than other SGD based optimizers.

The running time of the adaptation algorithm correlates with the number of parameters to adapt. For the time complexity analysis, SGD, Adam, and Amsgrad have similar complexity, while $\text{MEKF}_\lambda$ and $\text{MEKF}_{\text{EMA}}$ have similar complexity. So we only compare $\text{MEKF}_\lambda$ with SGD below. For adapting encoders hidden layers (12480 parameters), SGD takes 0.08 seconds per sample and $\text{MEKF}_\lambda$ takes 0.41 seconds per sample$^2$. For real-time adaptable prediction, as long as the number of parameters to adapt is not too big, $\text{MEKF}_\lambda$ can meet the real-time requirements.

$^2$ The running time is evaluated on the NGSIM dataset using a Ubuntu desktop with Intel Core i9-9940X CPU (3.30GHz) and GeForce RTX 2080 Ti GPU.
Effectiveness of extensions  This section studies the effectiveness of the proposed extensions to MEKF\(\lambda\) in sections 3.2 and 3.3. The hyperparameters are set as \(\mu_p = 0.3\) and \(\mu_v = 0.3\). The results are shown in table 2.

1. EMA-V or momentum barely improves the performance. Two potential reasons are: 1) The momentum does not help EKF-based optimizers. In every optimization step, EKF-based optimizers has already incorporated the historical data. Hence its step size \(V_t\) is already closer to optimum than that of SGD. The learning gain in SGD is not based on historical data but manually defined. 2) The moving average on the parameter or the step size is more applicable to offline training than to online adaptation. The inapplicability is due to the fact that online adaptation can only process data sequentially in time, which is significantly different from the shuffled, repetitive, and batched process in offline training.

2. EMA-P slightly improves the performance of MEKF\(\lambda\). Filtering of \(P_t\) can smooth the inner state and improve convergence.

3. Dynamic multi-epoch update improves the performance of MEKF\(\lambda\), and it has the best performance among all the proposed extensions.

We design the additional experiment about different criteria for DME in Appendix A.2 to compared the proposed criteria (Under the error spectrum, the first 50% are easy samples, the middle 50% to 99.9% are hard samples, and the last 0.1% are anomaly samples.) with fixed 2-epoch criteria (Each sample was used twice.) and random criteria (Which has same easy and hard ratio as the proposed criterion, but distinguishing easy, hard and anomaly samples randomly.). The results in Appendix A.2 show that: the proposed criterion outperforms other criteria, which justifies the effectiveness of the proposed error-based criterion. Nonetheless, we will investigate more reasonable and effective criterion for dynamic multi-epoch update in the future.

| Dataset | Metrics | MEKF\(\lambda\) | MEKF\(\lambda\) + EMA-V | MEKF\(\lambda\) + EMA-P | MEKF\(\lambda\) + DME |
|---------|---------|----------------|--------------------------|--------------------------|------------------------|
| Mocap   | accuracy| 0.985          | 0.985                    | 0.985                    | 0.985                  |
|         | MSE(dm²) | 2.788          | 2.790                    | 2.775                    | 2.749                  |
| NGSIM   | accuracy| 0.956          | 0.956                    | 0.956                    | 0.956                  |
|         | MSE(m²)  | 2.157          | 2.156                    | 2.123                    | 2.122                  |

5. Conclusions
This paper studied online adaptation of neural network-based prediction models for behavior prediction. An EKF-based adaptation algorithm MEKF\(\lambda\) was introduced as an effective base algorithm for online adaptation. In order to improve the performance and convergence of MEKF\(\lambda\), exponential moving average filtering was investigated, including momentum and EMA-P. Then this paper introduced a dynamic multi-epoch update strategy, which is compatible with any optimizer. By combining all extensions with the base MEKF\(\lambda\) algorithm, we introduced the robust online adaptation algorithm MEKF\(\text{EMA-DME}\). In the experiments, we demonstrated the effectiveness of the proposed adaptation algorithms.

