

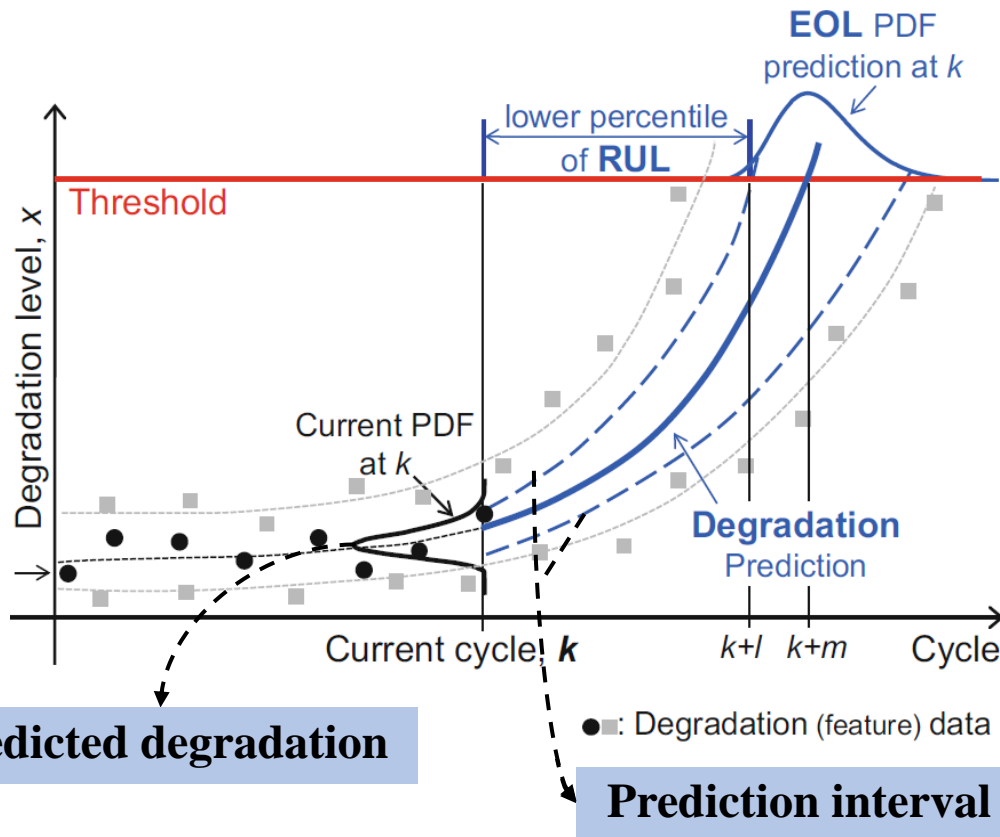
# Prognostics

## Degradation and RUL

# Prognostics

## Objective :

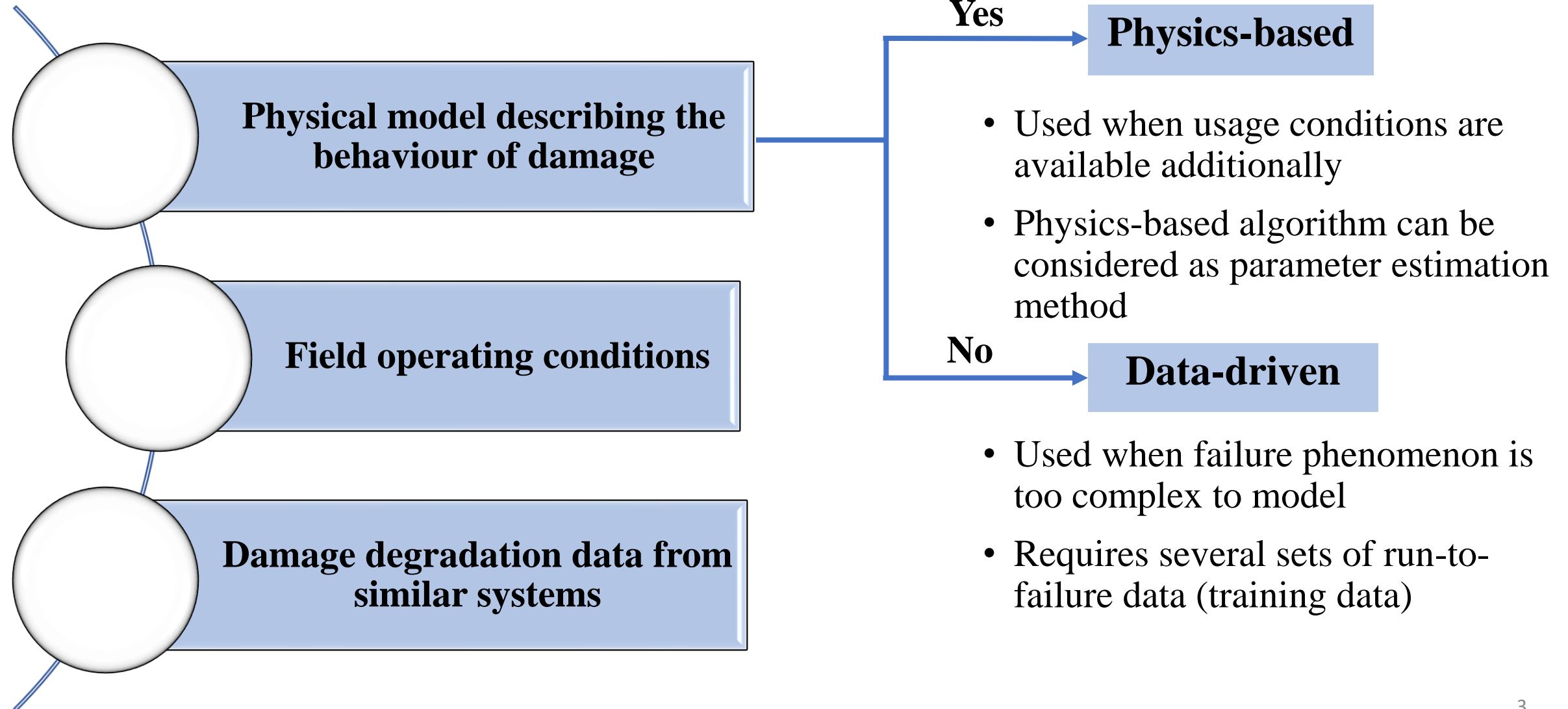
**Predict the remaining useful life (RUL) before the damage grows beyond the threshold**



- Due to uncertainty, predicted degradation and End of Life (EOL) is represented as distribution
- For a given time, the area under the EOL is considered as failure probability
- EOL signals the time for the system maintenance
- RUL, the remaining time to the maintenance from the current time is given by
$$t_{\text{RUL}} = t_{\text{EOL}} - t_k$$
- Lower bound of RUL is considered as the maintenance time

# Prognostics methods

Categorized based on the availability of



# Prediction of degradation behavior

## Least squares method - Parameter estimation

- Used to find unknown parameter or coefficients by minimizing the sum of square errors ( $SS_E$ ) between measured ( $y_k$ ) and simulation ( $z_k$ , from model/function) data
- Their relationship is given by

$$y_k = z_k + \varepsilon_k$$

$k$  is the time index

- The error  $\varepsilon_k$  can represent measurement error in  $y_k$  as well as in  $z_k$
- Assumptions
  - Error comes from measurement only
  - Measurement error does not include bias but unbiased noise
  - Simulation model  $z(t; \theta)$  is a linear function of input variable  $t$  (time) and parameter ( $\theta$ )

$$Z(t; \theta) = \theta_1 + \theta_2 t, \quad \theta = \{\theta_1 \ \theta_2\}^T$$

# Prediction of degradation behavior

## Sum of square errors ( $SS_E$ )

- The pair of input variable and measured degradation at data points is denoted as

$$(t_k, y_k), \quad k = 1, \dots, n_y \text{ and } y = \{y_1 \ y_2 \ \dots y_{n_y}\}^T$$

