# One-Step Condensed Forms for Square-Root Maximum Correntropy Criterion Kalman Filtering 

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#### Abstract

This paper suggests a few novel Cholesky-based square-root algorithms for the maximum correntropy criterion Kalman filtering. In contrast to the previously obtained results, new algorithms are developed in the so-called condensed form that corresponds to the a priori filtering. Square-root filter implementations are known to possess a better conditioning and improved numerical robustness when solving ill-conditioned estimation problems. Additionally, the new algorithms permit easier propagation of the state estimate and do not require a back-substitution for computing the estimate. Performance of novel filtering methods is examined by using a fourth order benchmark navigation system example.


Keywords-maximum correntropy, Kalman filter, Cholesky decomposition, one-step filtering, condensed form

## I. INTRODUCTION

The Kalman filtering (KF) like estimators developed under the maximum correntropy criterion (MCC) methodology are shown to outperform the classical KF for estimation quality in case of non-Gaussian and impulsive noise scenario in state' space models, both for linear and nonlinear systems [1]-[10]. Among nonlinear MCC filtering strategies, we may mention the MCC extended KF in [11], [12], the MCC unscented KF in [13], [14], the MCC-based Gauss-Hermite quadrature filter in [15] and the so-called accurate continuous-discrete extended KF [16], [17] explored under the MCC in [18]. Here, we focus on the MCC-KF estimators suggested in [6], [9].

The numerical stability issues of the cited MCC-KF estimators are of special interest in this paper. Recently, their reliable square-root implementation methods have been designed in [19]. Following the approaches existed in the KF community, the new MCC-KF implementations belong to three main factored-form classes: (i) the Cholesky-based algorithms; e.g., [20]-[22], (ii) the $U D U^{\top}$ factorization-based methods, e.g., [23], [24], and (iii) singular value decomposition solution, e.g., [25], [26]. All mentioned square-root strategies have been utilized for the robust MCC-KF filtering in [19]. In the cited paper, the implementation methods are developed in terms of covariance quantities and in the a posteriori form that yields the two-stage implementations. In contrast to the previously

[^0]obtained results, the goal of this paper is to suggest squareroot MCC-KF algorithms in the a priori form that results to the condensed implementations. We stress that the stated problem has never been solved in engineering literature, before.

More precisely, we develop the a priori covariance filtering approach for the original MCC-KF proposed in [6] and for its improved variant (IMCC-KF) published in [9]. Additionally, we discuss the square-root problem for the examined estimators and suggest a few condensed implementations for the robust filtering in terms of lower triangular Cholesky factors. As all square-root implementations, the new algorithms possess a better conditioning, improved numerical robustness (with respect to roundoff errors) and reliability when solving ill-conditioned estimation problems. Additionally, the new algorithms permit easier propagation of the state estimate and do not require a back-substitution for computing the estimate. It is achieved by utilizing orthogonal rotations for propagating the involved Cholesky factors as far as possible. Finally, we stress that derivation of new implementation methods is important for its own sake because it provides practitioners a diversity of methods with a fair possibility to choose any of them depending on practical application, complexity and accuracy requirements. Performance of novel filtering algorithms is examined by using a fourth order benchmark navigation system example.
The paper is organized as follows. Section 【 contains the problem statement and presents the derivation of the a priori MCC-KF recursions. The new square-root condensed implementation methods can be found in Section III together with their derivation. Section IV contains the results of numerical simulation study and Section $\square$ concludes the paper.

## II. PROBLEM STATEMENT

Consider a linear discrete-time stochastic system

$$
\begin{align*}
x_{k+1} & =F_{k} x_{k}+G_{k} w_{k},  \tag{1}\\
y_{k} & =H_{k} x_{k}+v_{k}, \quad k \geq 0 \tag{2}
\end{align*}
$$

where system matrices $F_{k} \in \mathbb{R}^{n \times n}, G_{k} \in \mathbb{R}^{n \times q}$ and $H_{k} \in \mathbb{R}^{m \times n}$ are known and constant over time. The vectors $x_{k} \in \mathbb{R}^{n}$ and $y_{k} \in \mathbb{R}^{m}$ are unknown dynamic state and
available measurement vector, respectively. We assume that random variables $x_{0}, w_{k}$ and $v_{k}$ satisfy

$$
\begin{array}{ll}
\mathbf{E}\left\{x_{0}\right\}=\bar{x}_{0}, & \mathbf{E}\left\{\left(x_{0}-\bar{x}_{0}\right)\left(x_{0}-\bar{x}_{0}\right)^{\top}\right\}=\Pi_{0}, \\
\mathbf{E}\left\{w_{k}\right\}=\mathbf{E}\left\{v_{k}\right\}=0, & \mathbf{E}\left\{w_{k} x_{0}^{\top}\right\}=\mathbf{E}\left\{v_{k} x_{0}^{\top}\right\}=0, \\
\mathbf{E}\left\{w_{k} v_{k}^{\top}\right\}=0, & \mathbf{E}\left\{w_{k} w_{j}^{\top}\right\}=Q_{k} \delta_{k j}, \\
& \mathbf{E}\left\{v_{k} v_{j}^{\top}\right\}=R_{k} \delta_{k j}
\end{array}
$$

where covariance matrices $Q_{k} \in \mathbb{R}^{q \times q}$ and $R_{k} \in \mathbb{R}^{m \times m}$ are known. The symbol $\delta_{k j}$ is the Kronecker delta function.

