

# Imprecise Software Reliability

Lev Utkin

Munich, September 2009

# Standard definitions of software reliability

- A **fault** in a software is an incorrect step, process, or data definition.
- A fault may cause a **failure**, or “the inability of a system or component to perform its required functions within specified performance requirements”; that is, a deviation from the stated or implied requirements.
- An **error** is a “discrepancy between a computed, observed, or measured value or condition and the true, specified, or theoretically correct value or condition.

# Two types of software models

There are two main software reliability models:

## ① Reliability growth model

- *statistical data are obtained during debugging process under condition that a detected error is removed or corrected.*

## ② Reliability growth model

- *statistical data are obtained during testing under condition that a detected error is not removed or corrected.*

# Standard initial information concerning software reliability

Three main kinds of reliability statistical data.

- 1 Calendar times (reals) between software failures  
 $\mathbf{T} = (t_1, \dots, t_n)$ .
- 2 Numbers of successful runs (a run is minimum execution unit of software, integers) between software failures  
 $\mathbf{K} = (k_1, \dots, k_n)$ .
- 3 Numbers of failures in certain periods of time  
 $(k_1, t_1), \dots, (k_n, t_n)$ .

# Main goals of modelling

*To predict the reliability measures or indices of a analyzed software after:*

- 1 **Debugging process** (growth models)
- 2 **Testing** (testing models)

In other words, we have to compute the probabilistic measures of the r.v.  $X_{n+1}$  (time or number of runs to failure, number of failures after  $n$  observations).

# Three types of models

- 1 Calendar times: continuous models (continuous random variables)
- 2 Numbers of successful runs: discrete models (discrete random variables).
- 3 Numbers of failures in certain periods of time: non-homogeneous Poisson process models.

# Assumptions in most well-known reliability growth models

- 1 The operational profile of the software remains constant, i.e., the software where the data comes from is operated in a similar manner as that in which reliability predictions are to be made.
- 2 Every model assume some precise probability distribution of random variables under consideration.
- 3 The failures, when the faults are detected, are independent, i.e., random variables under consideration are statistically independent.
- 4 Every fault has an equal chance of being encountered within a severity class as any other fault in that class.
- 5 After a failure, the fault causing it is corrected immediately and no new faults result from the correction.

# Assumptions in most well-known reliability testing models

- 1 The software doesn't change during testing and usage, except that faults are fixed.
- 2 The operational profile of the software remains constant, i.e., the software where the data comes from is operated in a similar manner as that in which reliability predictions are to be made.
- 3 Every model assume some precise probability distribution of random variables under consideration.
- 4 The failures, when the faults are detected, are independent, i.e., random variables under consideration are statistically independent.



# Assumptions in most well-known reliability models

*The assumptions are usually not fulfilled. As a result, the reliability may be too unreliable and risky.*

## A standard way for dealing the models

- 1 Let  $X_i$  be a random time interval between the  $(i - 1)$ -st and  $i$ -th software failures. The variable  $X_i$  is governed by a probability density function  $p_i(x|\theta)$  with a vector of parameters  $\theta_i$ . It is assumed that there holds  $\theta_i = f(i, \theta)$ .

# A standard way for dealing the models

- 1 Let  $X_i$  be a random time interval between the  $(i - 1)$ -st and  $i$ -th software failures. The variable  $X_i$  is governed by a probability density function  $p_i(x|\theta)$  with a vector of parameters  $\theta_i$ . It is assumed that there holds  $\theta_i = f(i, \theta)$ .
- 2 Let  $\mathbf{X} = (x_1, \dots, x_n)$  be the successive intervals between failures. The likelihood function is

$$L(\mathbf{X}|\theta) = \Pr\{X_1 = x_1, \dots, X_n = x_n|\theta\} = \prod_{i=1}^n p_i(x_i|\theta).$$

# A standard way for dealing the models

- 1 Let  $X_i$  be a random time interval between the  $(i - 1)$ -st and  $i$ -th software failures. The variable  $X_i$  is governed by a probability density function  $p_i(x|\theta)$  with a vector of parameters  $\theta_i$ . It is assumed that there holds  $\theta_i = f(i, \theta)$ .
- 2 Let  $\mathbf{X} = (x_1, \dots, x_n)$  be the successive intervals between failures. The likelihood function is

$$L(\mathbf{X}|\theta) = \Pr\{X_1 = x_1, \dots, X_n = x_n|\theta\} = \prod_{i=1}^n p_i(x_i|\theta).$$

