Predictive Modeling and Balance Property through Auto-calibration

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Part 1 - Global and local balances in insurance pricing

Setting

- Y = aggregate claim amount or its frequency component.
- $\boldsymbol{X} = (X_1, X_2, \dots, X_p) = \text{risk or rating factors.}$
- Pure premium= amount collected by the insurer to compensate for the claims, without loss nor profit.
- $\pi(\mathbf{X})$ candidate pure premium.
- $\mu(\mathbf{X}) = \mathsf{E}[\mathbf{Y}|\mathbf{X}]$ true pure premium.

Global balance

- **Global balance** is the very **basic requirement** for a pure premium in insurance.
- Formally, it requires that

$$\mathsf{E}[Y] = \mathsf{E}[\pi(\boldsymbol{X})].$$

• In practice, this means that

$$\frac{1}{n}\sum_{i=1}^{n}Y_{i}\approx\frac{1}{n}\sum_{i=1}^{n}\pi(\boldsymbol{X}_{i}) \text{ for large } n$$

• In words, global balance for $\pi(\cdot)$ ensures that

aggregate loss \approx total pure premium income, for large *n*.

Method of marginal totals (MMT)

- Early 1960s in North America, still reflects actuarial thinking.
- With $x_{i,j} \in \{0,1\}$ for all j, the **pure premium** is written as

$$\pi(\boldsymbol{x}_i) = e_i \gamma_0 \prod_{j=1}^p \gamma_j^{x_{i,j}}$$

where

- e_i = risk exposure for policy *i*
- $\gamma_0 =$ **base premium** $\pi(0)$ per unit of exposure
- γ_j = **relativity** associated with risk factor *j*.
- Considering $\boldsymbol{x}_{i_1} = \boldsymbol{x}_{i_2}$ except for $x_{i_1,j} = 1$, $x_{i_2,j} = 0$,

$$\gamma_j = \frac{\pi(\boldsymbol{x}_{i_1})}{\pi(\boldsymbol{x}_{i_2})}.$$

Method of marginal totals (MMT)

- MMT imposes
 - global balance

$$\sum_{i=1}^n y_i = \sum_{i=1}^n \pi(\boldsymbol{x}_i)$$

- and **local balance** over all contracts with $x_{i,j} = 1$

$$\sum_{i|\mathbf{x}_{i,j}=1} y_i = \sum_{i|\mathbf{x}_{i,j}=1} \pi(\mathbf{x}_i) \text{ for } j = 1, 2, \dots, p.$$

- Referring to contingency tables, this type of local balance is called **marginal balance**.
- Base premium γ_0 and relativities γ_j solve the system of balance equations.

From MMT to GLMs

- Assume that given **X**, the response Y obeys a distribution from the **ED family**: Binomial, Poisson, Gamma, Tweedie, etc.
- The pure premium rate for policy *i* is of the form

$$g^{-1}\left(\beta_0+\sum_{j=1}^p\beta_j x_{i,j}\right)$$

where g is the link function.

• With canonical link function

Minimizing deviance \Leftrightarrow Solving MMT equations.

⇒ GLM estimation with canonical link is equivalent to the actuarial MMT which predates GLMs.

 \Rightarrow MMT can be implemented with any tool developed for GLMs.

From MMT to GLMs

• If $x_{i,j} \in \{0,1\}$ for all j and $g = \log$ then

$$\pi(\boldsymbol{x}_i) = \exp\left(\ln e_i + \beta_0 + \sum_{j=1}^p \beta_j x_{i,j}\right) = e_i \exp(\beta_0) \prod_{j \mid x_{i,j}=1} \exp(\beta_j)$$

where

 $exp(\beta_0) = base premium per unit of exposure for the reference$ $class, for which <math>x_{i,1} = x_{i,2} = \ldots = x_{i,p} = 0$ $= \gamma_0$

$$exp(\beta_j)$$
 = relativity of risk factor *j*, i.e. effect of switching
from $x_{i,j} = 0$ to $x_{i,j} = 1$
= γ_j .

• We then **recover the multiplicative structure** of the premium.

Modern learning tools

- Linear score $\beta_0 + \sum_{j=1}^{p} \beta_j x_{i,j}$ in GLMs replaced with much more flexible ones.
- Premiums still determined by **minimizing deviance**.
- However, deviance now
 - **penalizes departures** of $\pi(\mathbf{x}_i)$ from y_i
 - maximizes correlation between $\pi(\mathbf{x}_i)$ and y_i

measured on different scales.