The classical KF produces the minimum linear expected mean square error (MSE) estimate of the unknown state vector $x_{k}$. It can be formulated in the a priori form as shown in [27, Theorem 9.2.1]; see also the derivation in [28, Section 5.3]. More precisely, the one-step ahead predicted estimate $\hat{x}_{k+1 \mid k}$ (a priori estimate) is calculated by

$$
\begin{align*}
\hat{x}_{k+1 \mid k} & =F_{k} \hat{x}_{k \mid k-1}+K_{p, k}\left(y_{k}-H_{k} \hat{x}_{k \mid k-1}\right),  \tag{3}\\
R_{e, k} & =R_{k}+H_{k} P_{k \mid k-1} H_{k}^{\top},  \tag{4}\\
K_{p, k} & =F_{k} P_{k \mid k-1} H_{k}^{\top} R_{e, k}^{-1} \tag{5}
\end{align*}
$$

where $e_{k}=y_{k}-H_{k} \hat{x}_{k \mid k-1}$ are innovations (residuals) of discrete-time KF with covariance $R_{e, k}=\mathbf{E}\left\{e_{k} e_{k}^{\top}\right\}$ and $K_{p, k}=\mathbf{E}\left\{\hat{x}_{k+1 \mid k} e_{k}^{\top}\right\}=F_{k} P_{k \mid k-1} H_{k}^{\top}$. The matrix $P_{k \mid k-1}$ is one-step ahead predicted error covariance $P_{k \mid k-1}=$ $\mathbf{E}\left\{\left(x_{k}-\hat{x}_{k \mid k-1}\right)\left(x_{k}-\hat{x}_{k \mid k-1}\right)^{\top}\right\}$ computed via the Riccati difference recursion:

$$
\begin{equation*}
P_{k+1 \mid k}=F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-K_{p, k} R_{e, k} K_{p, k}^{\top} \tag{6}
\end{equation*}
$$

with initial values $P_{0 \mid-1}=\Pi_{0}>0$ and $\hat{x}_{0 \mid-1}=\bar{x}_{0}$.
Being a linear estimator, the classical KF exhibits only sub-optimal behavior in non-Gaussian settings. To enhance its estimation quality and robustness with respect to outliers (impulsive noise), the maximum correntropy criterion has been utilized together with the KF estimation cost function in [5], [6]. In general, the correntropy is a similarity measure between two random variables [1]. Following [29, Chapter 5], it can be used in the estimation problem as follows: an estimator of unknown state $X \in \mathbb{R}$ is defined as a function of measurements $Y \in \mathbb{R}^{m}$, i.e., $\hat{X}=g(Y)$ where $g$ is solved by maximizing the correntropy between $X$ and $\hat{X}$ that is [30]

$$
\begin{equation*}
\arg \max _{g \in G} V(X, \hat{X})=\arg \max _{g \in G} \mathbf{E}\left\{k_{\sigma}(X-g(Y))\right\} \tag{7}
\end{equation*}
$$

where $G$ stands for the collection of all measurable functions of $Y, k_{\sigma}(\cdot)$ is a kernel function and $\sigma>0$ is the kernel size. The Gaussian kernel is the most widely used kernel function and it is given as follows:

$$
\begin{equation*}
k_{\sigma}(X-\hat{X})=\exp \left\{-(X-\hat{X})^{2} /\left(2 \sigma^{2}\right)\right\} . \tag{8}
\end{equation*}
$$

It is not difficult to see that the MCC cost (7) with Gaussian kernel (8) reaches its maximum if and only if $X=\hat{X}$.
In [6], the MCC-KF is developed by solving the following estimation problem with the Gaussian kernel:

$$
\begin{aligned}
\arg \max & \left\{k_{\sigma}\left(\left\|\hat{x}_{k \mid k}-F_{k-1} \hat{x}_{k-1 \mid k-1}\right\|\right)\right. \\
& \left.+k_{\sigma}\left(\left\|y_{k}-H_{k} \hat{x}_{k \mid k}\right\|\right)\right\}
\end{aligned}
$$

where weighted norms are additionally utilized.
Next, the resulted nonlinear equation arisen in the optimization problem above is solved by a fixed point rule that yields the following recursion [6, p. 503]:

$$
\begin{equation*}
\hat{x}_{k \mid k}=\hat{x}_{k \mid k-1}+K_{k}^{\lambda}\left(y_{k}-H_{k} \hat{x}_{k \mid k-1}\right) \tag{9}
\end{equation*}
$$

where $\hat{x}_{k \mid k-1}:=F_{k-1} \hat{x}_{k-1 \mid k-1}$ and the gain matrix is defined as $K_{k}^{\lambda}=\lambda_{k}\left(P_{k \mid k-1}^{-1}+\lambda_{k} H_{k}^{\top} R_{k}^{-1} H_{k}\right)^{-1} H_{k}^{\top} R_{k}^{-1}$ with a scalar adjusting weight $\lambda_{k}$ given as follows:

$$
\begin{equation*}
\lambda_{k}=\frac{k_{\sigma}\left(\left\|y_{k}-H_{k} \hat{x}_{k \mid k-1}\right\|_{R_{k}^{-1}}\right)}{k_{\sigma}\left(\left\|\hat{x}_{k \mid k-1}-F_{k-1} \hat{x}_{k-1 \mid k-1}\right\|_{P_{k \mid k-1}^{-1}}\right)} . \tag{10}
\end{equation*}
$$

Finally, the resulted MCC-KF estimator in [6] utilizes the recursion for state estimate in (9) together with a symmetric Joseph stabilized equation for calculating the error covariance matrix $P_{k \mid k}$ that is known to improve the numerical robustness of KF-like implementations because of ensured symmetric form of $P_{k \mid k}$ in the presence of roundoff errors [28], [31]. In summary, the MCC-KF estimator is given as follows [6].
MCC-KF estimator. Time Update $(k=\overline{1, K})$.