$n_y$  is measured degradation

- The simulation model can be evaluated at data points and corresponding error is given as

$$z = \begin{Bmatrix} z_1 \\ z_2 \\ \vdots \\ z_{n_y} \end{Bmatrix} = \begin{bmatrix} 1 & t_1 \\ 1 & t_2 \\ \vdots & \vdots \\ 1 & t_{n_y} \end{bmatrix} \begin{Bmatrix} \theta_1 \\ \theta_2 \end{Bmatrix} = X\theta$$

$X$  is the design matrix

$$e = \{\varepsilon_1 \ \varepsilon_2 \ \dots \ \varepsilon_{n_y}\}^T = y - z$$

$$SS_E = e^T e = \{y - z\}^T \{y - z\} = \{y - X\theta\}^T \{y - X\theta\}$$

$$\frac{d(SS_E)}{d\theta} = 2 \left[ \frac{de}{d\theta} \right]^T e = 2X^T \{y - X\theta\} = 0$$

$$\text{The estimated parameter} = \hat{\theta} = [X^T X]^{-1} \{X^T y\}$$

# Data-driven approach

- Functional relationship between input variables and output degradation has to be built from the data
- The quality of prediction depends on
  - Selection of mathematical function
  - Number of data
  - Level of noise in the measurement

## Quality of fitting

- Coefficient of determination,  $R^2$ , the ratio between the variation of function prediction to the data

$$R^2 = \frac{SS_R}{SS_T} = 1 - \frac{SS_E}{SS_T}$$

- $SS_T$  (total sum of squares) =  $\sum_{k=1}^{n_y} (y_k - \bar{y})^2$
- $SS_R$  (regression sum of squares) =  $\sum_{k=1}^{n_y} (z_k - \bar{y})^2$
- $SS_E$  (residual sum of squares) =  $\sum_{k=1}^{n_y} (y_k - z_k)^2$

# Data-driven approach

## Quality of fitting

- $SS_T$  is the variation of data with respect to the mean of data  $\bar{y}$
- $SS_R$  is the variation of the function prediction  $z_k$  with respect to the mean of data
- $SS_E$  is the sum of square of errors remaining after the fit
- When the sum of  $y_k$  is equal to the sum of  $z_k$

$$SS_T = SS_R + SS_E$$

- $R^2$  close to 1 is considered an accurate function
- However,  $R^2$  only measures accuracy in data points which can be unrelated to the true accuracy of the function prediction
- Therefore adjusted  $R^2$  denoted as  $\bar{R}^2$  is used by penalizing the number of coefficients as

$$\bar{R}^2 = 1 - (1 - R^2) \frac{(n_y - 1)}{(n_y - n_p)}$$

$n_y$  and  $n_p$  are the number of data and coefficients respectively

# Data-driven approach

## Overfitting

**When the number of unknown coefficients is larger than the number of data, the least square method tends to fit the noise rather than the trend**

- Overfitting is a modeling error occurring when function is overly complex and when the function has no conformability with the data shape
- Techniques available to avoid overfitting
  - Cross-validation
  - Regularization
  - Early stopping
  - Pruning
- Other approaches
  - Behavior of degradation is expressed with a simple function
  - More training data and usage conditions are used



# RUL prediction

- Remaining time until the degradation grows to a threshold
- The threshold of degradation is determined so that the system is still safe but needs maintenance
- To find RUL, it is necessary to find the time cycle when the level of degradation reached a threshold
- The non-linear equation has to be solved to find the time cycle  $t_{EOL}$

