An Integrated Framework for Model-Based Distributed Diagnosis and Prognosis

Anibal Bregon\textsuperscript{1}, Matthew Daigle\textsuperscript{2}, and Indranil Roychoudhury\textsuperscript{3}

\textsuperscript{1} University of Valladolid, Valladolid, Spain
anibal@infor.uva.es

\textsuperscript{2} NASA Ames Research Center, Moffett Field, CA 94035, USA
matthew.j.daigle@nasa.gov

\textsuperscript{3} SGT Inc., NASA Ames Research Center, Moffett Field, CA 94035, USA
indranil.roychoudhury@nasa.gov

\textbf{ABSTRACT}

Diagnosis and prognosis are necessary tasks for system reconfiguration and fault-adaptive control in complex systems. Diagnosis consists of detection, isolation and identification of faults, while prognosis consists of prediction of the remaining useful life of systems. This paper presents a novel integrated framework for model-based distributed diagnosis and prognosis, where system decomposition is used to enable the diagnosis and prognosis tasks to be performed in a distributed way. We show how different submodels can be automatically constructed to solve the local diagnosis and prognosis problems. We illustrate our approach using a simulated four-wheeled rover for different fault scenarios. Our experiments show that our approach correctly performs distributed fault diagnosis and prognosis in an efficient and robust manner.

\textbf{1. INTRODUCTION}

Systems health monitoring is essential to guaranteeing the safe, efficient, and reliable operation of engineering systems. Integrated systems health management methodologies include fault diagnosis and prognosis mechanisms, where diagnosis involves detecting when a fault has occurred, isolating the true fault, and identifying the true damage to the system; and prognosis involves predicting how much useful life remains in the different components, subsystems, or systems given the diagnosed fault conditions. The information on the fault size and its expected impact on system life can be used to initiate recovery and reconfiguration actions that mitigate the fault or extend system life.

A large body of research exists for both model-based diagnosis (Gertler, 1998; Blanke et al., 2006) and prognosis methods (Luo et al., 2008; Saha & Goebel, 2009; Orchard & Vachtsevanos, 2009). However, the integration of diagnosis and prognosis algorithms is seldom studied. In fact, many diagnosis methodologies leave out the fault identification step, which is necessary to perform a prediction from the current system state. Recently, we presented an integrated model-based framework for diagnosis and prognosis of complex systems, in which we made use of a common modeling framework for modeling both the nominal and faulty system behavior (Roychoudhury & Daigle, 2011).

In (Roychoudhury & Daigle, 2011), the nominal system behavior is estimated using an observer built with the nominal model. Faults are detected when a statistically significant deviation between the nominal estimates and the observed measurements is observed (Biswas et al., 2003). Fault isolation compares the observed measurement deviations against predictions of how the measurements would deviate for each possible fault (Mosterman & Biswas, 1999). Fault identification performs joint state-parameter estimation using multiple observers, where, for each fault, the faulty system model is constructed as the nominal model integrated with a hypothesized fault model (Roychoudhury, 2009). The prognosis module uses, for each fault hypothesis, a prediction model based on its faulty system model and the identified fault parameters, to predict the remaining useful life of the system (Daigle, Saha, & Goebel, 2012). However, this integrated solution performs the diagnosis and prognosis task in a centralized fashion, which is prone to single points of failure, and does not scale well as the size of the system increases.

To overcome such problems, in this work, we leverage recent results for distributed diagnosis (Bregon et al., 2011)
and distributed prognosis (Daigle, Bregon, & Roychoudhury, 2012), which make use of structural model decomposition techniques, to provide a systematic approach to distributing the different diagnosis and prognosis steps presented in (Roychoudhury & Daigle, 2011).

Distributed diagnosis is achieved by designing local distributed subsystems based on global diagnosability analysis of the system, thus computing globally correct distributed diagnosis results without the use of a centralized coordinator (Bregon et al., 2011). These local distributed subsystems are then used to construct local event-based distributed diagnosers for distributed fault isolation. Distributed fault identification is achieved by developing independent local state-parameter estimators for each hypothesized fault. Regarding distributed prediction, in (Daigle, Bregon, & Roychoudhury, 2012) we developed an architecture that enables a large prognosis problem to be decomposed into several independent local subproblems from which local results can be merged into a global result.

The main contribution of this paper is an integrated framework for distributed model-based diagnosis and prognosis of single faults based on structural model decomposition. The proposed framework scales well and the resulting subproblems are typically small and easy to solve, resulting in an efficient and scalable distributed solution to the combined diagnosis and prognosis problem. We perform a number of experiments on a simulated four-wheeled rover testbed (Balaban et al., 2011) to demonstrate and evaluate our approach.

The rest of the paper is organized as follows. Section 2 provides the problem formulation for our diagnosis and prognosis framework. Section 3 describes the distributed architecture and Section 4 briefly introduces its different components. Section 5 presents the case study and experimental results. Finally, Section 6 concludes the paper.