$$
\begin{align*}
& \hat{x}_{k \mid k-1}=F_{k-1} \hat{x}_{k-1 \mid k-1},  \tag{11}\\
& P_{k \mid k-1}=F_{k-1} P_{k-1 \mid k-1} F_{k-1}^{T}+G_{k-1} Q_{k-1} G_{k-1}^{\top} . \tag{12}
\end{align*}
$$

Measurement Update. Find a posteriori estimate:

$$
\begin{align*}
K_{k}^{\lambda} & =\lambda_{k}\left(P_{k \mid k-1}^{-1}+\lambda_{k} H_{k}^{\top} R_{k}^{-1} H_{k}\right)^{-1} H_{k}^{\top} R_{k}^{-1}  \tag{13}\\
\hat{x}_{k \mid k} & =\hat{x}_{k \mid k-1}+K_{k}^{\lambda} e_{k}, \quad e_{k}=y_{k}-H_{k} \hat{x}_{k \mid k-1}  \tag{14}\\
P_{k \mid k} & =\left(I-K_{k}^{\lambda} H_{k}\right) P_{k \mid k-1}\left(I-K_{k}^{\lambda} H_{k}\right)^{\top} \\
& +K_{k}^{\lambda} R_{k}\left[K_{k}^{\lambda}\right]^{\top} \tag{15}
\end{align*}
$$

where the scaling (inflation) parameter $\lambda_{k}$ is computed by (10).
In [9], the estimation quality of original MCC-KF estimator presented above has been boosted by deriving mathematically equivalent formulas for the gain and error covariance, similar to the KF equations in [28, pp. 128-129]. It yields the improved MCC-KF (IMCC-KF) estimator summarized below.
IMCC-KF estimator. Time Update $(k=\overline{1, K})$. Calculate a priori estimate $\hat{x}_{k \mid k-1}$ and $P_{k \mid k-1}$ by equations (11), (12).

Measurement Update. Find a posteriori estimate $\hat{x}_{k \mid k}$ by formula (14) with the scaling parameter $\lambda_{k}$ from (10) and

$$
\begin{align*}
K_{k}^{\lambda} & =\lambda_{k} P_{k \mid k-1} H_{k}^{\top}\left[R_{e, k}^{\lambda}\right]^{-1}  \tag{16}\\
R_{e, k}^{\lambda} & =\lambda_{k} H_{k} P_{k \mid k-1} H_{k}^{\top}+R_{k},  \tag{17}\\
P_{k \mid k} & =\left(I-K_{k}^{\lambda} H_{k}\right) P_{k \mid k-1} . \tag{18}
\end{align*}
$$

As can be seen, the original MCC-KF and IMCC-KF equations recursively updates a posteriori estimate $\hat{x}_{k \mid k}$ given the initial value $\hat{x}_{0 \mid 0}:=\bar{x}_{0}$. The discussed estimators can be re-formulated in the so-called a priori form similar to the classical KF equations (3) - (6). This yields one-step KFlike equations and allows for convenient condensed form of the resulted square-root implementation methods. Next, we derive one-step equations for both the MCC-KF and IMCCKF estimators and explore the possibility of designing their Cholesky factorization-based implementations.

## A. The MCC-KF estimator in a priori form

Following [28, Section 5.3], we derive a priori MCC-KF recursion related to the original $\mathrm{MCC}-\mathrm{KF}$ equations (11)-15). Having substituted (14) into (11), we get the recursion for onestep ahead predicted (a priori) estimate $\hat{x}_{k+1 \mid k}$ as follows:

$$
\begin{align*}
\hat{x}_{k+1 \mid k} & =F_{k} \hat{x}_{k \mid k}=F_{k}\left(\hat{x}_{k \mid k-1}+K_{k}^{\lambda} e_{k}\right), \\
& =F_{k} \hat{x}_{k \mid k-1}+F_{k} K_{k}^{\lambda} e_{k}=F_{k} \hat{x}_{k \mid k-1}+K_{p, k}^{\lambda} e_{k} \tag{19}
\end{align*}
$$

where we introduced the notation $K_{p, k}^{\lambda}=F_{k} K_{k}^{\lambda}$.
It is proved in [9] that equations (13) and (16) for gain calculation $K_{k}^{\lambda}$ are equivalent. Hence, we have

$$
\begin{align*}
K_{p, k}^{\lambda} & =\lambda_{k} F_{k}\left(P_{k \mid k-1}^{-1}+\lambda_{k} H_{k}^{\top} R_{k}^{-1} H_{k}\right)^{-1} H_{k}^{\top} R_{k}^{-1}  \tag{20}\\
& =\lambda_{k} F_{k} P_{k \mid k-1} H_{k}^{\top}\left[R_{e, k}^{\lambda}\right]^{-1} \tag{21}
\end{align*}
$$

Having substituted formula (15) into (12) and taking into account formula (21) and $K_{p, k}^{\lambda}=F_{k} K_{k}^{\lambda}$, we obtain

$$
\begin{aligned}
P_{k+1 \mid k} & =F_{k} P_{k \mid k} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top} \\
& =F_{k}\left(I-K_{k}^{\lambda} H_{k}\right) P_{k \mid k-1}\left(I-K_{k}^{\lambda} H_{k}\right)^{\top} F_{k}^{\top} \\
& +F_{k} K_{k}^{\lambda} R_{k}\left[K_{k}^{\lambda}\right]^{\top} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top} \\
& =\left(F_{k}-K_{p, k}^{\lambda} H_{k}\right) P_{k \mid k-1}\left(F_{k}-K_{p, k}^{\lambda} H_{k}\right)^{\top} \\
& +K_{p, k}^{\lambda} R_{k}\left[K_{p, k}^{\lambda}\right]^{\top}+G_{k} Q_{k} G_{k}^{\top} \\
& =F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-F_{k} P_{k \mid k-1} H_{k}^{\top}\left[K_{p, k}^{\lambda}\right]^{\top} \\
& -K_{p, k}^{\lambda} H_{k} P_{k \mid k-1} F^{\top} \\
& +K_{p, k}^{\lambda}\left(H_{k} P_{k \mid k-1} H_{k}^{\top}+R_{k}\right)\left[K_{p, k}^{\lambda}\right]^{\top} .
\end{aligned}
$$