- 3  $\theta_{opt} = \arg \max_{\theta} L(\mathbf{X}|\theta)$

# A standard way for dealing the models

- 1 Let  $X_i$  be a random time interval between the  $(i - 1)$ -st and  $i$ -th software failures. The variable  $X_i$  is governed by a probability density function  $p_i(x|\theta)$  with a vector of parameters  $\theta_i$ . It is assumed that there holds  $\theta_i = f(i, \theta)$ .
- 2 Let  $\mathbf{X} = (x_1, \dots, x_n)$  be the successive intervals between failures. The likelihood function is

$$L(\mathbf{X}|\theta) = \Pr\{X_1 = x_1, \dots, X_n = x_n|\theta\} = \prod_{i=1}^n p_i(x_i|\theta).$$

- 3  $\theta_{opt} = \arg \max_{\theta} L(\mathbf{X}|\theta)$
- 4 The software failure function after the  $n$ -th software failure is computed as follows:

$$F_{n+1}(t) = \int_0^t p_{n+1}(x|\theta_{opt}) dx.$$

# Jelinski-Moranda model

- 1 The initial number of faults in the software is  $N$ .
- 2 The time between failures follows an exponential distribution with a parameter that is proportional to the number of remaining faults in the software.
- 3 The mean time between the  $(i - 1)$ -st and  $i$ -th failures is  $1/\lambda(N - (i - 1))$ .

4

$$L(\mathbf{X}|\lambda, N) = \prod_{i=1}^n \lambda(N - (i - 1)) \exp(-\lambda(N - (i - 1))t) \rightarrow \max_{\lambda, N} .$$

5

$$F_{n+1}(t) = \exp(-\lambda(N - n)t) .$$

# Schick-Wolverton and Littliwood-Verrall models

- 1 **Schick-Wolverton model:** the time between failures is governed by the Rayleigh distribution with the pdf

$$f_i(t) = t\lambda \cdot \exp(-\lambda_i t^2/2).$$

- 2 **Littliwood-Verrall model:** the time between failures is governed by the Pareto distributions with the pdf

$$f_i(t) = \frac{\sigma (\psi(i))^\sigma}{[t + \psi(i)]^{\sigma+1}},$$

$$\psi(i) = \beta_0 + \beta_1 i, \quad \psi(i) = \beta_0 + \beta_1 i^2.$$

# Non-homogeneous Poisson process (NHPP) models

For any time points  $0 < t_1 < t_2 < \dots$ , the probability that the number of failures between  $t_{i-1}$  and  $t_i$  is  $k$  can be written as

$$\begin{aligned} & \Pr \{N(t_i) - N(t_{i-1}) = k\} \\ &= \frac{\{m(t_i) - m(t_{i-1})\}^k}{k!} \exp \{- (m(t_i) - m(t_{i-1}))\}. \end{aligned}$$

Here  $m(t)$  is the mean number of failures occurring up to time  $t$ . The NHPP models differ by a mean value function  $m(t)$ .



## Specific NHPP models

- $m(t) = at^b$  (Duan model),
- $m(t) = a(1 - (1 - b)^t)$  (Yamada-Ohba-Osaki model),
- $m(t) = a(1 - \exp(-bt))$  (Goel-Okumoto model),
- $m(t) = a \ln(1 + bt)$  (Musa-Okumoto model).

## General class of NHPP models

It has been shown by Pham *et al* (Pham, Nordmann, Zhang 1999) that a general class of NHPP models can be obtained by solving the differential equation

$$\frac{dm(t)}{dt} = b(t) [a(t) - m(t)]$$

with suitably chosen  $a(t)$  and  $b(t)$ .

The parameters can be estimated using the maximum likelihood method based on the number of failures per interval of testing and debugging.

