Optimization heuristics for determining internal grading scales

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Optimization Heuristics for Determining Internal Rating Grading Scales

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Abstract

Basel II imposes regulatory capital on banks related to the default risk of their credit portfolio. Banks using an internal rating approach compute the regulatory capital from pooled probabilities of default. These pooled probabilities can be calculated by clustering credit borrowers into different buckets and computing the mean PD for each bucket. The clustering problem can become very complex when Basel II regulations and real-world constraints are taken into account.

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into account. Search heuristics have already proven remarkable performance in tackling this problem as complex as it is. A Threshold Accepting algorithm is proposed, which exploits the inherent discrete nature of the clustering problem. This algorithm is found to outperform alternative methodologies already proposed in the literature, such as standard \( k \)-means and Differential Evolution. Besides considering several clustering objectives for a given number of buckets, we extend the analysis further by introducing new methods to determine the optimal number of buckets in which to cluster banks’ clients.

**Keywords:** credit risk, probability of default, clustering, Threshold Accepting, Differential Evolution.

1 **Introduction**

The second Basel Accord on Banking Supervision requires banks to hold a minimum level of shareholders’ capital in excess of provisions. This regulatory capital (\( RC \)) may be regarded as some form of self-insurance (in excess of provisions) against the consequences of an unexpectedly high number of defaults. This amount of capital depends on the exposure to risk of the bank. Financial intermediaries then have to assess the clients’ riskiness by evaluating their probability of default (\( PD \)), i.e., the probability that a borrower will default over the subsequent 12 months. Afterwards, clients are pooled together in buckets (\( PD \)-buckets) and are assigned the same “pooled” \( PD \).
While many studies have been devoted to the phases of rating assignment, quantification, and validation, the problem of determining the width and the number of $PD$ buckets has received much less attention. We propose to fill the gap in the literature by proposing an error-based statistical methodology to determine the optimal structure of $PD$-buckets. Thereby, we consider the problem of determining the $PD$-buckets as a clustering problem, where the aim is to find the cluster structure that allows to minimize a given error measure under the relevant real-world constraints. Previous related work can be found, e.g., in Foglia et al. (2001), Krink et al. (2007) and Krink et al. (in press). We extend the analysis mainly in two directions.

First, we propose a methodology not only to tackle the problem of determining the $PD$ buckets width, but also to determine the optimal number of buckets in which to partition the banks' clients. This problem is complex to tackle since there is a trade-off between having a small number of large buckets and a high number of small buckets. In fact, clients belonging to the same buckets are assigned the same pooled $PD$. Hence, we would like to have a large number of buckets in order to minimize the loss of precision. However, in such a case it would be difficult to validate the consistency of the rating scheme ex post, since the number of defaults in each bucket would probably be too low for statistical validation. On the contrary, if the number in which to partition the clients is small, buckets tend to be too wide which might lead to an overstatement of the capital charge, given the concave shape of the capital function (Kiefer and Larson, 2004), and to opportunistic behavior.
and adverse selection of clients.

Second, we introduce the Threshold Accepting (TA) algorithm (Dueck and Scheuer, 1990; Winker, 2001) in order to determine the optimal PD buckets structure. Compared to the Differential Evolution (DE) heuristic used in previous studies (Krink et al., 2007), TA is particularly suited for discrete search spaces. By exploiting the discreteness of the search space of the PD bucketing problem, it avoids to search on plateaus of the objective function, but can still deal with local minima. Our extensive investigation on a real-world dataset shows that TA can be a faster and more robust alternative to DE.

The paper is structured as follows. Section 2 introduces the formal framework for the error-based approach to PD bucketing by considering the regulations put forward by the Basel II accord and some other real-world constraints. Several objective functions and constraints for the optimization problem are presented. Section 3 describes the two optimization heuristics, namely Differential Evolution and Threshold Accepting. Empirical results and performance comparison are then reported in Section 4. Section 5 extends our formal framework by introducing the endogenous choice of the optimum number of buckets and discusses some results. Finally, Section 6 concludes and suggests further research perspectives.
2 Basel II and Clustering of Credit Risk

The framework of Basel II puts a strong emphasis on the adequacy of banks’ equity for a given risk profile. Thereby, a core risk measure is the value at risk. A bank’s value at risk (\(VaR_i\)) associated with some borrower \(i\) is equal to this debtor’s exposure at default (\(EAD_i\)) times the fraction (loss given default, \(LGD_i\)) of \(EAD_i\) that may not be recovered. A bank may account for the expected part of \(VaR\) (i.e., \(VaR\) times borrower \(i\)’s probability of default \(PD_i\)) by provisioning. However, under sufficiently negative economic conditions the conditional (also called stressed or through the cycle) probability of default (\(PD_{c,i}\)) is likely to exceed \(PD_i\) and thus may cause losses in excess of provisions. In order to ensure the stability of the banking system, banks are required by the second Basel Capital Accord (Basel II) to hold regulatory capital (\(RC\)) that is related to these unexpected losses.