- Documented in Denuit, Sznajder and Trufin (2019), Denuit, Charpentier and Trufin (2021).
- Balance is no more imposed and optimizing deviance may favor correlation.
- Lack of balance is not **problematic** in many applications (like credit scoring or fraud detection) but well **in pure premium calculation**.

Local balance by autocalibration

• Autocalibration by Krüger and Ziegel (2021) imposes

$$\mathsf{E}[Y|\pi(X) = s] = s$$
 for all s .

- Balance thus operates within each group of contracts charged the same premium, preventing any transfer.
- **Example**: two values for $\pi(\mathbf{X})$: π_1 and π_2 . Then, $\pi(\mathbf{X})$ autocalibrated means that

$$\frac{1}{n_1}\sum_{i:\pi(\boldsymbol{X}_i)=\pi_1}Y_i=\pi_1$$
 and $\frac{1}{n_2}\sum_{i:\pi(\boldsymbol{X}_i)=\pi_2}Y_i=\pi_2.$

where $n_j = \sum_{i=1}^n I[\pi(X_i) = \pi_j], j = 1, 2.$

Part 2 - Testing for local balances using the concept of auto-calibration

Lorenz and concentration curves

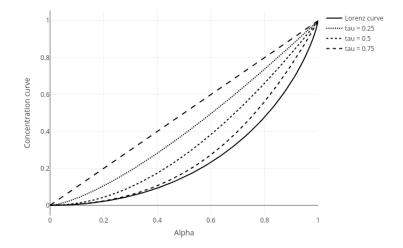
• The Lorenz curve associated to the predictor $\widehat{\pi}$ is defined by

$$\operatorname{LC}[\widehat{\pi}(\boldsymbol{X});\alpha] = \frac{\mathsf{E}\big[\widehat{\pi}(\boldsymbol{X})\mathsf{I}[\widehat{\pi}(\boldsymbol{X}) \leq F_{\widehat{\pi}}^{-1}(\alpha)]\big]}{\mathsf{E}[\widehat{\pi}(\boldsymbol{X})]}, \quad \alpha \in (0,1).$$

• The concentration curve of the response Y with respect to the predictor $\widehat{\pi}$ is defined as

$$\operatorname{CC}[\boldsymbol{Y},\widehat{\boldsymbol{\pi}}(\boldsymbol{X});\alpha] = \frac{\mathsf{E}\big[\boldsymbol{Y}\mathsf{I}[\widehat{\boldsymbol{\pi}}(\boldsymbol{X}) \leq F_{\widehat{\boldsymbol{\pi}}}^{-1}(\alpha)]\big]}{\mathsf{E}[\boldsymbol{Y}]}, \quad \alpha \in (0,1).$$

Lorenz and concentration curves



A new characterization of auto-calibration

• **Proposition:** (Denuit et al., 2024) We have

 $\operatorname{CC}[\mu(\boldsymbol{X}), \widehat{\pi}(\boldsymbol{X}); \alpha] = \operatorname{LC}[\widehat{\pi}(\boldsymbol{X}); \alpha] \quad \text{ for all } \alpha \in (0, 1)$

if, and only if $\widehat{\pi}_{\text{unbiased}}(\mathbf{X}) = \frac{\mathsf{E}[Y]}{\mathsf{E}[\widehat{\pi}(\mathbf{X})]}\widehat{\pi}(\mathbf{X})$ is auto-calibrated.

Statistical test for auto-calibration

- Denuit et al. (2024) introduced a statistical test for auto-calibration.
- The authors test the null hypothesis

$$\mathcal{H}_0: \mathrm{CC}[\mu(\boldsymbol{X}), \widehat{\pi}(\boldsymbol{X}); lpha] = \mathrm{LC}[\widehat{\pi}(\boldsymbol{X}); lpha]$$
 for all $lpha \in (0, 1)$

against the alternative

 $\mathcal{H}_1 : \mathrm{CC}[\mu(\boldsymbol{X}), \widehat{\pi}(\boldsymbol{X}); \alpha] \neq \mathrm{LC}[\widehat{\pi}(\boldsymbol{X}); \alpha] \text{ for some } \alpha \in (0, 1).$

Testing procedure: based on the difference between sample versions

$$\widehat{\mathrm{CC}}[\mu(\boldsymbol{X}),\widehat{\pi}(\boldsymbol{X});\alpha] = \frac{1}{n\bar{Y}}\sum_{i=1}^{n}Y_{i} \mathsf{I}\left[\widehat{\pi}(\boldsymbol{X}_{i}) \leq F_{\widehat{\pi}}^{-1}(\alpha)\right], \ \alpha \in (0,1),$$

and

$$\widehat{\mathrm{LC}}[\widehat{\pi}(\boldsymbol{X}); lpha] = rac{1}{n ar{\pi}} \sum_{i=1}^n \widehat{\pi}(\boldsymbol{X}_i) \mid \left[\widehat{\pi}(\boldsymbol{X}_i) \leq F_{\widehat{\pi}}^{-1}(lpha)
ight], \ \ lpha \in (0, 1).$$