# Fuzzy software reliability models (Cai's models)

- The time between the  $(i - 1)$ -st and  $i$ -th failures  $X_i$  is a fuzzy variable governed by a membership function  $\mu_i(x)$ , for example  $\mu_i(x) = \exp(-(x - a_i)^2)$  (Kai-Yuan Cai et al 1991,1993).
- $a_i = f(i, \theta)$ , for example,  $f(i) = (A + Bi)^{-\alpha} + C$ ,  $A, B, C, \alpha \in \theta$ .
- The possibilistic likelihood function is

$$\begin{aligned} L(x_1, \dots, x_n | \theta) &= \text{Pos} \{X_1 = x_1, \dots, X_n = x_n\} \\ &= \min \{\mu_1(x_1), \dots, \mu_n(x_n)\}. \end{aligned}$$

- The reliability after  $n$ -th software failure is

$$R(t) = \sup_{x \geq t} \mu_{n+1}(x).$$

- The main difficulty is how to interpret  $R(t)$ .

# The first idea: maximum of the likelihood function over the set of CDFs (discrete case)

- Every  $X_i$  is governed by an unknown CDF belonging to a set  $\mathcal{M}_i(\mathbf{d})$  depending on a vector of parameters  $\mathbf{d}$  and defined by **lower and upper CDFs**:

$$\underline{F}_i(k | \mathbf{d}) = \inf_{F(k) \in \mathcal{M}_i(\mathbf{d})} F(k), \quad \bar{F}_i(k | \mathbf{d}) = \sup_{F(k) \in \mathcal{M}_i(\mathbf{d})} F(k).$$

- The likelihood function  $L(\mathbf{K} | \mathbf{d}, F)$  is maximized over all distributions  $F$  from  $\mathcal{M}_i(\mathbf{d})$  and the resulting “modified” likelihood function depends on  $\mathbf{d}$ :

$$L(\mathbf{K} | \mathbf{d}) = \max_{F \in \mathcal{M}_1(\mathbf{d}), \dots, F \in \mathcal{M}_n(\mathbf{d})} L(\mathbf{K} | \mathbf{d}, F).$$

# The maximized likelihood function

## Proposition

*If random variables  $X_1, \dots, X_n$  are independent and discrete, then there holds*

$$\max_{\mathcal{M}} \Pr \{X_1 = k_1, \dots, X_n = k_n\} = \prod_{i=1}^n \{\bar{F}_i(k_i) - \underline{F}_i(k_i - 1)\}.$$

## “Precise” case

### Corollary

If  $\bar{F}_i(k) = \underline{F}_i(k) = F_i(k)$ , then

$$\max_{\mathcal{M}} \Pr \{X_1 = k_1, \dots, X_n = k_n\} = \prod_{i=1}^n p_i(k_i) = L(\mathbf{K} \mid \mathbf{d}).$$

Here  $p_i(k)$  is the probability mass function corresponding to the distribution function  $F_i(k)$ .

We have the standard likelihood function.

# The maximized likelihood function (the lack of independence)

## Proposition

*If there is no information about independence of random variables  $X_1, \dots, X_n$ , then there holds*

$$\max_{\mathcal{M}} \Pr \{X_1 = k_1, \dots, X_n = k_n\} = \min_{i=1, \dots, n} \{\bar{F}_i(k_i) - \underline{F}_i(k_i - 1)\}.$$

## “Precise” case

### Corollary

If  $\bar{F}_i(k) = \underline{F}_i(k)$ , then

$$\max_{\mathcal{M}} \Pr \{X_1 = k_1, \dots, X_n = k_n\} = \min_{i=1, \dots, n} p_i(k_i).$$

We have the possibilistic likelihood function.



## Justification

The likelihood function is the probability

$$L(\mathbf{K} \mid \mathbf{d}) = \Pr \{X_1 = k_1, \dots, X_n = k_n \mid \mathbf{d}\}.$$

Our final goal is to maximize this probability over a set of parameters  $\mathbf{d}$ . But we have a set of probabilities. Therefore, we choose the largest probability in the set, i.e., we maximize the likelihood function over the set of probabilities depending on  $\mathbf{d}$ .

## The second idea: imprecise Bayesian inference

- Every set  $\mathcal{M}_i(\mathbf{d})$  is defined by boundary **lower and upper CDFs**  $\underline{F}_i(k | \mathbf{d})$  and  $\overline{F}_i(k | \mathbf{d})$  which can be determined by using imprecise Bayesian inference.
- Before debugging process: we do not have information about  $\underline{F}_i(k | \mathbf{d})$  and  $\overline{F}_i(k | \mathbf{d})$ . They are 0 and 1.
- After debugging process: we have  $n$  observations and by taking one of the imprecise Bayesian models (corresponding to a specific probability distribution), we determine the lower and upper CDFs depending on  $\mathbf{d}$ .