2. Problem Formulation

The nominal system model is represented as follows:

\[ \dot{x}(t) = f(t, x(t), \theta(t), u(t), v(t)), \]
\[ y(t) = h(t, x(t), \theta(t), u(t), n(t)), \]

where \( x(t) \in \mathbb{R}^{n_x} \) is the state vector, \( \theta(t) \in \mathbb{R}^{n_\theta} \) is the parameter vector, \( u(t) \in \mathbb{R}^{n_u} \) is the input vector, \( v(t) \in \mathbb{R}^{n_v} \) is the process noise vector, \( f \) is the state equation, \( y(t) \in \mathbb{R}^{n_y} \) is the output vector, \( n(t) \in \mathbb{R}^{n_n} \) is the measurement noise vector, and \( h \) is the output equation.\(^1\)

Faults in the system are represented as changes in the above nominal system model. In this work, we only consider single faults occurring as changes in system parameters, \( \theta(t) \). We denote a fault, \( f \in F \), as a tuple, \((\theta, g_f)\), where \( \theta \in \Theta \) is the fault parameter, and \( g_f \) denotes the fault progression function, which models the way fault \( f \) is manifested in parameter \( \theta \), i.e.,

\[ \dot{\theta}(t) = g_f(t, x_f(t), \theta(t), u(t), m_f(t)), \]

where \( x_f(t) = [x(t), \theta(t)]^T \), \( \theta_f(t) = [\theta(t) \setminus \{\theta(t)\}] \), \( \phi_f(t) \in \mathbb{R}^{n_\phi} \) is a vector of fault progression parameters, and \( m_f(t) \in \mathbb{R}^{n_m} \) is a process noise vector associated with the fault progression function.

To develop our integrated diagnosis and prognosis framework, the faulty system model for fault \( f = (\theta, g_f) \) is constructed from the nominal system model by including the parameter as a state and augmenting the state equation by including the fault progression function, i.e.,

\[ \dot{x}_f(t) = f_f(t, x_f(t), \theta_f(t), u(t), v(t)), \]
\[ y(t) = h(t, x(t), \theta(t), u(t), n(t)), \]

where,

\[ f_f(\cdot) = \begin{bmatrix} f(t, x(t), \theta(t), u(t), v(t)) \n g_f(t, x_f(t), \theta_f(t), u(t), m_f(t)) \end{bmatrix} = \begin{bmatrix} x(t) \n \theta(t) \end{bmatrix} \]

The goal of diagnosis is to: (i) detect a change in some \( \theta \in \Theta \); (ii) isolate, under the single fault assumption, the true fault \( f \in F \), i.e., both the parameter \( \theta \) that has changed, and its fault progression function \( g_f \); and (iii) identify (i.e. estimate) the fault by computing \( p(x_f(t), \theta_f(t)|y(0:t)) \), where \( y(0:t) \) denotes all measurements observed up to time \( t \).

The goal of prognosis is to determine the end of (useful) life (EOL) of a system, and/or its remaining useful life (RUL). For a given fault \( f \), using the fault estimate, \( p(x_f(t), \theta_f(t)|y(0:t)) \), a probability distribution of EOL, \( p(EOL_f(t_P)|y(0:t_P)) \), and/or RUL, \( p(RUL_f(t_P)|y(0:t_P)) \) is computed at a given time point \( t_P \) (Daigle, Saha, & Goebel, 2012). Since the prognosis problem is stochastic, EOL/RUL are random variables and we represent them by probability distributions. The acceptable behavior of the system is expressed through a set of \( n_c \) constraints, \( C_{EOL_f} = \{c_i\}_{i=1}^{n_c} \), where \( c_i : \mathbb{R}^{n_x} \times \mathbb{R}^{n_\theta} \times \mathbb{R}^{n_u} \rightarrow \mathbb{B} \) maps a given point in the joint state-parameter space given the current inputs, \( (x_f(t), \theta_f(t), u(t)) \), to the Boolean domain \( \mathbb{B} \triangleq [0,1] \), where \( c_i(x_f(t), \theta_f(t), u(t)) = 1 \) if the constraint is satisfied (Daigle, Saha, & Goebel, 2012). If \( c_i(x_f(t), \theta_f(t), u(t)) = 0 \), then the constraint is not satisfied, and the behavior of the system is deemed to be unacceptable. These individual constraints are combined into a single threshold function \( T_{EOL_f} : \mathbb{R}^{n_x} \times \mathbb{R}^{n_\theta} \times \mathbb{R}^{n_u} \rightarrow \mathbb{B} \),

\[^1\]Here, we use boldface typeface to denote vectors, and use \( n_a \) to denote the length of a vector \( a \).
of (Submodel) Definition 2 from observed measurement deviations. A submodel in our sufficient analytical redundancy to generate fault hypotheses compose the global system model into a set of submodels sat-
al., 2006; Pulido & Alonso-González, 2004) to compute lo-
cal independent subproblems, which may be solved in paral-
lel, thus providing scalability and efficiency. Model decom-
position problem, constraints are that submodels use
the global model inputs and some measured values as local
inputs, i.e., $U_i \subseteq U \cup (Y - Y_i)$. An algorithm for computing the set of minimal submodels that satisfies these properties is given in (Daigle et al., 2011), which is based on the model decomposition algorithms presented in (Pulido & Alonso-González, 2004; Bregon et al., 2012).