Taking into account that $H_{k} P_{k \mid k-1} H_{k}^{\top}+R_{k}=R_{e, k}$, formula (21) and symmetric form of $R_{e, k}$ and $R_{e, k}^{\lambda}$, we get

$$
F_{k} P_{k \mid k-1} H_{k}^{\top}=\lambda_{k}^{-1} K_{p, k}^{\lambda} R_{e, k}^{\lambda}
$$

we obtain the recursion for a priori error covariance $P_{k+1 \mid k}$ :

$$
\begin{aligned}
P_{k+1 \mid k} & =F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-2 \lambda_{k}^{-1} K_{p, k}^{\lambda} R_{e, k}^{\lambda}\left[K_{p, k}^{\lambda}\right]^{\top} \\
& +K_{p, k}^{\lambda} R_{e, k}\left[K_{p, k}^{\lambda}\right]^{\top}=F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top} \\
& -K_{p, k}^{\lambda}\left[2 \lambda_{k}^{-1} R_{e, k}^{\lambda}-R_{e, k}\right]\left[K_{p, k}^{\lambda}\right]^{\top} \\
& =F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top} \\
& -K_{p, k}^{\lambda}\left(H_{k} P_{k \mid k-1} H_{k}^{\top}+\left(2 \lambda_{k}^{-1}-1\right) R_{k}\right)\left[K_{p, k}^{\lambda}\right]^{\top} .
\end{aligned}
$$

As can be seen, the MCC-KF equation for matrix $P_{k+1 \mid k}$ computation does not allow the related square-root form because the term $2 \lambda_{k}^{-1}-1$ might be negative, although the adjusting weight $\lambda_{k}$ defined in (10) is a nonnegative value. Meanwhile, for the improved MCC-KF counterpart the efficient square-root implementations do exist as shown below.

## B. The IMCC-KF estimator in a priori form

Similarly, a priori recursion for the IMCC-KF equations (11), (12), (14), (16)-(18) has been recently derived in [32]. For readers' convenience, we briefly discuss it here. First, one-step ahead predicted estimate $\hat{x}_{k+1 \mid k}$ is calculated by equation (19) as in the MCC-KF examined above.

Next, having substituted formula (18) into (12) and taking into account formula (21) and symmetric form of any covariance matrix, we obtain

$$
\begin{aligned}
P_{k+1 \mid k} & =F_{k} P_{k \mid k} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top} \\
& =F_{k}\left[\left(I-K_{k}^{\lambda} H_{k}\right) P_{k \mid k-1}\right] F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top} \\
& =F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-F_{k} K_{k}^{\lambda} H_{k} P_{k \mid k-1} F_{k}^{\top} \\
& =F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-K_{p, k}^{\lambda}\left(F_{k} P_{k \mid k-1} H_{k}^{\top}\right)^{\top} \\
& =F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-\frac{1}{\lambda_{k}} K_{p, k}^{\lambda} R_{e, k}^{\lambda}\left[K_{p, k}^{\lambda}\right]^{\top}
\end{aligned}
$$

or, alternatively, we can avoid a scalar division by introducing

$$
\begin{equation*}
K_{p, k}=\frac{1}{\lambda_{k}} F_{k} K_{k}^{\lambda}=F_{k} P_{k \mid k-1} H_{k}^{\top}\left[R_{e, k}^{\lambda}\right]^{-1} \tag{22}
\end{equation*}
$$

i.e., $K_{p, k}^{\lambda}=\lambda_{k} K_{p, k}$ defined in (21). Thus, the equation for error covariance computation is re-formulated as follows:

$$
\begin{align*}
P_{k+1 \mid k}= & F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-\lambda_{k} K_{p, k} R_{e, k}^{\lambda} K_{p, k}^{\top}  \tag{23}\\
\text { where } & K_{p, k}=F_{k} P_{k \mid k-1} H_{k}^{\top}\left[R_{e, k}^{\lambda}\right]^{-1}  \tag{24}\\
& \hat{x}_{k+1 \mid k}=F_{k} \hat{x}_{k \mid k-1}+\lambda_{k} K_{p, k} e_{k} \tag{25}
\end{align*}
$$

Formulas (23)-(25) are the a priori form of the IMCC-KF estimator, which we are looking for. It is important that the adjusting weight $\lambda_{k}$ defined in (10) is a nonnegative value and, hence, a square root exists (for real nonnegative numbers).

## III. MAIN RESULTS: CONDENSED FORMS

The resulted one-step IMCC-KF equations (23) - (25) allow for convenient condensed form of stable square-root implementation methods that we derive in this section. Our square-root solution is based on the Cholesky decomposition of a symmetric positive definite matrix $A$ in the form $A=A^{1 / 2} A^{\top / 2}$ where the factor $A^{1 / 2}$ is a lower triangular matrix with positive diagonal elements.

The condensed form implies that one-step IMCC-KF equations (23) - (25) are summarized into unique pre-array and, next, a stable orthogonal rotation is applied for obtaining the resulted post-array with the propagated filter quantities. Such implementation methods are additionally suitable for parallel implementation. Thus, we formulate the first condensed implementation method for the IMCC-KF estimator.