The null hypothesis is rejected for large values of the test statistic

$$\mathcal{T} = \sup_{\alpha \in (0,1)} |T_n(\alpha)|,$$

where

$$\begin{aligned} T_n(\alpha) &= \sqrt{n} \left(\widehat{\mathrm{CC}}[\mu(\boldsymbol{X}), \widehat{\pi}(\boldsymbol{X}); \alpha] - \widehat{\mathrm{LC}}[\widehat{\pi}(\boldsymbol{X}); \alpha] \right) \\ &= n^{-1/2} \sum_{i=1}^n \left(\frac{Y_i}{\bar{Y}} - \frac{\widehat{\pi}(\boldsymbol{X}_i)}{\bar{\pi}} \right) \mathsf{I}[\widehat{\pi}(\boldsymbol{X}_i) \leq F_{\widehat{\pi}}^{-1}(\alpha)], \quad \alpha \in (0, 1). \end{aligned}$$

Proposition: Assume that (Y_i, π̂(X_i)), i = 1, 2, ..., n, are such that E[Y_i] ≠ 0, E[π̂(X_i)] ≠ 0, E[π̂²(X_i)] < ∞ and E[Y_i²] < ∞. Then, under the null hypothesis, T_n(α) converges weakly to a Gaussian process.

The proposed test rejects the null hypothesis when

$$\mathcal{T} = \sup_{\alpha \in (0,1)} |T_n(\alpha)| > c_\beta,$$

for some critical value c_{β} such that $P[\sup_{\alpha \in (0,1)} | \mathcal{T}_n(\alpha)| > c_{\beta}] = \beta.$

However, the analyst cannot compute the critical value c_β since the underlying distribution of (Y_i, π̂(X_i)) is unknown. Given the asymptotic Gaussian behavior of the process, the non-parametric Monte-Carlo methods of Zhu et al. (2016) can be used.

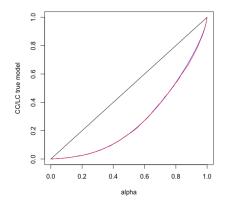
Case study - Data set

- Swiss motor insurance database with 500 000 insurance policies used in Wüthrich and Buser (2016) and in Wüthrich (2020).
- For each policy i :
 - Numbers of claims Y_i;
 - **Exposure-to-risk** $e_i \leq 1$ (i.e. the duration of observation expressed in years)
 - Features $X_i = (X_{i1}, ..., X_{i8})$:
 - X_{i1} = the age of the driver (age);
 - X_{i2} = the age of the car (ac);
 - X_{i3} = the power of the car (power);
 - X_{i4} = the fuel type of the car (fuel);
 - X_{i5} = the vehicle brand (vehicle brand);
 - X_{i6} = the area code of the living place of the driver (area);
 - X_{i7} = the population density at the living place of the driver (dens);
 - X_{i8} = the Swiss canton of the license plate of the car (ct).

- Important: The values e_iµ(X_i) of the true model are also provided (the Y_i have been generated from expected true model frequencies e_iµ(X_i)).
- We partition the data set into a training set D (80%) and a validation set D (20%).

Case study - Data set

Estimates of the CC (in red) and LC (in blue) on D for the true model:



▶ Applying our testing procedure to the true model on D
, we get p
_{equal} = 0.29, meaning that the null hypothesis is not rejected at the level 5%.

Case study - Models under consideration

- Given X = x and the exposure-to-risk e, Y is assumed to be **Poisson with mean** $e\mu(x)$.
- ▶ D is divided into D₁ and D₂, where D₁ includes 80% of the observations of D and D₂ gathers the remaining 20%.
- We first fit **two GAMs** on \mathcal{D}_1 using the R package mgcv:
 - $\hat{\pi}^{\text{GAM1}}(\boldsymbol{x})$, with only the feature X_1 (age);
 - $\hat{\pi}^{\text{GAM2}}(\mathbf{x})$, using all 8 available features.

The covariates age, ac, power and dens are captured by splines and we do not consider interaction terms.