**Now we can construct SRGMs by taking the corresponding probability distributions for imprecise Bayesian inference.**

## The imprecise beta-binomial model (1)

**Assumption 1:** The  $i$ -th run lifetime of software  $X_j$  is governed by the geometric distribution with parameter  $p_j$ ,  $j = 1, \dots, n$ .

The prior Beta distribution of the random variable  $p$  is:

$$\pi(p) = \text{BetaDen}(a, b) = \frac{1}{\text{Beta}(a, b)} p^{a-1} (1-p)^{b-1}, \quad 0 \leq p \leq 1.$$

Here  $a > 0$ ,  $b > 0$  are parameters.

The posterior beta distribution  $\pi(p|k)$  after  $k$  events under consideration from the total number of  $n$  event:

$$\pi(p|k) = \text{BetaDen}(a + k, b + n - k).$$

## The imprecise beta-binomial model (2)

**Assumption 2:** Introduce the growth function  $\psi(j) = (j - 1)\varphi$  (with parameter  $\varphi$ ) such that parameters of the posterior beta distribution are

$$a^* = a + j - 1, \quad b^* = b + D_j,$$

$$D_j = K_j + \psi(j), \quad K_j = \sum_{i=1}^{j-1} (k_i - 1), \quad \psi(j) = (j - 1)\varphi.$$

The predictive CDF  $F_j(m)$  for the  $j$ -th step of the debugging is

$$F_j(m|\varphi) = 1 - \frac{\Gamma(\alpha^* + \beta^*)}{\Gamma(\beta^*)} \frac{\Gamma(\beta^* + m)}{\Gamma(\alpha^* + \beta^* + m)}.$$

## The imprecise beta-binomial model (3)

Introduce new parameters  $s > 0$  and  $\gamma \in [0, 1]$  such that

$$a = s\gamma, \quad b = s - s\gamma.$$

Then

$$\underline{F}_j^{(s)}(m|\varphi) = 1 - \frac{\text{Beta}(s + j - 1 + D_j, m)}{\text{Beta}(s + D_j, m)},$$

$$\overline{F}_j^{(s)}(m|\varphi) = 1 - \frac{\text{Beta}(s + j - 1 + D_j, m)}{\text{Beta}(D_j, m)}.$$

## The imprecise beta-binomial model (4)

$$\begin{aligned} \varphi_0 &= \arg \max_{\varphi} L^{(s)}(\mathbf{K}|\varphi) \\ &= \prod_{j=1}^n \left( \frac{\text{Beta}(s + j - 1 + D_j, k_j - 1)}{\text{Beta}(s + D_j, k_j - 1)} - \frac{\text{Beta}(s + j - 1 + D_j, k_j)}{\text{Beta}(D_j, k_j)} \right). \end{aligned}$$

$$\underline{F}_{n+1}^{(s)}(m) = 1 - \frac{\text{Beta}(s + n + K_{n+1} + n\varphi_0, m)}{\text{Beta}(s + K_{n+1} + n\varphi_0, m)},$$

$$\overline{F}_{n+1}^{(s)}(m) = 1 - \frac{\text{Beta}(s + n + K_{n+1} + n\varphi_0, m)}{\text{Beta}(K_{n+1} + n\varphi_0, m)}.$$

## The imprecise negative binomial model

The number of failures has a Poisson distribution with the parameter  $\lambda$ . If we observed  $K$  failures during a period of time  $T$ , then the predictive probability of  $k$  failures during time  $t$  under condition that  $K$  failures were observed during time  $T$  is

$$\begin{aligned} P(k, t) &= \int_0^\infty \frac{(\lambda t)^k e^{-\lambda t}}{k!} \text{Gamma}(\alpha^*, \beta^*) d\lambda \\ &= \frac{\Gamma(\alpha^* + k)}{\Gamma(\alpha^*) k!} \left( \frac{\beta^*}{\beta^* + t} \right)^{\alpha^*} \left( \frac{t}{\beta^* + t} \right)^k. \end{aligned}$$

Here  $\alpha^* = \alpha + K$ ,  $\beta^* = \beta + T$ .