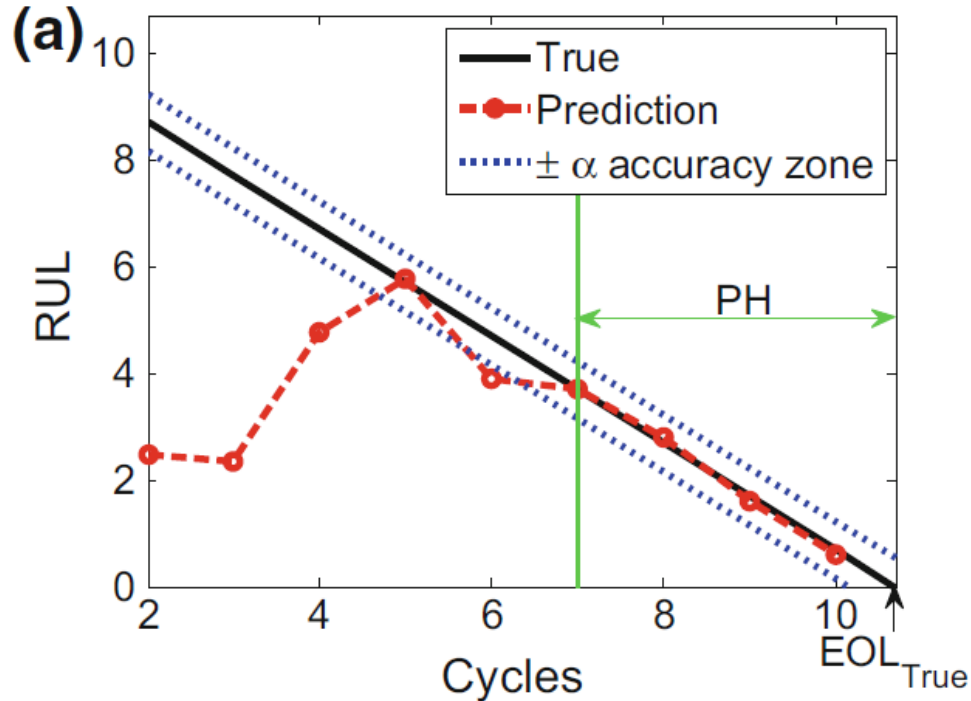
$$y_{threshold} - z(t_{EOL}; \hat{\theta}) = 0$$

- The above equation is solved numerically using Newton-Raphson iterative method
- Maintenance has to be ordered when the RUL becomes zero
- The performance of different method in predicting RUL is compared using metrics such as
  - Prognostic horizon (PH)
  - $\alpha$ - $\lambda$  accuracy
  - Relative accuracy (RA)
  - Cumulative relative accuracy (CRA)
  - Convergence

# Prognostics metrics

## Prognostic Horizon (PH)

Difference between the EOL and the first time when the prediction result continuously resides in the pha accuracy zone

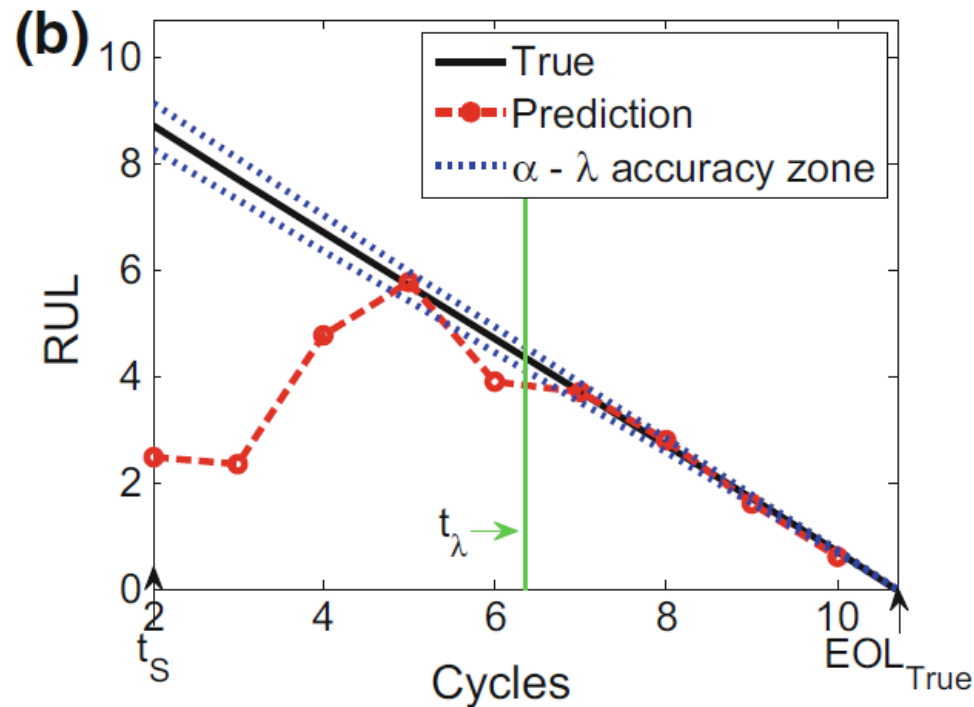


- The two parallel dotted lines indicates the constant bound of the accuracy zone with a magnitude of  $\pm \text{pha}$  error with respect to true EOL.
- Here  $\text{pha} = 5\%$
- The prognostics method with a larger PH indicates the better performance