The basic idea of the model decomposition problem is to de-
compose the global system model into a set of submodels sat-
ifying given constraints, such that each submodel contains sufficient analytical redundancy to generate fault hypotheses from observed measurement deviations. A submodel in our framework is defined as follows.

Definition 2 (Submodel). A submodel $M_i$ of a system model $M = (X, \Theta, U, Y, C)$ is a tuple $M_i = (X_i, \Theta_i, U_i, Y_i, C_i)$, where $X_i \subseteq X$, $\Theta_i \subseteq \Theta$, $U_i \subseteq X \cup U \cup Y$, and $Y_i \subseteq Y$ are the state, parameter, input, and output variables, respectively, and $C_i \subseteq C$ are the submodel constraints.$^2$

Next, we discuss the fundamental ideas of our model decomposition approach and show the constraints needed to obtain the different submodels for distributed diagnosis and prognosis. Then, we propose our integrated approach.

3. Distributed Architecture

For a large system, both the diagnosis and prognosis problems are correspondingly large. A centralized approach does not scale well, can be computationally expensive, and prone to single points of failure. Therefore, we propose to decom-
pose the global integrated diagnosis and prognosis problem into independent local subproblems. In this work, we build on the ideas from structural model decomposition (Blanke et al., 2006; Pulido & Alonso-González, 2004) to compute local independent subproblems, which may be solved in parallel, thus providing scalability and efficiency. Model decomposition is not a new concept, and several approaches have been developed for purposes of system identification, estimation, learning, and diagnosis (Staroswiecki & Declerck, 1989; Pulido & Alonso-González, 2004; Williams & Millar, 1998). Structural model decomposition allows decomposing a global model into a set of local submodels for which local diagnosis and prognosis problems can be directly defined. The global model of the system, denoted as $M$, is defined as follows.

Definition 1 (Model). The model of a system, $M$, is a tuple $M = (X, \Theta, U, Y, C)$, where $X$ is the set of state variables of $x$, $\Theta$ is the set of unknown parameters of $\theta$, $U$ is the set of input variables of $u$, $Y$ is the set of output variables of $y$, and $C$ is the set of model constraints of $f$, $h$, and EOL constraints of $C_{EOL_f}$.

The third model decomposition problem computes submod-
els for residual generation and fault isolation. For this model decomposition problem, the constraints are that sub-
models are constructed by merging the minimal submodels, to fulfill global diagnosability conditions as in (Bregon et al., 2011).$^3$ Once the globally diagnosable subsystems have been designed, the merged submodels are used for distributed residual generation and to compute event-based local diagnosers for fault isolation. These design and diagnoser computation processes are detailed in (Bregon et al., 2011). Because the subsystems are designed to be globally diagnosable, the resulting local diagnosers are independent, and can pro-
vide globally correct diagnosis results without a centralized coordinator.

The third model decomposition problem computes submod-
els for distributed fault identification. For each consistent fault hypothesis $f$, the joint state-parameter estimators are computed from the minimal submodels of the faulty system model $M(f)$ with the constraints that $U_i \subseteq U \cup (Y - Y_i)$ with

A faulty submodel for a fault $f$ is denoted as $M_i(f)$.

$^2$A faulty submodel for a fault $f$ is denoted as $M_i(f)$.

$^3$In this work, a subsystem is globally diagnosable if all faults in the subsys-
tem are distinguishable from every other fault in the system using only local measurements.
\( Y_i \) as a singleton. It will be shown later that the fault identification module is the central part of our diagnosis-prognosis integration approach and provides the joint state-parameter estimations for the prediction module.

Finally, for distributed prediction, the model decomposition problem starts off from the faulty system model, and, as detailed in (Daigle, Bregon, & Roychoudhury, 2012), it fulfills the following constraints: (i) the submodels use \( U_i \subseteq U_P \), where \( U_P \subseteq X \cup U \) (here, \( U_P \) is a set of variables whose future values can be predicted \textit{a priori}, which depends on the hypothesized faults); and (ii) each computed submodel has at least one \( c \in C_{EOL} \) belonging to \( C_i \), and over all submodels, all constraints in \( C_{EOL} \) are covered. This ensures that \( T_{EOL} \) may be computed for the system from the local constraints.