## Algorithm 1. SR IMCC-KF (a priori, one-step form)

Initialization: $(k=0)$
Apply Cholesky factorization: $\Pi_{0}=\Pi_{0}^{1 / 2} \Pi_{0}^{\top / 2}$;
2 Set initial values: $\hat{x}_{0 \mid-1}=\bar{x}_{0}, P_{0 \mid-1}^{1 / 2}=\Pi_{0}^{1 / 2}$;
Filter Recursion: $(k=\overline{0, K})$
Compute $\lambda_{k}$ and residual $e_{k}$ by formulas (10), (14);
4 Build pre-array and block-triangularize it
$\underbrace{\left[\begin{array}{ccc}R_{k}^{1 / 2} & \lambda_{k}^{1 / 2} H_{k} P_{k \mid k-1}^{1 / 2} & 0 \\ 0 & F_{k} P_{k \mid k-1}^{1 / 2} & G_{k} Q_{k}^{1 / 2}\end{array}\right]}_{\text {Pre-array } \mathbb{A}_{k}} \mathbb{Q}_{k}=\underbrace{\left[\begin{array}{ccc}{\left[R_{e, k}^{\lambda}\right]^{1 / 2}} & 0 & 0 \\ \bar{K}_{p, k}^{\lambda} & P_{k+1 \mid k}^{1 / 2} & 0\end{array}\right]}_{\text {Post-array } \mathbb{R}_{k}} ;$
5 Read-off from post-array: $P_{k+1 \mid k}^{1 / 2},\left[R_{e, k}^{\lambda}\right]^{1 / 2}, \bar{K}_{p, k}^{\lambda}$;
$6 \quad$ Find $\hat{x}_{k+1 \mid k}=F_{k} \hat{x}_{k \mid k-1}+\lambda_{k}^{1 / 2}\left[\bar{K}_{p, k}^{\lambda}\right]\left[R_{e, k}^{\lambda}\right]^{-1 / 2} e_{k}$.

Following [21], it is not difficult to prove that the transformation in line 4 of Algorithm 1 implies the calculation by formulas (23) and (24) of the conventional one-step implementation. Let's consider the pre-array in line 4 of Algorithm 1

$$
\left[\begin{array}{ccc}
R_{k}^{1 / 2} & \lambda_{k}^{1 / 2} H_{k} P_{k \mid k-1}^{1 / 2} & 0 \\
0 & F_{k} P_{k \mid k-1}^{1 / 2} & G_{k} Q_{k}^{1 / 2}
\end{array}\right]
$$

and block-triangularize it by any orthogonal rotation, i.e.,

$$
\mathbb{A}_{k} \mathbb{Q}_{k}=\left[\begin{array}{lll}
X & 0 & 0 \\
Y & Z & 0
\end{array}\right]
$$

It is worth noting here that the orthogonal transformation $\mathbb{Q}_{k}$ sets up a conformal (i.e., a norm- and angle-preserving) mapping between the (block) rows of the pre-array $\mathbb{A}_{k}$ and post-array. Thus, we have

$$
\begin{gathered}
<\left[R_{k}^{1 / 2}, \lambda_{k}^{1 / 2} H_{k} P_{k \mid k-1}^{1 / 2}, 0\right],\left[R_{k}^{1 / 2}, \lambda_{k}^{1 / 2} H_{k} P_{k \mid k-1}^{1 / 2}, 0\right]> \\
=<[X, 0,0],[X, 0,0]> \\
<\left[R_{k}^{1 / 2}, \lambda_{k}^{1 / 2} H_{k} P_{k \mid k-1}^{1 / 2}, 0\right],\left[0, F_{k} P_{k \mid k-1}^{1 / 2}, G_{k} Q_{k}^{1 / 2}\right]> \\
=<[X, 0,0],[Y, Z, 0]> \\
<\left[0, F_{k} P_{k \mid k-1}^{1 / 2}, G_{k} Q_{k}^{1 / 2}\right],\left[0, F_{k} P_{k \mid k-1}^{1 / 2}, G_{k} Q_{k}^{1 / 2}\right]> \\
=<[Y, Z, 0],[Y, Z, 0]>
\end{gathered}
$$

From the first equality above, we obtain

$$
X X^{\top}=\lambda_{k} H_{k}\left(P_{k \mid k-1}^{1 / 2} P_{k \mid k-1}^{\top / 2}\right) H_{k}^{\top}+R_{k}
$$

Having compared with equation (17), we conclude that $X X^{\top}=R_{e, k}^{\lambda}$ and, hence, $X:=\left[R_{e, k}^{\lambda}\right]^{1 / 2}$.

At the same way, from the second equality we get

$$
X Y^{\top}=\lambda_{k}^{1 / 2} H_{k}\left(P_{k \mid k-1}^{1 / 2} P_{k \mid k-1}^{\top / 2}\right) F_{k}^{\top}
$$

and, hence, we define $Y=\lambda_{k}^{1 / 2} F_{k} P_{k \mid k-1} H_{k}^{T}\left[R_{e, k}^{\lambda}\right]^{-\top / 2}:=$ $\bar{K}_{p, k}^{\lambda}$. Having compared the normalized quantity $\bar{K}_{p, k}^{\lambda}$ naturally appeared in square-root Algorithm 1 with the value $K_{p, k}$ defined by equation (24) in the conventional one-step method, we obtain the relationship

$$
\bar{K}_{p, k}^{\lambda}=\lambda_{k}^{1 / 2} K_{p, k}\left[R_{e, k}^{\lambda}\right]^{1 / 2} \text { or } K_{p, k}=\lambda_{k}^{-1 / 2} \bar{K}_{p, k}^{\lambda}\left[R_{e, k}^{\lambda}\right]^{-1 / 2}
$$