For determining \(RC\) borrowers have to be assigned to at least seven internal borrower grades \(b\) (also called groups or buckets) for non-defaulted borrowers based on their creditworthiness. Then, \(RC\) can be computed by e.g. treating the mean \(PD\) (\(PD_b\)) of all borrowers in bucket \(b\) as a proxy of an individual borrower’s \(PD\). We assume that a bank employs a statistical default prediction model so that an estimator for each borrower’s individual \(PD\) is available. Then, \(RC\) for an individual borrower (\(RC(PD_i)\)), when no maturity adjustment is considered, is given by equation (1) where the stressed \(PD\) (\(PD_{c,i}\)) is given by equation (2). \(\Phi\) and \(\Phi^{-1}\) denote the cumu-
lative standard normal density function and its inverse, respectively. The asset correlation $R$ reflects how the individual PDs are linked together by the general state of the economy, the firm’s size (as measured by sales) and the size of their $EAD$.

$$RC(PD_i) = 1.06 \cdot EAD_i \cdot LGD \cdot (PD_{c,i} - PD_i).$$ \hspace{1cm} (1)

$$PD_{c,i} = \Phi \left( \frac{\Phi^{-1}(PD_i) - \sqrt{R_i} \cdot \Phi^{-1}(0.001)}{\sqrt{1 - R_i}} \right).$$ \hspace{1cm} (2)

If a borrower $i$ is assigned to bucket $b$ her conditional PD ($PD_{c,i,b}$) can be determined by replacing $PD_i$ with $\overline{PD}_b$ in equation (2). The sum of $RC$ for all borrowers $i$ over all buckets $b$ may be computed as (3):\(^1\)

$$RC = \sum_b \sum_i 1.06 \cdot EAD_i \cdot LGD \cdot (PD_{c,i,b} - \overline{PD}_b).$$ \hspace{1cm} (3)

Computing $RC$ from pooled PDs as shown above results in an approximation error. Therefore, Basel II requires banks to perform credit risk rating, i.e., assigning borrowers to buckets meaningfully. On the one hand, this means to maximize the homogeneity of borrowers within a given bucket. This may be done by grouping borrowers in minimizing some objective func-

\(^1\)In our implementation we compute the asset correlation according to paragraph 273 of the Basel II framework by normalizing debtors’ sales to EUR 5 million if they are below that threshold and to EUR 50 million if they are above this threshold. Consequently, we do not treat small firms’ exposures as retail exposures as stated in paragraph 232. 1.06 is an empirically derived scaling factor that prevents $RC$ calculated under Basel II to drop below $RC$ under the Basel I framework.
tion using an optimization technique as described in Section 3. On the other hand, adjacent buckets must be clearly distinguishable, i.e., heterogeneous. There is a trade-off between homogeneity and heterogeneity since increasing the number of buckets is likely to decrease heterogeneity within buckets but raise homogeneity between buckets. This trade-off as well as the necessity to ex post validate the meaningfulness of the credit risk rating system leads us to the question which number of buckets to choose. We will address this issue in Section 5.

The goal of maximizing within-buckets homogeneity may be operationalized by different objective functions. First, one may minimize the squared error that arises from substituting a borrower’s individual $PD$ by the mean of its bucket. This may be done using unconditional $PD$s (point-in-time approach) resulting in the following objective function:

$$\min \sum_b \sum_{i \in b} (PD_i - \overline{PD}_b)^2.$$  

(4)

However, if a bank’s portfolio is strongly affected by overall business conditions the use of conditional $PD$s in a through-the-cycle approach may be more appropriate:

$$\min \sum_b \sum_{i \in b} (PD_{c,i} - \overline{PD}_{c,b})^2.$$  

(5)

It may be supposed that banks grant higher loans to good borrowers than to borrowers with a relatively high $PD$. Thus, VaR that arises from a good borrower may be comparatively high, as well. Consequently, it might be more
reasonable to use weighted versions of the objective functions (4) and (5) using the $EAD$s as weights, e.g., for the conditional $PD$s (5):

$$min \sum_b \sum_{i \in b} EAD_i : (PD_{c,i} - \overline{PD}_{c,b})^2 .$$

(6)