### 3.2. Distributed Architecture

Figure 1 illustrates an example architecture for our distributed diagnosis and prognosis scheme. At each discrete time step, \( k \), the system takes as input both \( u_k \) and \( y_k \) and splits them into local inputs \( u^i_k \) and local outputs \( y^i_k \) for the local diagnosers. Within each \( M_i \) local diagnoser, nominal tracking is performed, computing estimates of nominal measurements, \( \hat{y}_k \). The fault detector compares the estimated measurements against the observed measurements, to determine statistically significant deviations for the residual, \( r^i_k = y^i_k - \hat{y}_k \). Qualitative values of the deviations in the residuals are used by the event-based diagnoser to isolate faults. The set of isolated fault candidates \( \mathbf{F}^i \) together with the estimated nominal states, \( \hat{x}^i_k \), parameters, \( \hat{\theta}^i_k \), and the measurements, \( y^i_k \), are used as input for the corresponding identification module. Identification is performed for each hypothesized fault in a distributed way, e.g., for the isolated faults \( f_1 \) and \( f_2 \) in Figure 1, we run an instantiation of the identification submodel for each one the faults, i.e., \( M_1(f_1) \) and \( M_1(f_2) \). Fault identification uses the minimal submodels from the faulty system model, and computes local state-parameter estimates \( p(\mathbf{x}^i_{f,k}; \mathbf{\theta}^i_{f,k}|\mathbf{y}^i_{0:k}) \). These local estimates are then used as input to the prediction submodels. In some cases, the local estimates have to be split or merged with other estimates according to the prediction submodels. For example, in the figure, estimates from \( M_1(f_1) \) are used by both local prediction submodels \( M_2(f_1) \) and \( M_5(f_1) \), and those submodels may also need estimates not included within submodel \( M_1(f_1) \). These estimates are typically obtained from the local diagnosers or other fault identification blocks.\(^4\)

Distributed prediction modules compute, for each hypothesized fault, local EOL/RUL predictions, \( p(EOL^i_{f,k,p}|\mathbf{y}^i_{0:k}) \) and \( p(RUL^i_{f,k,p}|\mathbf{y}^i_{0:k,p}) \), at given prediction time \( k_p \) based on the local EOL constraints. Finally, local predictions are combined into global predictions \( p(EOL_{f,k,p}|\mathbf{y}^i_{0:k}) \) and \( p(RUL_{f,k,p}|\mathbf{y}^i_{0:k}) \) for each hypothesized fault. The next section describes the details of the different modules of the distributed integrated diagnosis and prognosis architecture.

### 4. Diagnosis and Prognosis Approach

Figure 1 shows the basic modules of our distributed integrated approach. In this section we give details on how each module is implemented, and establish the integration between the diagnosis and prognosis tasks.

#### 4.1. Distributed Diagnosis

For distributed diagnosis, each local diagnoser first takes a subset of the local inputs \( u_k^i \) and local outputs \( y_k^i \), to compute an estimate of its output measurements \( \hat{y}^i_k \). Tracking is performed in discrete time using a robust filtering scheme, e.g., the extended or unscented Kalman filter (Julier & Uhlmann, 2004), which provides accurate tracking in the presence of sensor noise, process noise, and discretization error.

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\(^4\)Since prediction submodels are constructed by using any variable which value can be hypothesized as input, in some cases, prediction submodels cannot always be formed by just merging the minimal estimation submodels. To indicate this, we named our prediction submodels differently from the estimation submodels, e.g., \( M_4(f_1) \) and \( M_5(f_1) \) instead of \( M_1(f_1) \).
For fault detection, a statistical test is used to look for significant deviations in the residual signal $r^i_k$, which is computed as the difference between $\hat{y}^i_k$ and the $y^i_k$. In our approach, we use a Z-test as described in (Biswas et al., 2003).

Fault isolation is performed using local event-based diagnosters, constructed as detailed in the previous section (Bregon et al., 2011; Daigle et al., 2009). Fault isolation is triggered when a fault is detected, and it works as follows. Initially, all event-based local diagnosers start in their initial state, and the set of faulty candidates is empty. Local residual deviations cause the local diagnosers to move from one state to another. These residual deviations are abstracted to a tuple of qualitative symbols $(\sigma_1, \sigma_2)$ for each residual signal, where $\sigma_1$ represents magnitude changes and $\sigma_2$ represents slope changes. A $(\sigma_1, \sigma_2)$ value indicates a change above normal for a measurement residual or a positive (resp. below) normal for a measurement residual or a flat residual slope. The symbols are utilized to abstract the qualitative deviations in the measurement value or a flat residual slope. A $0$ implies no change in the measurement value or a flat residual slope. The symbols are generated using a sliding window technique as described in detail in (Biswas et al., 2003). If there is a match between an event from the current state and a tuple of qualitative symbols generated by any residual, the local diagnoser moves to the next state and remains active. If not, the local diagnoser blocks. This process continues until a local diagnoser reaches an accepting state, which corresponds to a unique isolation result.

In our distributed diagnosis approach, identification submodels, $M_i(f)$, are obtained, as explained in the previous section, as minimal submodels from the faulty system model. A local state-parameter estimator is constructed for each identification submodel $M_i(f)$, and produces a local estimate $p(x_{f,k}^i, \theta_{f,k}^i | y_{0:k}^i)$ by using an appropriate algorithm. In this paper, we use an unscented Kalman filter (UKF) (Julier & Uhlmann, 2004) with a variance control algorithm (Daigle, Saha, & Goebel, 2012).