# Prognostics metrics

## $\alpha$ - $\lambda$ accuracy

Determines whether a prediction result falls within the  $\alpha$  accuracy zone at a specific cycle  $t_\lambda$



- The accuracy zone varies with  $\pm\alpha$  ratio to the true RUL
- The dotted line shows the accuracy zone when  $\alpha = 0.05$
- The accuracy zone shrinks with more data suggesting an increase of prediction accuracy
- Specific cycle  $t_\lambda$  is expressed with a fraction of  $\lambda$  between 0 (starting cycle of RUL prediction) and 1 (true EOL)

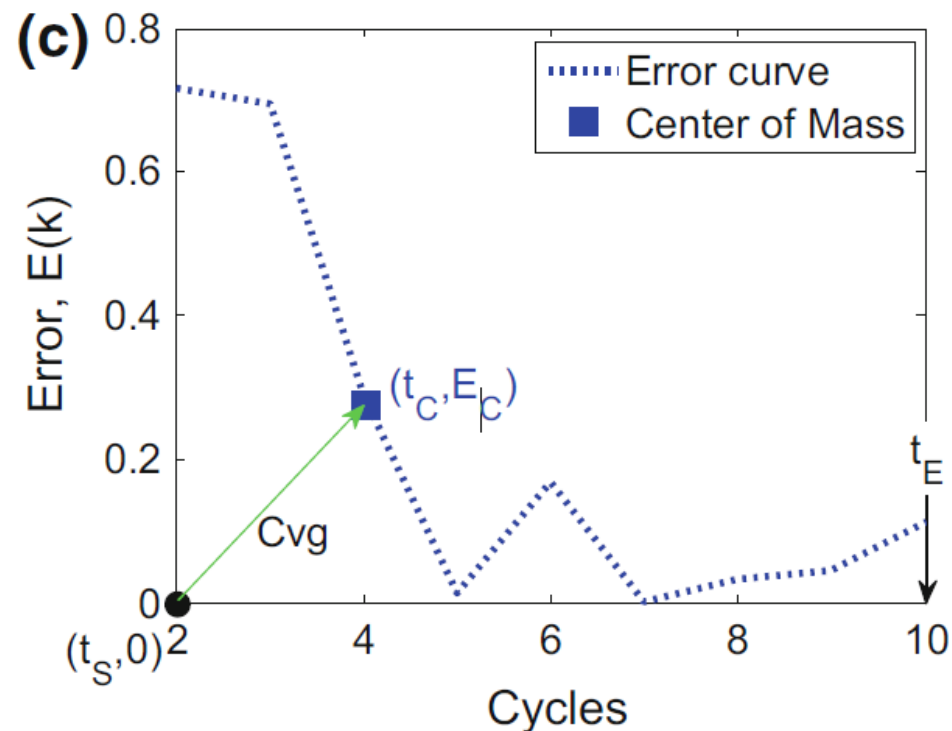
$$t_\lambda = t_s + \lambda(EOL_{True} - t_s)$$

# Prognostics metrics

## (Cumulative) Relative accuracy (RA, CRA)

RA is the relative accuracy between the true and prediction RUL at  $t_\lambda$

CRA is same as the average of RA values accumulated at every cycle from  $t_s$  to  $t_E$