In the future, mathematical analysis of the proposed online adaptation algorithm MEKF\(\text{EMA-DME}\) will be performed in order to provide theoretical guarantees on stability, convergence, and boundedness. In addition, we will apply the proposed algorithm on a wider range of problems in addition to behavior prediction problems.
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Appendix A. Detailed Experiments

A.1. Experimental design

A.1.1. Multi-task prediction

In experiment we consider a multi-task prediction problem for simultaneous intention and trajectory prediction. Intentions are discrete representations of future trajectories. For example, in vehicle behavior prediction, intention can be acceleration and deceleration in a certain time window in the future.

The transition models for trajectory and intention prediction of the target agent are formulated as

\[ Y_t = f(\theta, X_t) \]  \hspace{1cm} (7)
\[ z_t = g(\theta, X_t) \]  \hspace{1cm} (8)

where the input vector \( X_t \) denotes the stack of \( n \)-step current and past measurement at time steps \( t, t-1, \ldots, t-n+1 \). The measurement \( x_t \) can include the position and velocity of the target agent as well as the state of the environment. For human behavior prediction, this paper uses raw position and velocity measurements of the human. For vehicle behavior prediction, this paper additionally uses environment features. The output vector \( Y_t \) denotes the stack of the \( m \)-step future trajectory at time steps \( t+1, t+2, \ldots, t+m \). Another output vector \( z_t \) is a probability distribution over different intentions at time step \( t \). The function \( f \) maps current and past measurements to the future trajectory, while the function \( g \) maps current and past measurements to the current intention.

One possible design of the multi-task prediction model is to use an encoder-decoder-classifier architecture. The encoder serves as a common part for all sub-tasks, which maps the input vector \( X_t \) to a hidden representation \( h_t \). The decoder works for trajectory prediction, which maps the hidden representation \( h_t \) to the predicted future trajectory \( Y_t \). The classifier aims to predict the intention \( z_t \) from the hidden representation \( h_t \). Mathematically, the relationships among the encoder, the decoder, and the classifier are:

\[ h_t = Encoder(\theta^E, X_t), \]  \hspace{1cm} (9)
\[ Y_t = Decoder(\theta^D, h_t), \]  \hspace{1cm} (10)
\[ z_t = Classifier(\theta^C, h_t), \]  \hspace{1cm} (11)
where $\theta^E$ is the parameter for the encoder, which affects all sub-tasks, $\theta^D$ is the parameter for the decoder, and $\theta^C$ is the parameter for the classifier. The total (ground truth) parameter of the model is $\theta := \{\theta^E, \theta^D, \theta^C\}$.

In online adaptation of multi-task learning, the adaptation algorithm updates the prediction model only considering the error measured between the predicted trajectory and the ground truth trajectory. The ground truth intention is not available for online adaptation since it is not directly observable. Figure 1 illustrates an online adaptation theme which only adapts the encoder’s parameter.

![Figure 1: The online adaptation framework for a multi-task model.](image1)

![Figure 2: The neural network architecture for the RNN-based multi-task prediction model.](image2)
A.1.2. NEURAL NETWORK ARCHITECTURE

We construct Recurrent Neural Network Salehinejad et al. (2017) (RNN) based architectures in our experiments to evaluate the effectiveness of MEKF_λ and MEKF_{EMA-DME}, as shown in Fig. 2. The neural networks follow the encoder-decoder-classifier structure for simultaneous intention and trajectory prediction as shown in Fig. 1. Trajectory prediction is based on encoder-decoder Sutskever et al. (2014) structure and intention prediction is based on encoder-classifier structure. Both the encoder and the decoder are composed of single layer Gated Recurrent Units (GRU’s) Cho et al. (2014) and the classifier is composed of two-layer FC neural networks. In order to improve the performance of offline trained model, an attention mechanism Bahdanau et al. (2014) is applied to the output vectors of the encoder.

A.1.3. DATASET

We used Mocap dataset and NGSIM dataset in our experiment. In each dataset, we randomly split the dataset as 80% offline training, 10% offline validation and 10% testing according to different trials.

1. Mocap dataset. This is a human-motion capture dataset collected by researchers from CMU^3. We chose the wrist trajectories of three actions (walking, running, and jumping) of all subjects in the Mocap datasets^4. The intentions are identified with the labeled actions. There are 543 trials for all three actions.