# The imprecise negative binomial growth model (1)

- **Assumption:**  $m(t; a, b) = a \cdot \tau(t, b)$  (the parameter  $a$  can be written separately).
- The parameter  $\lambda$  and the argument  $t$  of the Poisson distribution are replaced by the parameter  $a$  and the discrete time  $\tau(t_i, b) - \tau(t_{i-1}, b)$ , respectively.



## The imprecise negative binomial growth model (2)

The predictive CDF of the number of failures in the interval between  $t_i$  and  $t$  ( $t \in [t_i, t_{i+1}]$ ) after  $n$  periods is

$$\begin{aligned} F_i(k, t | \mathbf{c}, b) &= 1 - \frac{B_{q(i,t)}(k+1, \alpha + K_n)}{B(k+1, \alpha + K_n)} \\ &= 1 - I(q(i, t), k+1, \alpha + K_n). \end{aligned}$$

$$\begin{aligned} q(i, t) &= \frac{T_i(t, b)}{\beta + \tau(t_n, b) + T_i(t, b)}, \\ T_i(t, b) &= \tau(t, b) - \tau(t_i, b), \quad K_n = \sum_{j=1}^n k_j, \end{aligned}$$

$B_q(k+1, r)$  is the incomplete Beta-function with  $I(q, k, r)$  the regularized incomplete Beta-function.

## The imprecise negative binomial growth model (3)

- Choose all vectors  $(\alpha, \beta)$  within the triangle  $(0, 0)$ ,  $(s_1, 0)$ ,  $(0, s_2)$ . All possible prior rates of occurrence of failures are represented, as the prior allows interpretation of  $\alpha/\beta = \gamma$  as this rate, hence this would include all such rates in  $(0, \infty)$ .



$$\underline{F}_i(k, t | s_1, s_2, b) = 1 - I \left( \frac{T_i(t, b)}{\tau(t_n, b) + T_i(t, b)}, k + 1, s_1 + K_n \right),$$

$$\overline{F}_i(k, t | s_1, s_2, b) = 1 - I \left( \frac{T_i(t, b)}{s_2 + \tau(t_n, b) + T_i(t, b)}, k + 1, K_n \right).$$

- The likelihood function is

$$L(\mathbf{K} | b, s) = \prod_{i=1}^n (\overline{F}_i(k_i, t_i | s_1, s_2, b) - \underline{F}_i(k_i - 1, t_i | s_1, s_2, b)).$$

## About two caution parameters

- ① The lower bound  $\underline{\mathbb{E}}_i^{(s)} X$  is

$$\underline{\mathbb{E}}_i^{(s)} X = K \frac{t}{s + T}$$

In fact, the parameter  $s$  here increases the time of testing on the value  $s$  (hidden time).

- ② The upper bound  $\overline{\mathbb{E}}_i^{(s)} X$  is

$$\overline{\mathbb{E}}_i^{(s)} X = (s + K) \frac{t}{T}$$

In fact, the parameter  $s$  here increases the number of failures on the value  $s$  (hidden number of failure).

## The imprecise negative binomial growth model (3)

The cumulative probability distribution of the number of failures in time interval  $[t_n, t]$  after  $n$  periods of debugging

$$\underline{F}_{n+1}(k, t | s_1, s_2) = 1 - I \left( \frac{T_n(t, b)}{\tau(t_n, b) + T_n(t, b)}, k + 1, s_1, + K_n \right),$$

$$\overline{F}_{n+1}(k, t | s_1, s_2) = 1 - I \left( \frac{T_n(t, b)}{s_2 + \tau(t_n, b) + T_n(t, b)}, k + 1, K_n \right).$$

# Imprecise modifications of NHPP models

- 1 **Imprecise Bayesian modifications of Musa-Okumoto model:**  $m(t) = a \ln(1 + bt)$ .

For the model:

$$\tau(t, b) = \ln(1 + bt), \quad T_j(t, b) = \frac{\ln(1 + bt)}{\ln(1 + bt_j)}.$$

# Imprecise modifications of NHPP models

- 1 **Imprecise Bayesian modifications of Musa-Okumoto model:**  $m(t) = a \ln(1 + bt)$ .