Taking into account the formula above, we express equation (25) of the conventional method in terms of values available in the square-root Algorithm 1, i.e., in terms of $\left[\bar{K}_{p, k}^{\lambda}\right]$ and $\left[R_{e, k}^{\lambda}\right]^{1 / 2}$ that are directly read-off from the post-array

$$
\begin{aligned}
\hat{x}_{k+1 \mid k} & =F_{k} \hat{x}_{k \mid k-1}+\lambda_{k} K_{p, k} e_{k} \\
& =F_{k} \hat{x}_{k \mid k-1}+\lambda_{k}^{1 / 2}\left[\bar{K}_{p, k}^{\lambda}\right]\left[R_{e, k}^{\lambda}\right]^{-1 / 2} e_{k}
\end{aligned}
$$

that is exactly the formula in line 6 of Algorithm 1.
Finally, we consider the last equality, i.e.,

$$
F_{k}\left(P_{k \mid k-1}^{1 / 2} P_{k \mid k-1}^{\top / 2}\right) F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}=Y Y^{\top}+Z Z^{\top}
$$

Having substituted $Y=\bar{K}_{p, k}^{\lambda}$ and $\bar{K}_{p, k}^{\lambda}=\lambda_{k}^{1 / 2} K_{p, k}\left[R_{e, k}^{\lambda}\right]^{1 / 2}$, we get

$$
\begin{aligned}
Z Z^{\top} & =F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-\bar{K}_{p, k}^{\lambda}\left[\bar{K}_{p, k}^{\lambda}\right]^{\top} \\
& =F_{k} P_{k \mid k-1} F_{k}^{\top}+G_{k} Q_{k} G_{k}^{\top}-\lambda_{k} K_{p, k} R_{e, k}^{\lambda} K_{p, k}^{\top}
\end{aligned}
$$

i.e., from (23) we conclude $Z:=P_{k+1 \mid k}^{1 / 2}$. Thus, the squareroot Algorithm 1 is justified.

The newly-derived square-root condensed form in Algorithm 1 is equivalent to conventional one-step formulas (23) (25), but it is convenient for practical application and more suited to parallel computations. Besides, the square-root formulation makes the new Algorithm 1 is inherently more stable (with respect to roundoff errors) than any conventional onestep implementation; e.g., the one-step formulas (23) - (25).
We further suggest one more stable square-root implementation developed in the condensed form with the following additional benefit. As can be seen, Algorithms 1 allows for simple propagation of the filter quantities $\left[\bar{K}_{p, k}^{\lambda}\right]$ and $\left[R_{e, k}^{\lambda}\right]^{1 / 2}$ at each iteration step. Indeed, these values are simply read-off from the resulted post-array obtained from block-triangularization of the pre-array. Although stable orthogonal rotations are used for $\left[\bar{K}_{p, k}^{\lambda}\right]$ and $\left[R_{e, k}^{\lambda}\right]^{1 / 2}$ calculation, the state estimate computation still requires the $m \times m$ matrix inversion that is the inverse of $\left[R_{e, k}^{\lambda}\right]^{1 / 2}$; see line 6 of Algorithm 1. We stress that $\left[R_{e, k}^{\lambda}\right]^{1 / 2}$ is a triangular matrix and, hence, the inversion can be performed in efficient way (i.e., with the reduced computational cost) by solving the related linear equations through backward substitution. However, we can avoid the matrix inversion at all by the following trick. We augment the pre-array by a specially designed "data" column and apply the same orthogonal rotation to get the augmented post-array. Such algorithms are called the extended implementations and they skip the matrix inversion for computing the state estimate.

Following [21], consider the extended pre-array

$$
\left[\begin{array}{ccc}
R_{k}^{1 / 2} & \lambda_{k}^{1 / 2} H_{k} P_{k \mid k-1}^{1 / 2} & 0 \\
0 & F_{k} P_{k \mid k-1}^{1 / 2} & G_{k} Q_{k}^{1 / 2} \\
\hline-\lambda_{k}^{1 / 2} y_{k}^{\top} R_{k}^{-\top / 2} & \hat{x}_{k \mid k-1}^{\top} P_{k \mid k-1}^{-1 / 2} & 0
\end{array}\right]
$$

and lower-triangularize the first two (block) rows

$$
\tilde{\mathbb{A}}_{k} \mathbb{Q}_{k}=\left[\begin{array}{ccc}
{\left[R_{e, k}^{\lambda}\right]^{1 / 2}} & 0 & 0 \\
\bar{K}_{p, k}^{\lambda} & P_{k+1 \mid k}^{1 / 2} & 0 \\
\hline \alpha & \beta & \gamma
\end{array}\right] .
$$

Similar to above presented derivation of Algorithm 1, we need to find $\alpha, \beta$ and $\gamma$. Taking into account properties of orthogonal matrices, from the inner product of the first and the last (block) rows of the pre-array, we get

$$
\begin{aligned}
& -\lambda_{k}^{1 / 2} R_{k}^{1 / 2} R_{k}^{-1 / 2} y_{k}+\lambda_{k}^{1 / 2} H_{k} P_{k \mid k-1}^{1 / 2} P_{k \mid k-1}^{-1 / 2} \hat{x}_{k \mid k-1} \\
& \quad=\left[R_{e, k}^{\lambda}\right]^{1 / 2} \alpha^{\top}=-\lambda_{k}^{1 / 2}\left(y_{k}-H_{k} \hat{x}_{k \mid k-1}\right)=-\lambda_{k}^{1 / 2} e_{k}
\end{aligned}
$$

i.e., $\alpha^{\top}=-\lambda_{k}^{1 / 2}\left[R_{e, k}^{\lambda}\right]^{-1 / 2} e_{k}$. We introduce notation for the normalized residuals

$$
\bar{e}_{k}=\left[R_{e, k}^{\lambda}\right]^{-1 / 2} e_{k} \text { and } \bar{e}_{k}^{\lambda}=\lambda_{k}^{1 / 2} \bar{e}_{k}=\lambda_{k}^{1 / 2}\left[R_{e, k}^{\lambda}\right]^{-1 / 2} e_{k} .
$$