4.2. Distributed Prognosis

The local state-parameter estimates for each local distributed prediction module are constructed from the local estimates of the distributed fault identification submodels. Each prediction submodel is made up of a set of states $X_i$ and parameters $\Theta_i$, and constructs a local distribution $p(x_{f,k}^i, \theta_{f,k}^i | y_{0:k}^i)$, by assuming that the local state-parameter estimates are sufficiently represented by a mean $\mu^i$ and covariance $\Sigma^i$. For each prediction submodel $M_i(f)$, we combine the estimates of the local identification submodels that estimate states and parameters in $X_i \cup \Theta_i$, into $\mu^i$ and $\Sigma^i$. If two submodels estimate the same state variable or parameter, then many different techniques can be applied depending on the desired performance of the prediction submodels, e.g., taking the estimate with the smallest variance, or taking an average.

Several approaches can be used to perform prediction for each prediction submodel. In this work, given the mean and covariance information, we represent the distribution with a set of sigma points derived using the unscented transform. Then, each sigma point is simulated forward to EOL, and we recover the statistics of the EOL distribution given by the sigma points (Daigle & Goebel, 2010).

Algorithm 1 (Daigle, Saha, & Goebel, 2012), shows the pseudocode for the prediction procedure. The algorithm is executed for each submodel $i$, deriving local EOL predictions using its local threshold function based on the local EOL constraints. For a given submodel, each sample $j$ is propagated forward until $T_{EOL}^i(x_{f,k}^i, \theta_{f,k}^i)$ evaluates to 1. The algorithm hypothesizes future inputs $\hat{u}_{f,k}$. Then, the EOL/RUL is determined by the minimum of the local EOL/RUL distributions for each prediction submodel, i.e., $p(\text{EOL}_{f,k}^i, \text{RUL}_{f,k}^i | y_{0:k}^i)$. To compute this, we sample from each local EOL distribution and take the minimum of the local samples. This is repeated many times and the statistics of the global EOL distribution are computed (Daigle, Bregon, & Roychoudhury, 2012).

5. CASE STUDY

In this section, we apply our distributed diagnosis and prognosis approach to a four-wheeled rover testbed developed at NASA Ames Research Center. We develop a model of the rover, and demonstrate the approach using simulated scenarios.

5.1. Nominal System Modeling

The rover model was originally presented in (Balaban et al., 2011). In this section we summarize the main features and include some extensions to the model.

The rover consists of a symmetric rigid frame with four
where \( m \) is the body mass, \( \tau \) is the torque, \( \tau_{gl} \) is the torque due to slippage, \( \mu \) is a friction coefficient, and \( \omega \) is the rotational velocity of the rover body; and \( \tau_{gr} \) is the torque due to the translational movement of the rover body, where \( \mu_{gr} \) is a friction coefficient, \( \rho \) is the rotational velocity of the rover body, and \( \gamma = \arctan \frac{l}{b} \) with \( l \) being the rover’s length and \( b \) being its width.

The translational velocity \( v \) of the rover is described by

\[
\dot{v} = \frac{1}{m} \left( F_{pFL} + F_{pFR} + F_{pBL} + F_{pBR} \right),
\]

where \( m \) is the rover mass, and for wheel \( w, F_{glw} = \mu_{gl} (v_w - v) \) is the force due to slippage. The rotational velocity \( \omega \) is described by

\[
\dot{\omega} = \frac{1}{J} \left( d \cos \gamma F_{pFR} + d \cos \gamma F_{pBR} - d \cos \gamma F_{pFL} - d \cos \gamma F_{pBL} - dF_{pFR} - dF_{pBR} - dF_{pFL} - dF_{pBL} \right),
\]

where \( J \) is the rotational inertia of the rover and \( d \) is the distance from the center of the rover to each wheel.

The wheels are driven by DC motors with PI control that sets the voltages \( V \) applied to the motors. The motor currents \( i \) are governed by

\[
\dot{i}_{FL} = \frac{1}{L} \left( V_{FL} - i_{FL} R_{FL} - k_e \omega_{FL} \right),
\]

\[
\dot{i}_{FR} = \frac{1}{L} \left( V_{FR} - i_{FR} R_{FR} - k_e \omega_{FR} \right),
\]

\[
\dot{i}_{BL} = \frac{1}{L} \left( V_{BL} - i_{BL} R_{BL} - k_e \omega_{BL} \right),
\]

\[
\dot{i}_{BR} = \frac{1}{L} \left( V_{BR} - i_{BR} R_{BR} - k_e \omega_{BR} \right),
\]

where \( L \) is the motor inductance, \( R \) is the motor resistance, and \( k_e \) is an energy transformation term. The motor torque is \( \tau_{mw} = k_e i_{mw} \), where \( k_e \) is an energy transformation gain. The voltages applied to the motors are determined by the controllers, where for wheel \( w, V_w = P (u_w - \omega_w) + I \cdot e_{iw} \), where \( P \) is a proportional gain, \( u_w \) is the commanded wheel speed, \( I \) is an integral gain, and \( e_{iw} \) is the integral error term.