For the model:

$$\tau(t, b) = \ln(1 + bt), \quad T_j(t, b) = \frac{\ln(1 + bt)}{\ln(1 + bt_j)}.$$

- 2 **Imprecise Bayesian modification of Goal-Okumoto model:**  $m(t) = a(1 - \exp(-bt))$ .

$$\tau(t, b) = 1 - \exp(-bt).$$

# Validation of models

**Algorithm:** We predict the  $(i + 1)$ -st mean time to failure  $\underline{\mathbb{E}}_{i+1}^{(s)} X_{i+1}$ ,  $\overline{\mathbb{E}}_{i+1}^{(s)} X_{i+1}$ ,  $\mathbb{E}_{i+1}^{(s)} X_{i+1}$ , starting from  $i = 3$ . Then we compare these values with the actual times to failure  $k_{i+1}$ .

**Measures of model quality:**

$$R_1 = M^{-1} \cdot \left| \underline{\mathbb{E}}_{i+1}^{(s)} X_{i+1} - k_{i+1} \right|, \quad R_2 = M^{-1} \cdot \left| \overline{\mathbb{E}}_{i+1}^{(s)} X_{i+1} - k_{i+1} \right|,$$

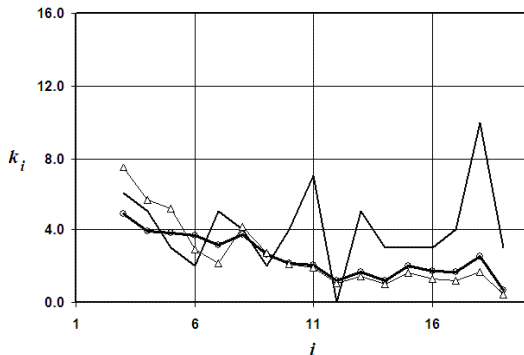
$$R_3 = M^{-1} \cdot \left| \mathbb{E}_{i+1}^{(s)} X_{i+1} - k_{i+1} \right|, \quad R_4 = M^{-1} \cdot \left| \mathbb{E}_{i+1} X_{i+1} - k_{i+1} \right|,$$

where  $M$  is the number of predicted times to failure,

$$\mathbb{E}_{i+1}^{(s)} X_{i+1} = \gamma \underline{\mathbb{E}}_{i+1}^{(s)} X_{i+1} + (1 - \gamma) \overline{\mathbb{E}}_{i+1}^{(s)} X_{i+1}, \quad \gamma = 0.5.$$

# Validation of the imprecise Goal-Okumoto model (1)

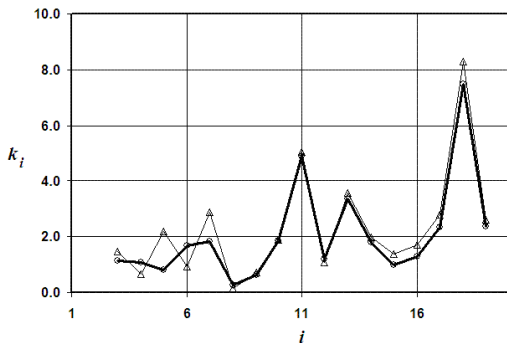
Predicted expected values of numbers of failures for different models based on the Goal-Okumoto model





## Validation of the imprecise Goal-Okumoto model (2)

Deviations of the predicted expected values of numbers of failures from the given values in data sets for the Goal-Okumoto models



## Validation of the imprecise Goal-Okumoto model (3)

- The measures of quality by  $s = 1$  after predicting 17 numbers of failures (from the 3-rd test till 19-th test) by means of the imprecise Bayesian Goel-Okumoto model and the standard Goel-Okumoto model