Hence, we conclude that $\alpha^{\top}=-\bar{e}_{k}^{\lambda}$. Next, equation (25) of conventional one-step method is expressed in terms of available $\left[\bar{K}_{p, k}^{\lambda}\right]$ and the normalized residuals as follows:

$$
\hat{x}_{k+1 \mid k}=F_{k} \hat{x}_{k \mid k-1}+\lambda_{k}^{1 / 2}\left[\bar{K}_{p, k}^{\lambda}\right] \bar{e}_{k}=F_{k} \hat{x}_{k \mid k-1}+\left[\bar{K}_{p, k}^{\lambda}\right] \bar{e}_{k}^{\lambda}
$$

where no matrix inversion is required for calculating the state, in contrast to the formula in line 6 of Algorithm 1.

It is worth noting here that the state vector can be calculated in alternative way. For that, we consider the inner product of the second and the last (block) rows of the pre-array in order to define $\beta$ value

$$
\begin{aligned}
\bar{K}_{p, k}^{\lambda} \alpha^{\top}+P_{k+1 \mid k}^{1 / 2} \beta^{\top} & =-\bar{K}_{p, k}^{\lambda} \bar{e}_{k}^{\lambda}+P_{k+1 \mid k}^{1 / 2} \beta^{\top} \\
& =F_{k} P_{k \mid k-1}^{1 / 2} P_{k \mid k-1}^{-1 / 2} \hat{x}_{k \mid k-1}=F_{k} \hat{x}_{k \mid k-1}
\end{aligned}
$$

and, next, we conclude

$$
\beta^{\top}=P_{k+1 \mid k}^{-1 / 2}\left(F_{k} \hat{x}_{k \mid k-1}+\bar{K}_{p, k}^{\lambda} \bar{e}_{k}^{\lambda}\right)=P_{k+1 \mid k}^{-1 / 2} \hat{x}_{k+1 \mid k}
$$

Thus, an alternative state computation way is to read-off values $\left[P_{k+1 \mid k}^{-1 / 2} \hat{x}_{k+1 \mid k}\right]$ and $\left[P_{k+1 \mid k}^{1 / 2}\right]$ from the extended postarray, and simply multiply these blocks to obtain

$$
\hat{x}_{k+1 \mid k}=\left[P_{k+1 \mid k}^{1 / 2}\right]\left[P_{k+1 \mid k}^{-1 / 2} \hat{x}_{k+1 \mid k}\right] .
$$

In summary, the extended square-root condensed form implementation is presented in Algorithm 2.

```
Algorithm 2. ESR IMCC-KF (a priori, one-step, extended)
    InITIALIZATION: \((k=0)\)
        Apply Cholesky factorization: \(\Pi_{0}=\Pi_{0}^{1 / 2} \Pi_{0}^{\top / 2}\);
        Set initials: \(P_{0 \mid-1}^{1 / 2}=\Pi_{0}^{1 / 2}, P_{0 \mid-1}^{-1 / 2} \hat{x}_{0 \mid-1}=\Pi_{0}^{-1 / 2} \bar{x}_{0}\);
    Filter Recursion: \((k=\overline{0, K})\)
        Compute \(\lambda_{k}\) by equation (10);
        Build and triangularize the first two (block) rows
            \(\underbrace{\left[\begin{array}{ccc}R_{k}^{1 / 2} & \lambda_{k}^{1 / 2} H_{k} P_{k \mid k-1}^{1 / 2} & 0 \\ 0 & F_{k} P_{k k \mid k-1}^{1 / /} & G_{k} Q_{k}^{1 / 2} \\ \hline-\lambda_{k}^{1 / 2} y_{k}^{\top} R_{k}^{-\top / 2} & {\left[P_{k \mid k-1}^{-1 / 2} \hat{x}_{k \mid k-1}\right]^{\top}} & 0\end{array}\right]}_{\text {Extended pre-array } \tilde{\mathbb{A}}_{k}} \mathbb{Q}_{k}\)
        \(=\underbrace{\left[\begin{array}{ccc}{\left[R_{e, k}^{\lambda}\right]^{1 / 2}} & \text { Extended pre-array } & 0 \\ \tilde{\mathbb{R}}_{p, k}^{\lambda} & P_{k+1 \mid k}^{1 / 2} & 0 \\ \hline-\left[\bar{e}_{k}^{\lambda}\right]^{\top} & {\left[P_{k+1 \mid k}^{-1 / 2} \hat{x}_{k+1 \mid k}{ }^{\top}\right.} & (*)\end{array}\right]}_{\text {Extended post-array } \tilde{\mathbb{R}}_{k}} ;\)
        Read-off from post-array: \(\left[P_{k+1 \mid k}^{1 / 2}\right],\left[P_{k+1 \mid k}^{-1 / 2} \hat{x}_{k+1 \mid k}\right]\);
\(6 \quad\) Compute \(\hat{x}_{k+1 \mid k}=\left[P_{k+1 \mid k}^{1 / 2}\right]\left[P_{k+1 \mid k}^{-1 / 2} \hat{x}_{k+1 \mid k}\right]\).
```

Finally, we note that value $\gamma$ is of no interest for the filtering algorithm above. Hence, we use notation $(*)$ for the block that is not used in the filter. However, it is worth noting here that this quantity turns to be utilized in the square-root Rauch-Tung-Striebel formulas for the smoothed estimate $\hat{x}_{k \mid K}$, if one derives the smoother under the MCC methodology. In this case, the suggested Algorithm 2 yields a simple computation of this value.