2. US 101 human driving data from Next Generation SIMulation (NGSIM) dataset. It is a widely used benchmark dataset for autonomous driving Colyar and Halkias (2007). We extract three actions from the dataset, which are driving with constant speed, acceleration, and deceleration respectively^5. At time step t, if the vehicle will accelerate (or decelerate) in the next three seconds(t ∼ t + 3s), we label the intention as acceleration (or deceleration) at time step t. Otherwise, we label it as constant speed. In our experiment, we used a subset of the dataset which contains 100 trials for all three actions.

A.1.4. EVALUATION METRICS

We used accuracy to evaluate the intention prediction and average mean square error (MSE) for the trajectory prediction. The average MSE is computed as,

\[
\text{MSE} = \frac{1}{T} \sum_{t=1}^{T} \text{MSE}_t, \quad \text{MSE}_t = \frac{1}{m} \| \mathbf{Y}_t - \hat{\mathbf{Y}}_t \|_2, \quad (12)
\]

Where T is total number of timesteps in the testing set. To maintain similar orders of magnitude on different datasets, we used dm^2 unit for the trajectory in Mocap dataset, and used m^2 unit for the trajectory in NGSIM dataset.

3. http://mocap.cs.cmu.edu/motcat.php
4. We didn’t perform full body motion prediction, since it requires special design of the neural network model to encode the geometric constraints, which is out of the scope of this paper.
5. In our experiment, value of acceleration \( a > 0.5 \text{ m/s}^2 \) was denoted as acceleration, and value of acceleration \( a < -0.5 \text{ m/s}^2 \) was denoted as deceleration.
A.1.5. Offline Training

Before online adaptation, the prediction models are trained offline. We used an Adam optimizer with a 128 batch size and a 0.01 learning rate. For the Mocap dataset, past $n = 20$ steps input information was used to predict the trajectories of the future $m = 10$ steps and the intention. We used a concatenation of raw trajectory and speed as the input information. For the NGSIM dataset, past $n = 20$ steps input information was used to predict the trajectories of the future $m = 50$ steps and the intention. We used a concatenation of raw trajectories and extracted features as input information. The extracted features were similar to the features used in the parameter sharing generative adversarial imitation learning Bhattacharyya et al. (2018). Table 3 shows the prediction performance after offline learning. In experiments of the adaptation, we studied the performance of various adaptation algorithms on hidden weights of encoder of these offline-trained models (with online adaptation on testing set).

| Metrics | CMU Mocap dataset | NGSIM dataset |
|---------|-------------------|---------------|
| accuracy | 0.984             | 0.951         |
| MSE     | 3.271 (dm$^2$)    | 2.559 (m$^2$) |

A.2. Additional experiment

In order to demonstrate the effectiveness of the proposed discrimination criterion in dynamic multi.epoch update strategy, we design the following experiment on the NGSIM dataset. We compared three different criteria for DME.

1. the proposed criterion as discussed in section 3.3. In particular, we set $\xi_1$ as the 50% quantile value of the errors, and $\xi_2$ as the 99.9% quantile value of the errors. Under the error spectrum, the first 50% are “easy” samples, the middle 50% to 99.9% are “hard” samples, and the last 0.1% are “anomaly” samples.
2. fixed criterion: we set $\kappa_t = 2$ for all samples. That means, we run fixed 2-epoch update strategy and use each sample twice.
3. random criterion: for each sample, we set $\kappa_t = 1$ with the probability of 50%, set $\kappa_t = 2$ with the probability of 49.9%, and set $\kappa_t = 0$ with the probability of 0.1%. That means, the random criterion has same “easy” and “hard” ratio as the proposed criterion, but distinguishing “easy”, “hard” and “anomaly” samples randomly.

|                     | w/o DME | fixed criterion | random criterion | proposed criterion |
|---------------------|---------|-----------------|------------------|--------------------|
| accuracy            | 0.956   | 0.957           | 0.957            | 0.958              |
| MSE (m$^2$)         | 2.157   | 2.136           | 2.140            | 2.122              |

The results in table 4 show that: the proposed criterion outperforms other criteria, which justifies the effectiveness of the proposed error-based criterion. Nonetheless, we will investigate more reasonable and effective criterion for dynamic multi-epoch update in the future.