Case study - Models under consideration

• We fit **four GBTs** on \mathcal{D}_1 producing the following estimators:

- $\hat{\pi}^{\text{GBT1}}(\mathbf{x})$, with only the feature X_1 (age);
- $\hat{\pi}^{\text{GBT2}}(\mathbf{x})$, using all **8 available features** and ID = 1;
- $\hat{\pi}^{\text{GBT3}}(\mathbf{x})$, using all 8 available features and ID = 2;
- $\hat{\pi}^{\text{GBT4}}(\mathbf{x})$, using all 8 available features and ID = 3.

The optimal values for the number of trees T are determined by minimizing the out-of-sample Poisson deviance loss computed on \mathcal{D}_2 :

•
$$T = 507$$
 for $\hat{\pi}^{\text{GBT1}}$;

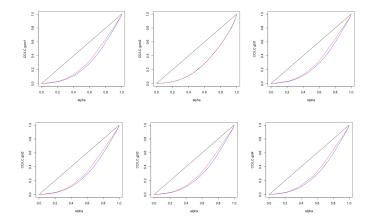
•
$$T = 2977$$
 for $\hat{\pi}^{\text{GBT2}}$;

•
$$T = 2984$$
 for $\hat{\pi}^{\text{GBT3}}$;

▶
$$T = 2943$$
 for $\hat{\pi}^{\text{GBT4}}$.

Case study - Testing for equality of Concentration and Lorenz curves

• Estimates of the CC (in red) and LC (in blue) on $\overline{\mathcal{D}}$:



Case study - Testing for equality of Concentration and Lorenz curves

► Values of \hat{p}_{equal} :

$\widehat{\pi}$	\widehat{p}_{equal}
$\widehat{\pi}^{GAM1}$	0.000
$\widehat{\pi}^{GAM2}$	0.744
$\widehat{\pi}^{GBT1}$	0.000
$\widehat{\pi}^{GBT2}$	0.000
$\widehat{\pi}^{GBT3}$	0.000
$\widehat{\pi}^{GBT4}$	0.000

Part 3 - Balance correction as a means for auto-calibration

Balance correction as a way to restore financial equilibrium

Start from any predictor π(X) strongly correlated with the response Y.

 \Rightarrow The **ranks** produced by $\pi(\cdot)$ **are informative** to order contracts from the cheapest to the most expensive one.

- An accurate premium can then be obtained by averaging claim amounts over neighborhoods induced by $\pi(\cdot)$.
- The balance-corrected premium $\pi_{\rm bc}(\pmb{X})$ obtained from $\pi(\pmb{X})$ by

$$\pi_{\mathsf{bc}}(\boldsymbol{X}) = \mathsf{E}\big[Y\big|\pi(\boldsymbol{X})\big]$$

satisfies autocalibration.

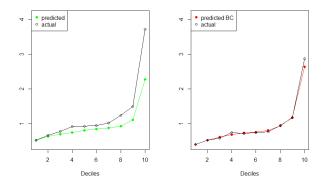
• In practice, balance correction can be implemented by

- local regression (Denuit, Charpentier and Trufin, 2021, Ciatto et al. 2023);

- isotonic regression (Wüthrich and Ziegel, 2023).

Balance correction: Example (Ciatto et al. (2023)

- freMTPL2freq data set (in the CASdatasets package in R).
- 678 013 observations of the number of claims (response Y) in a French MTPL portfolio, with 9 features $\boldsymbol{X} = (X_1, \dots, X_9)$.
- $\hat{\pi}$: boosted Poisson model (training set = 60% of the data).
- $\hat{\pi}_{bc}$: obtained by local regressions (on 20% of the data= validation set).
- Lift chart (on the test set = 20% of the data) :



Impact of balance correction on Bregman divergence

- For a convex function $\ell(\cdot),$ define the error measure for the mean as

$$L(y,m) = \ell(y) - \ell(m) - \ell'(m)(y-m).$$

Functions $L(\cdot, \cdot)$ are called **Bregman loss functions**.

- Bregman loss functions are **important in insurance ratemaking** since any Bregman loss function is a strictly consistent loss function for the mean functional.
- Then,

$$\mathsf{E}\big[L\big(Y,\widehat{\pi}_{\mathrm{bc}}(\boldsymbol{X})\big)\big] \leq \mathsf{E}\big[L\big(Y,\widehat{\pi}(\boldsymbol{X})\big)\big].$$

Impact of balance correction on concentration curves

• **Proposition:** (Denuit and Trufin, 2024) We have

$$\operatorname{CC}[\boldsymbol{Y}, \widehat{\pi}_{\mathrm{bc}}(\boldsymbol{X}); \alpha] \leq \operatorname{CC}[\boldsymbol{Y}, \widehat{\pi}(\boldsymbol{X}); \alpha]$$

for any probability level α .