## IV. NUMERICAL EXPERIMENTS

The goal of this section is to substantiate the theoretical derivation and correctness of the suggested condensed squareroot Algorithms 1 and 2 on practical example.

Example 1 (see [6]). Consider a benchmark navigation problem where the vehicle dynamics is given as follows:

$$
x_{k}=\left[\begin{array}{cccc}
1 & 0 & T & 0 \\
0 & 1 & 0 & T \\
0 & 0 & 1 & 0 \\
0 & 0 & 0 & 1
\end{array}\right] x_{k-1}+\left[\begin{array}{c}
w_{k-1}^{1} \\
w_{k-1}^{2} \\
w_{k-1}^{3} \\
w_{k-1}^{4}
\end{array}\right]
$$

where $T=0.01$ is the sampling period, the first two state components are the north and east positions of a land vehicle, and the last two components are the north and east velocities. A position-measuring device provides a noisy measurement of the north and east positions as follows:

$$
y_{k}=\left[\begin{array}{llll}
1 & 0 & 0 & 0 \\
0 & 1 & 0 & 0
\end{array}\right] x_{k}+\left[\begin{array}{c}
v_{k}^{1} \\
v_{k}^{2}
\end{array}\right]
$$

The MCC-KF methods are tested at the presence of impulsive noise, i.e., in the presence of outliers. More precisely, the process and measurement noises are generated as follows:

$$
\begin{aligned}
w_{k} & \sim \mathcal{N}(0, Q)+\text { Shot noise }, & & Q=0.1 I_{4} \\
v_{k} & \sim \mathcal{N}(0, R)+\text { Shot noise }, & & R=0.1 I_{2}
\end{aligned}
$$

where $\bar{x}_{0}=[1,1,0,0]^{T}$ and $\Pi_{0}=\operatorname{diag}([4,4,3,3])$.
To simulate the impulsive noise (the shot noise), we follow the approach suggested in [6]. The Matlab routine Shot_noise recently published in [33, Appendix] can be used for generating the process and measurement noise in this set of experiments with $20 \%$ of outliers distributed randomly in the time interval [21, $K-1$ ] (where $K=300$ ) and with the randomly chosen outliers' magnitude from $[0,5]$.

In our numerical simulations, we test all square-root IMCCKF implementations existed nowadays [9], [19]. As mentioned in Introduction, they are developed in two-step a posteriori form, only. The new one-step condensed Algorithms 1 and 2 are examined against a priori one-step IMCC-KF conventional equations (23) - (25). All filtering algorithms under examination utilize the same initial filtering conditions, the same measurements and the same noise covariances. The experiment is repeated for $M=100$ Monte Carlo runs and the root mean square error (RMSE) is calculated for a posteriori and a priori estimates. The results of numerical experiments are summarized in Table [I]
We observe that all examined implementation methods produce the same RMSEs both in a priori and a posteriori forms, respectively. In other words, they work with the same estimation accuracy. For a priori one-step methods derived in this paper, this means that the results of practical example substantiate the correctness of their theoretical derivation presented in Section III This also justifies an algebraic equivalence of new square-root condensed algorithms and the conventional way of calculations by equations (23) - (25).

## V. Conclusion

This paper proposes one-step filtering strategy for the maximum correntropy criterion Kalman filters that corresponds to $a$ priori filtering. It has been shown that the resulted $a$

TABLE I
The RMSE ERRORS FOR IMCC-KF implementations in Example $1, M=100$ Monte Carlo simulations.

| Estimator Implementation method | Covariance factorization | $\mathrm{RMSE}_{x_{1}}$ | $\mathrm{RMSE}_{x_{2}}$ | $\mathrm{RMSE}_{x_{3}}$ | $\mathrm{RMSE}_{x_{4}}$ | $\\|\mathrm{RMSE}\\|_{2}$ |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  | The errors are computed for a priori (predicted) estimates |  |  |  |  |
| IMCC-KF: conventional eqs. (23)-25) | - | 6.7524 | 6.3572 | 38.7909 | 38.5804 | 55.4905 |
| SR IMCC-KF: new Algorithm 1 | Cholesky | 6.7524 | 6.3572 | 38.7909 | 38.5804 | 55.4905 |
| eSR IMCC-KF: new Algorithm 2 | Cholesky | 6.7524 | 6.3572 | 38.7909 | 38.5804 | 55.4905 |
|  |  | The errors are computed for a posteriori (filtered) estimates |  |  |  |  |
| IMCC-KF: eqs. (11), 12), 14, (16)-18) | - | 6.6162 | 6.2099 | 38.7753 | 38.5790 | 55.4455 |
| SR IMCC-KF: Algorithm 2 in [9] | Cholesky | 6.6162 | 6.2099 | 38.7753 | 38.5790 | 55.4455 |
| eSR IMCC-KF: Algorithm 3 in [9] | Cholesky | 6.6162 | 6.2099 | 38.7753 | 38.5790 | 55.4455 |
| UD IMCC-KF: Algorithm 2b in 19] | modified Cholesky | 6.6162 | 6.2099 | 38.7753 | 38.5790 | 55.4455 |
| SVD IMCC-KF: Algorithm 2c in [19] | SVD | 6.6162 | 6.2099 | 38.7753 | 38.5790 | 55.4455 |

priori recursion for the original MCC-KF does not allow for robust square-root implementations because of the possibly negative coefficient involved. In contrast, the improved MCCKF estimator (IMCC-KF) easily permits stable square-root implementations for its a priori recursion in the condensed form. Two Cholesky-based methods of this kind are derived in this paper. Finally, the square-root solution might be derived via $U D U^{\top}$ factorization as well. This problem is still open.

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