The integral error terms are governed by

\[
\dot{e}_{FL} = u_{FL} - \omega_{FL},
\]

\[
\dot{e}_{FR} = u_{FR} - \omega_{FR},
\]

\[
\dot{e}_{BL} = u_{BL} - \omega_{BL},
\]

\[
\dot{e}_{BR} = u_{BR} - \omega_{BR}.
\]

The batteries, which are connected in series, are described by a simple electrical circuit equivalent model that includes a large capacitance \( C_b \) in parallel with a resistance \( R_p \), together in series with another resistance \( R_s \). The battery charge variables \( q_i \) are governed by

\[
\dot{q}_1 = -V_1/R_{p1} - (i_{FL} + i_{FR} + i_{BL} + i_{BR}),
\]

\[
\dot{q}_2 = -V_2/R_{p2} - (i_{FL} + i_{FR} + i_{BL} + i_{BR}),
\]

\[
\dot{q}_3 = -V_3/R_{p3} - (i_{FL} + i_{FR} + i_{BL} + i_{BR}),
\]

\[
\dot{q}_4 = -V_4/R_{p4} - (i_{FL} + i_{FR} + i_{BL} + i_{BR}).
\]

The available sensors measure the voltages of the batteries,

\[
V_{1}^* = q_1/C_b0 - R_s * (i_{FL} + i_{FR} + i_{BL} + i_{BR}),
\]

\[
V_{2}^* = q_2/C_b1 - R_s * (i_{FL} + i_{FR} + i_{BL} + i_{BR}),
\]

\[
V_{3}^* = q_3/C_b2 - R_s * (i_{FL} + i_{FR} + i_{BL} + i_{BR}),
\]

\[
V_{4}^* = q_4/C_b3 - R_s * (i_{FL} + i_{FR} + i_{BL} + i_{BR}),
\]

the motor currents,

\[
i_{FL}^* = i_{FL},
\]

\[
i_{FR}^* = i_{FR},
\]

\[
i_{BL}^* = i_{BL},
\]

\[
i_{BR}^* = i_{BR},
\]

and the wheel speeds,

\[
\omega_{FL}^* = \omega_{FL},
\]

\[
\omega_{FR}^* = \omega_{FR},
\]

\[
\omega_{BL}^* = \omega_{BL},
\]

\[
\omega_{BR}^* = \omega_{BR}.
\]

Here, the * superscript indicates a measured value.

**5.2. Faulty System Modeling**

In this work, we consider different faults in the motors and the batteries. First, we consider friction-based damage progression in the motors, resulting in an increase in motor friction over time. For wheel \( w \), the fault progression function is defined as:

\[
\dot{\mu}_{FL} = V_{FL} \mu_{FL} \omega_{FL}^2,
\]

\[
\dot{\mu}_{FR} = V_{FR} \mu_{FR} \omega_{FR}^2,
\]

\[
\dot{\mu}_{BL} = V_{BL} \mu_{BL} \omega_{BL}^2,
\]

\[
\dot{\mu}_{BR} = V_{BR} \mu_{BR} \omega_{BR}^2,
\]

We use a simple model here only for demonstration purposes. More detailed battery models for prognosis can be found in the literature, e.g., (Saha & Goebel, 2009).
where the charge threshold is given by $q$. Table 2 shows the submodels for residual generation represented as an abrupt change in parameter $R$. For the batteries, we consider abrupt capacitance decreases, of the subsystems. In (Bregon et al., 2011) determines that we need to merge the minimal submodels in Table 1 by defining the fault progression parameter to become the parameter. Regarding progression function (if necessary), and by making the fault parameter, and $v_{fw}$ is the fault progression parameter.

We also consider abrupt resistance increases in the motors, represented as an abrupt change in parameter $R_w$ for wheel $w$, with $\Delta R_w$ as the fault progression parameter.

For the batteries, we consider abrupt capacitance decreases, represented as an abrupt change in parameter $C_{bi}$ for capacity $i$. $\Delta C_{bi}$ is the fault progression parameter.

We are interested in predicting when any of the rover batteries are at their charge threshold, beyond which the batteries will be damaged. These faults can cause the charge thresholds to be reached earlier since they will affect current draw. The constraints are given as

$$q_1 > q^-,$$  \hspace{1cm} (c_{35})

$$q_2 > q^-,$$  \hspace{1cm} (c_{36})

$$q_3 > q^-,$$  \hspace{1cm} (c_{37})

$$q_4 > q^-,$$  \hspace{1cm} (c_{38})

where the charge threshold is given by $q^- = 2 \times 10^4$ C. The rover cannot be operated when any of the constraints $c_{35}$-$c_{38}$ are violated.

### 5.3. Results

To demonstrate the validity of the approach, we describe two different faulty scenarios of the rover. In the first, friction damage is progressing on one motor, and in the second, a capacitance decrease occurs in one battery. In all cases, the rover travels between various waypoints, moving at an average speed of 0.5 m/s. Table 1 shows the minimal submodels for the rover derived by using measured values as local inputs. Table 2 shows the submodels for residual generation and fault isolation. These submodels have been designed to obtain globally diagnosable subsystems by using the design algorithm in (Bregon et al., 2011). In this work, we have considered five subsystems, one for each wheel components and another one for the batteries. For example, the subsystem for the front left wheel components is not globally diagnosable if we only consider submodel $M_5$ (which includes the front left wheel friction wear parameter, $v_{fFL}$). The design algorithm in (Bregon et al., 2011) determines that we need to merge submodels $M_5$ and $M_9$ to make the front left wheel subsystem globally diagnosable. The process is similar for the rest of the subsystems.