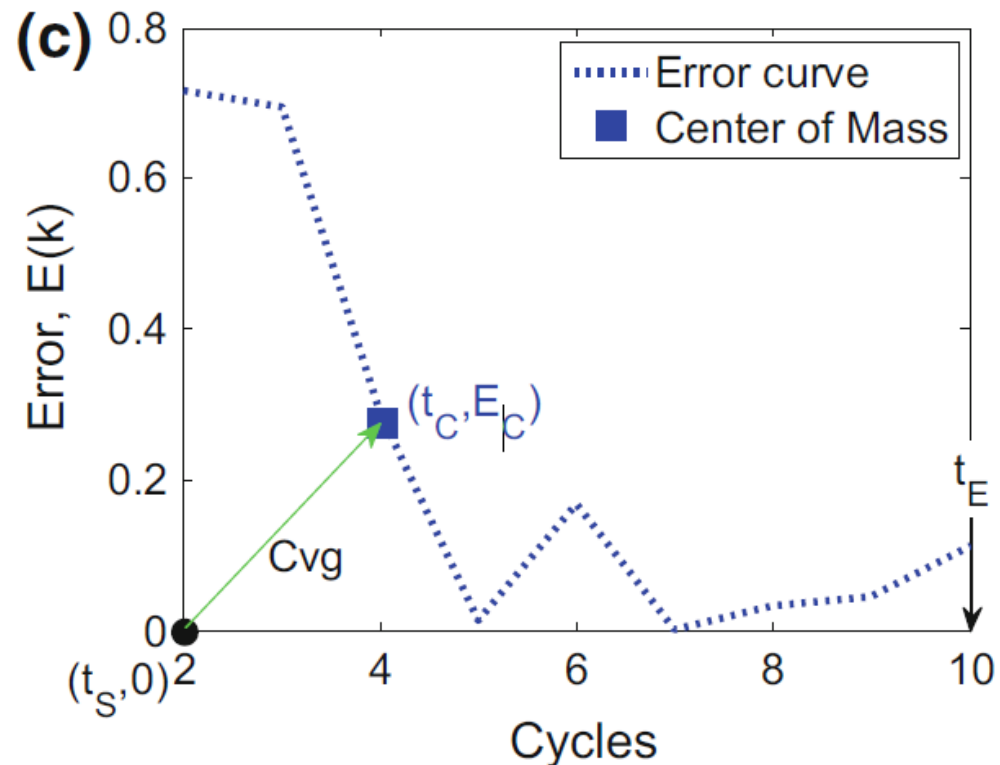
$$RA = 1 - \frac{|RUL_{True} - RUL|}{RUL_{True}} \text{ at } t_\lambda$$

- The relative error shown in the plot (dotted curve) can be used to calculate RA ( $RA = 1 - \text{relative error}$ )
- When RA and CRA are close to 1, prediction accuracy is high

# Prognostics metrics

## Convergence

Euclidean distance between the point  $(t_s, 0)$  and the center of mass of the area under the error curve  $(t_c, E_c)$



$$\text{Center of mass} = CoM = \sqrt{(t_c - t_s)^2 + E_c^2}$$

Where

$$t_c = \frac{1}{2} \frac{\sum_{k=s}^E (t_{k+1}^2 - t_k^2) (E(k))}{\sum_{k=s}^E (t_{k+1} - t_k) E(k)}$$

$$E_c = \frac{1}{2} \frac{\sum_{k=s}^E (t_{k+1} - t_k) (E^2(k))}{\sum_{k=s}^E (t_{k+1} - t_k) E(k)}$$

- The lower the distance is, the faster the convergence is

# Uncertainty

- Noise in measured data is caused by measurement variability, which represent uncertain measurement environment
- Noise is random in nature and thus it is assumed to be Gaussian, *i.e.*, statistical noise having a probability density function equal to that of the normal distribution

**Objective: Estimate the level of uncertainty in estimated model parameters and remaining useful life when measured data have noise**

## Quantification using LS regression

- Assumption:
  - Noise is normally distributed with zero mean
  - They are independent and identically distributed (i.i.d)
- Variance of error is given by

$$\hat{\sigma}^2 = \frac{SS_E}{n_y - n_p}$$

$(n_y - n_p)$  represents the degree of freedom.  $n_y$  and  $n_p$  are the number of data and unknown parameters respectively