$$R_1 = 7.491, R_2 = 2.054, R_3 = 1.786,$$

$$R_1^* = 8.311, R_2^* = 2.312, R_3^* = 1.959.$$

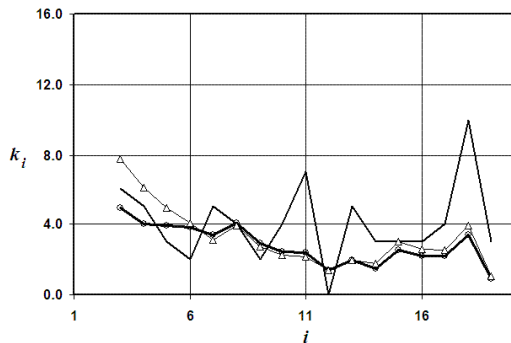
- After predicting 6 numbers of failures (from the 3-rd test till 8-th test)

$$R_1 = 1.827, R_2 = 1.133, R_3 = 0.568,$$

$$R_1^* = 2.854, R_2^* = 1.373, R_3^* = 1.007.$$

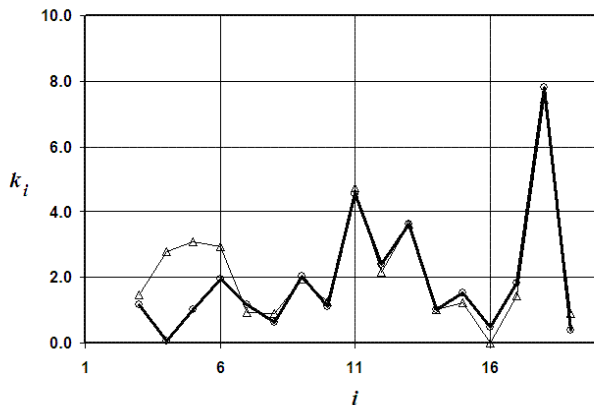
# Validation of the imprecise Musa-Okumoto model (1)

Predicted expected values of numbers of failures for different models based on the Musa-Okumoto model



## Validation of the imprecise Musa-Okumoto model (2)

Deviations of the predicted expected values of numbers of failures from the given values in data sets for the Musa-Okumoto models



## Validation of the imprecise Musa-Okumoto model (3)

- The measures of quality by  $s = 1$  after predicting 17 numbers of failures (from the 3-rd test till 19-th test) by means of the imprecise Bayesian Musa-Okumoto model and the standard Musa-Okumoto model

$$R_1 = 7.817, R_2 = 1.924, R_3 = 1.905,$$

$$R_1^* = 7.445, R_2^* = 2.222, R_3^* = 1.800.$$

- After predicting 6 numbers of failures (from the 3-rd test till 8-th test)

$$R_1 = 1.932, R_2 = 0.990, R_3 = 0.620,$$

$$R_1^* = 3.060, R_2^* = 2.005, R_3^* = 1.020.$$

# Open problems and ideas (1)

- 1 The lack of independence of times to failures:

$$\max_{\mathcal{M}} \Pr \{X_1 = k_1, \dots, X_n = k_n\} = \min_{i=1, \dots, n} \{ \bar{F}_i(k_i) - \underline{F}_i(k_i - 1) \} .$$

**How to realize Bayesian approach in this case?**

# Open problems and ideas (1)

- 1 The lack of independence of times to failures:

$$\max_{\mathcal{M}} \Pr \{X_1 = k_1, \dots, X_n = k_n\} = \min_{i=1, \dots, n} \{ \bar{F}_i(k_i) - \underline{F}_i(k_i - 1) \} .$$

**How to realize Bayesian approach in this case?**

- 2 **The method of generalized moments in place of the imprecise Bayesian approach.** The probability  $\Pr \{X_1 = k_1, \dots, X_n = k_n\}$  can be found by using the natural extension with sample moments (or sample generalized moments) as the initial previsions. The number of moments defines the imprecision like the parameter  $s$ .

## Open problems and ideas (2)

- 3 **Hypothesis:**  $k$  moments produce a set  $\mathcal{M}$  of probability mass functions such that

$$\underline{F}_i(k | \mathbf{d}) = \inf_{F(k) \in \mathcal{M}_i(\mathbf{d})} F(k), \quad \overline{F}_i(k | \mathbf{d}) = \sup_{F(k) \in \mathcal{M}_i(\mathbf{d})} F(k).$$

Suppose that  $\pi^*(k) = \max_{\mathcal{M}} \Pr \{X = k\}$ . Is it true that

$$\pi^*(k) = \overline{F}(k) - \underline{F}(k-1)?$$

**It is true for two first moments!**



# Questions

?