Impact of balance correction on Lorenz curves

• **Proposition:** (Denuit and Trufin, 2024) Let $t \to m(t) = E[Y|\hat{\pi}(X) = t]$. Assume that $m(\cdot)$ is continuous and strictly increasing. Then,

$$t \mapsto \frac{m(t)}{t} \text{ non-decreasing} \Rightarrow \operatorname{LC}[\widehat{\pi}_{\operatorname{bc}}(\boldsymbol{X}); \alpha] \leq \operatorname{LC}[\widehat{\pi}(\boldsymbol{X}); \alpha]$$

while

$$t \mapsto \frac{m(t)}{t} \text{ non-increasing} \Rightarrow \operatorname{LC}[\widehat{\pi}_{\operatorname{bc}}(\boldsymbol{X}); \alpha] \ge \operatorname{LC}[\widehat{\pi}(\boldsymbol{X}); \alpha].$$

Part 4 - Performance criteria for auto-calibrated predictors

Context

- There are many tools for model selection in machine learning : deviance criteria, Gini index, concentration and Lorenz curves, correlation coefficients, ...
- Interesting facts:
 - **Deviance is a consistent scoring rule** for the mean;
 - Not the case in general for the other measures listed before.
 - \Rightarrow Working with these tools for model selection may thus lead to a wrong model choice.
- Restricting the Gini index to the class of autocalibrated regression models makes it a consistent scoring rule (Wüthrich, 2023).

Pearson's linear correlation

- ► Correlation coefficients are often used to assess the strength of dependence within the pair (Y, π̂(X)).
- The most elementary correlation coefficient is Pearson's linear one:

$$r(Y,\widehat{\pi}(\boldsymbol{X})) = rac{\mathbb{C}\mathsf{ov}[Y,\widehat{\pi}(\boldsymbol{X})]}{\sqrt{\mathsf{Var}[Y]\mathsf{Var}[\widehat{\pi}(\boldsymbol{X})]}}.$$

Proposition 4.1: (Denuit and Trufin, 2023) Let π̂(X) be an autocalibrated predictor. Then, r(Y, π̂(X)) is maximum if, and only if, π̂(X) = μ(X).

Gini index and ICC

The integral of the concentration curve (ICC) over the whole interval [0, 1] is then given by

$$\begin{split} \mathrm{ICC}[\boldsymbol{Y}, \widehat{\boldsymbol{\pi}}(\boldsymbol{X})] &= \int_{0}^{1} \mathrm{CC}[\boldsymbol{Y}, \widehat{\boldsymbol{\pi}}(\boldsymbol{X}); \boldsymbol{\alpha}] \mathrm{d}\boldsymbol{\alpha} \\ &= \frac{\mathsf{E}\left[\boldsymbol{Y} \int_{0}^{1} \mathsf{I}[\widehat{\boldsymbol{\pi}}(\boldsymbol{X}) \leq F_{\widehat{\boldsymbol{\pi}}}^{-1}(\boldsymbol{\alpha})] \mathrm{d}\boldsymbol{\alpha}\right]}{\mathsf{E}[\boldsymbol{Y}]}. \end{split}$$

The Gini index can be defined as

$$\operatorname{Gini}[\boldsymbol{Y}, \widehat{\boldsymbol{\pi}}(\boldsymbol{X})] = \frac{\frac{1}{2} - \operatorname{ICC}[\boldsymbol{Y}, \widehat{\boldsymbol{\pi}}(\boldsymbol{X})]}{\frac{1}{2} - \int_0^1 \operatorname{CC}[\boldsymbol{Y}, \boldsymbol{Y}; \alpha] \mathrm{d}\alpha}.$$

Gini index and ICC

- The Gini index is not in general a consistent scoring rule for the mean, so neither is the ICC (Wüthrich, 2023).
- BUT the Gini index gives a consistent scoring rule within the class of autocalibrated regression functions (Wüthrich, 2023).

Concentration curve

Proposition 4.3: (Denuit and Trufin, 2023) For any autocalibrated predictor \(\hat{\alpha}\), we have

 $\operatorname{CC}[Y, \widehat{\pi}(X); \alpha] \ge \operatorname{CC}[Y, \mu(X); \alpha] \quad \text{ for any } \alpha \in (0, 1),$

with an identity if, and only if, $\widehat{\pi}(\mathbf{X}) = \mu(\mathbf{X})$.

- Since the LC coincides with the CC for autocalibrated predictors, Proposition 4.3 shows that the true model also has the smallest Lorenz curve.
- This legitimates model assessment based on Lorenz curve under autocalibration, as it is often performed in practice!

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