# Uncertainty (Model parameters)

## Quantification using LS regression

- For the linear model with 2 parameters ( $z(x) = \theta_1 + \theta_2 x$ ), the variance of parameters is derived using the estimated parameter and the variance of error equations

$$\text{Design matrix} = \mathbf{X} = \begin{bmatrix} 1 & x_1 \\ 1 & x_2 \\ \vdots & \vdots \\ 1 & x_{n_y} \end{bmatrix}, \quad \text{Therefore, } \mathbf{X}^T \mathbf{X} = \begin{bmatrix} n_y & \sum_{i=1}^{n_y} x_i \\ \sum_{i=1}^{n_y} x_i & \sum_{i=1}^{n_y} x_i^2 \end{bmatrix}, \quad \mathbf{X}^T \mathbf{y} = \begin{bmatrix} \sum_{i=1}^{n_y} y_i \\ \sum_{i=1}^{n_y} x_i y_i \end{bmatrix}$$

$$\hat{\boldsymbol{\theta}} = \begin{Bmatrix} \hat{\theta}_1 \\ \hat{\theta}_2 \end{Bmatrix} = \begin{bmatrix} n_y & \sum_{i=1}^{n_y} x_i \\ \sum_{i=1}^{n_y} x_i & \sum_{i=1}^{n_y} x_i^2 \end{bmatrix}^{-1} \begin{bmatrix} \sum_{i=1}^{n_y} y_i \\ \sum_{i=1}^{n_y} x_i y_i \end{bmatrix} = \begin{Bmatrix} \bar{y} - \hat{\theta}_2 \bar{x} \\ \frac{\sum_{i=1}^{n_y} (x_i - \bar{x})(y_i - \bar{y})}{\sum_{i=1}^{n_y} (x_i - \bar{x})^2} \end{Bmatrix}$$

$$\hat{\boldsymbol{\theta}} = \begin{Bmatrix} \hat{\theta}_1 \\ \hat{\theta}_2 \end{Bmatrix} = \begin{Bmatrix} \bar{y} - \bar{x} S_{xy}/S_{xx} \\ S_{xy}/S_{xx} \end{Bmatrix}$$

# Uncertainty (Model parameters)

## Quantification using LS regression

- By using a theorem for variance calculation, the variance of the two parameters is obtained as

$$\text{Var}(\hat{\theta}_2) = \text{Var}\left(\frac{\sum(x_i - \bar{x})(y_i - \bar{y})}{S_{xx}}\right) = \frac{\sum(x_i - \bar{x})^2 \sigma^2}{S_{xx}^2} = \frac{S_{xx} \sigma^2}{S_{xx}^2} = \frac{\sigma^2}{S_{xx}}$$

$$\text{Var}(\hat{\theta}_1) = \text{Var}(\bar{y} - \hat{\theta}_2 \bar{x}) = \frac{\sigma^2}{n_y} + \bar{x}^2 \frac{\sigma^2}{S_{xx}}$$

- The terms related with x are constants because the source of uncertainty is from the noise in the data and the terms related with data  $y_i$  are random variables
- The variance of two parameters in a linear model is

$$\text{Var}(\hat{\theta}) = \begin{bmatrix} \frac{\sigma^2}{n_y} + \bar{x}^2 \frac{\sigma^2}{S_{xx}} \\ \frac{\sigma^2}{S_{xx}} \end{bmatrix}, \quad S_{xx} = \sum_{i=1}^{n_y} (x_i - \bar{x})^2$$



# Uncertainty (Model parameters)

## Quantification using LS regression

- The variance in model parameters is linearly proportional to the variance of data
- Large number of data reduces uncertainty in model parameters and eventually makes them deterministic ( $n_y \rightarrow \infty$ ,  $S_{xx} \rightarrow \infty$ )
- In case of many unknown parameters, the derivation is complicated and the following can be employed to describe the correlation between the parameters

$$\Sigma_{\hat{\theta}} = \sigma^2 [X^T X]^{-1}$$

- Since the variance of error in data is unknown, usually estimated value can be used instead

# Issues in practical prognostics

- In practical cases, simple polynomial functions and Gaussian noise does not suffice
- Bayesian-based approaches are employed instead of linear regression method
- The noise and bias have an effect on the prognostics results
- It is difficult to identify model parameters accurately when they are correlated
- Loading conditions can be correlated with the parameters too
- Physical degradation models are rare in practice
- In case of data-driven approach, it is not easy to obtain several sets of degradation data due to expensive time and costs
- Degradation data cannot be directly measured in most case and need to be extracted from sensor signals