Minimal submodels for identification can be computed from the minimal submodels in Table 1 by defining the fault progression function (if necessary), and by making the fault parameter to become a state and the corresponding fault progression parameter to become the parameter. Regarding prediction, the correct prediction submodels to use depend on the scenario, as will be shown later.

### 5.3.1. Friction Damage Progression

We first consider a scenario in which, for the front-left motor, the friction begins to increase. The friction damage progression begins at $t = 50$ s with friction wear parameter $v_{fFL} = 1 \times 10^{-3}$ s. A fault is detected by the local di-
As a result of the continuously increasing friction, the current drawn by the motor increases as well in order for the motor controller to maintain the same desired wheel speed (Figure 3 shows this increase in the current through time). Hence, the total current drawn from the batteries is increased, and EOL occurs around half an hour. Because \( i_{FL} \) is constantly changing, and in a way that is dependent on the motor state, it is incorrect to use it as a local input for prediction and to decompose the prediction problem into independent local prediction problems for the batteries and motors, i.e., it is not known a priori. Therefore, we compute submodels using as local inputs average values for the remaining motor currents, average commanded wheel speeds, and average rover translational velocity \( v \) and rotational velocity \( \omega \). The prediction submodels for this case are shown in Table 3. EOL for this fault is computed by merging the local EOL from those submodels in the table. Note that the prediction submodels used in this case do not correspond directly to those used for estimation. So, when constructing the estimate for \( \mathcal{M}_{13} \), for example, it takes the estimates from \( \mathcal{M}_1 \) and \( \mathcal{M}_9 \) as a particular example, but, in general, both approaches obtain equivalent results.

The prediction results are shown in Figure 4. The increased friction causes the batteries to discharge faster, and EOL occurs around 1650 s. Here, we used the relative accuracy (RA) as a measure of prediction accuracy, and the relative standard deviation (RSD) as a measure of spread. Each prediction metric is averaged over multiple prediction points (one every 100 s of usage) (see (Saxena et al., 2010; Daigle, Saha, & Goebel, 2012) for the mathematical definitions of these metrics). For this experiment, RA averages to 91.63\% and RSD averages to 16.26\%.

For the sake of comparison, we also ran this experiment using the centralized approach. Figure 5 shows the prediction results obtained. Looking at the prediction metrics, we see that the centralized approach behaved very similar to the distributed approach but a little bit worse, with RA averaging 90.90\% and RSD averaging 17.72\%. However, this is just a particular example, but, in general, both approaches obtain equivalent results.

### 5.3.2. Capacitance Decrease

As a second scenario, we consider a capacitance decrease fault in battery 3 of the rover, \( C_{b3} \). The fault begins at \( t = 50 \) s with an abrupt decrease from 2000 to 1800 in the capacity of the battery. The fault is detected immediately by the local diagnoser computed from submodel \( \mathcal{M}_{1,2,3,4} \) at 50.0 s, via an increase in the voltage \( V_i \). The fault candidate is immediately isolated, \( \{C_{b3}\} \), based on the signatures and orderings, thus starting the fault identification. For the capacitance fault, the estimated value of the capacitance averaged \( C_{b3} = 1798.6 \) C with very small output error. As a result of the decrease in capacitance, the battery discharges at a faster rate, and so reaches end of discharge more quickly. The prediction sub-

<table>
<thead>
<tr>
<th>Submodel</th>
<th>( \Theta )</th>
<th>( U )</th>
<th>( Y )</th>
</tr>
</thead>
<tbody>
<tr>
<td>( \mathcal{M}<em>{13}(\nu</em>{FL}) )</td>
<td>( q_{FL}, i_{FL}, e_{FL}, \omega_{FL}, \mu_{FL}, C_{FL}, R_{FL} )</td>
<td>( u_{FL}, v, \omega, i_{FR}, i_{BL}, i_{BR} )</td>
<td>( c_{1, c_{13}, c_{7}, c_{11}, c_{15}, c_{35}} )</td>
</tr>
<tr>
<td>( \mathcal{M}<em>{12}(\nu</em>{FL}) )</td>
<td>( q_{3}, i_{FL}, e_{FL}, \omega_{FL}, \mu_{FL}, C_{b3}, \nu_{FL}, R_{FL} )</td>
<td>( u_{FL}, v, \omega, i_{FR}, i_{BL}, i_{BR} )</td>
<td>( c_{1, c_{13}, c_{7}, c_{11}, c_{16}, c_{36}} )</td>
</tr>
<tr>
<td>( \mathcal{M}<em>{10}(\nu</em>{FL}) )</td>
<td>( q_{4}, i_{FL}, e_{FL}, \omega_{FL}, \mu_{FL}, C_{b4}, \nu_{FL}, R_{FL} )</td>
<td>( u_{FL}, v, \omega, i_{FR}, i_{BL}, i_{BR} )</td>
<td>( c_{1, c_{13}, c_{7}, c_{11}, c_{16}, c_{38}} )</td>
</tr>
</tbody>
</table>

Table 3. Prediction submodels using commanded wheel speeds and rover velocities as local inputs.
models for faults in the capacity of the batteries are shown in Table 4. For this scenario, with a fault in $C_{b3}$, we used submodel $M_{19}$, obtaining RA average to 98.25% and RSD average to 10.12%.

6. CONCLUSIONS

This paper presented a distributed integrated model-based diagnosis and prognosis framework. Our approach starts off with a common modeling paradigm to model both the nominal behavior and fault progression, and then proposes a framework where the global system model is decomposed into smaller independent submodels. These submodels are then used to distribute the different diagnosis and prognosis tasks. Model decomposition is carried out based on the requirements and constraints of each task. We demonstrated our approach on a four-wheeled rover testbed, where we diagnosed faults and prognosed the EOL/RUL accurately. We compared results obtained by using our distributed approach against those obtained using a centralized approach, showing that both approaches obtain the same results.

Most approaches in the literature focus in either the diagnosis or the prognosis task. Some works have proposed the integration of both tasks within a common framework (Patrick et al., 2007; Orchard & Vachtsevanos, 2009; Roychoudhury & Daigle, 2011), however, unlike our approach, these approaches perform the diagnosis and prognosis tasks in a centralized way, thus suffering from scalability issues due to the large number of states and parameters in real-world systems. To the best of our knowledge, there is no approach in the literature which combines, in a distributed way, the diagnosis and prognosis tasks. Our approach is limited by the number and location of the sensors in the system. Since our decomposition algorithm is guided by the set of available sensors, the distribution capabilities of the approach is determined by them.

In future, we will apply this approach to larger systems, to study the scalability of our diagnosis and prognosis scheme, and will perform a more detailed comparison against the results obtained by using a centralized approach. We will also extend the capability of this approach to hybrid systems, as well as diagnosis and prognosis of multiple faults.

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BIOGRAPHIES

Anibal Bregon received his B.S., M.S., and Ph.D. degrees in Computer Science from the University of Valladolid, Spain, in 2005, 2007, and 2010, respectively. From September 2005 to June 2010, he was Graduate Research Assistant
with the Intelligent Systems Group at the University of Valladolid, Spain. He has been visiting researcher at the Institute for Software Integrated Systems, Vanderbilt University, Nashville, TN, USA; the Dept. of Electrical Engineering, Linkoping University, Linkoping, Sweden; and the Diagnostics and Prognostics Group, NASA Ames Research Center, Mountain View, CA, USA. Since September 2010, he has been Assistant Professor and Research Scientist at the Department of Computer Science from the University of Valladolid.

Dr. Bregon is a member of the Prognostics and Health Management Society and the IEEE. His current research interests include model-based reasoning for diagnosis, prognostics, health-management, and distributed diagnosis of complex physical systems.

Matthew Daigle received the B.S. degree in Computer Science and Computer and Systems Engineering from Rensselaer Polytechnic Institute, Troy, NY, in 2004, and the M.S. and Ph.D. degrees in Computer Science from Vanderbilt University, Nashville, TN, in 2006 and 2008, respectively.

From September 2004 to May 2008, he was a Graduate Research Assistant with the Institute for Software Integrated Systems and Department of Electrical Engineering and Computer Science, Vanderbilt University, Nashville, TN. During the summers of 2006 and 2007, he was an intern with Mission Critical Technologies, Inc., at NASA Ames Research Center. From June 2008 to December 2011, he was an Associate Scientist with the University of California, Santa Cruz, at NASA Ames Research Center. Since January 2012, he has been with NASA Ames Research Center as a Research Computer Scientist. His current research interests include physics-based modeling, model-based diagnosis and prognosis, simulation, and hybrid systems.

Dr. Daigle is a member of the Prognostics and Health Management Society and the IEEE. He is a recipient of a University Graduate Fellowship from Vanderbilt University, a best paper award in the Annual Conference of the Prognostics and Health Management Society 2011, a NASA Ames Group Achievement Award in 2011, and an Ames Contractor Council Excellence Award in 2011. He has published over 40 peer-reviewed papers in the area of Systems Health Management.

Indranil Roychoudhury received the B.E. (Hons.) degree in Electrical and Electronics Engineering from Birla Institute of Technology and Science, Pilani, Rajasthan, India in 2004, and the M.S. and Ph.D. degrees in Computer Science from Vanderbilt University, Nashville, Tennessee, USA, in 2006 and 2009, respectively. Since August 2009, he has been with SGT, Inc., at NASA Ames Research Center as a Computer Scientist.

Dr. Roychoudhury is a member of the Prognostics and Health Management Society and the IEEE. His research interests include hybrid systems modeling, model-based diagnostics and prognostics, distributed diagnostics and prognostics, and Bayesian diagnostics of complex physical systems.