Moreover, banks may want to minimize excess regulatory capital resulting in objective function 7:

$$min \sum_b \sum_{i \in b} |RC (PD_i) - RC (\overline{PD}_b)| .$$

(7)

Apart from the selection of an appropriate objective function, several constraints imposed by the Basel II framework have to be taken into account when rating credit risk. First, according to paragraph 285 of the framework the pooled $PD$ for corporate and bank exposures must be no smaller than 0.03%. Second, paragraphs 403 and 406 of the framework require banks to have a meaningful distribution of exposures without excessive concentrations. Thus, following Krink et al. (2007), we assume that no bucket may contain more than 35% of a bank’s total exposure:

$$\frac{\sum_{i \in b} EAD_{b,i}}{\sum_b \sum_{i \in b} EAD_{b,i}} \leq 35\% .$$

(8)

Third, in order to avoid buckets that are too small, the number of borrowers

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\[2\]This constraint is not binding in our application since no $PD$ in our dataset is smaller than 0.03%.
in a bucket \((N_b)\) should be larger than some percentage \(x\) of the entire number of borrowers \(N\):

\[
N_b \geq x \cdot N.
\]  

(9)

Again following Krink et al. (2007), we will assume \(x = 1\%\) for our application in Section 4. However, we will define \(x\) based on statistical criteria when endogenizing the number of buckets in Section 5.

Fourth, the clustering algorithm must be set up such that buckets do not overlap and the union of buckets is the set of all borrowers. Furthermore, paragraph 404 of the framework requires banks to have at least seven borrower grades for non-defaulted borrowers.

3 Two Optimization Heuristics for Credit Risk Bucketing

We tackle the \(PD\) bucketing problem as a clustering one, i.e., we want to determine the optimal partition of \(N\) bank clients in \(B\) buckets with respect to a given objective function and subject to some constraints (see Section 2). Since clustering problems are NP-hard when the number of cluster exceeds three (Brucker, 1978), stochastic search heuristics, such as Differential Evolution and Threshold Acceptance, can be a valid tool to tackle such problems. Furthermore, the presence of constraints narrows and segments further the search space. DE and TA allow to explore the whole search space, not focus-
ing on the borders resulting from the constraints as conventional approaches often do. Following Krink et al. (2007), we build candidate solutions in TA or DE to encode the thresholds of buckets. Hence, when considering the problem in a continuous domain, the fitness landscape has large plateaus given that a change in the threshold of one bucket modifies the categorization only if there are some clients in the \( PD \) interval between the old and the new thresholds, e.g., if a threshold varies from 0.2 to 0.21, the \( PD \)-bucketing partition would vary only if there are clients with \( PD \) in the interval \([0.2, 0.21]\). Then, the fitness value of each individual will vary across generation only when the new bucketing thresholds correspond to a new categorization. Given this inherent discrete nature of the problem, we expect TA to be a better alternative than DE.

### 3.1 Differential Evolution

Differential Evolution (DE) is a search heuristic, introduced by Storn and Price (1997), which has shown remarkable performance in continuous numerical problems when compared with other heuristics (e.g., Corne et al. (1999), Price et al. (2005), and Vesterstom and Thomsen (2004)). The main advantage of DE is its robustness in converging towards the optimal solution and the insensitiveness to parameter tuning. Recent investigations have shown that such heuristics can be a reliable tool in tackling real-world financial problems, such as index tracking (Maringer and Oyewumi (2007)) and multi-objective portfolio optimization (Krink and Paterlini (2008)). Even if
DE is specialized on continuous numerical problems, DE has already shown better performance than GA and PSO in tackling the credit risk bucketing Krink et al. (2007).

Algorithm 1 Pseudocode for Differential Evolution.

1: Initialize parameters $n_p, n_G, F$ and $CR$
2: Initialize population $P_{j,i}^{(1)}$, $j = 1, \cdots, B - 1$, $i = 1, \cdots, n_p$
3: for $k = 1$ to $n_G$ do
   4: $P^{(0)} = P^{(1)}$
   5: for $i = 1$ to $n_p$ do
      6: Generate $r_1, r_2, r_3 \in 1, \cdots, n_p$, $r_1 \neq r_2 \neq r_3 \neq i$
      7: Compute $P_{j,i}^{(v)} = P_{j,i}^{(0)} + F \times (P_{r_1,j}^{(0)} - P_{r_2,j}^{(0)})$
      8: for $i = 1$ to $B - 1$ do
         9: if $u < CR$ then
            10: $P_{j,i}^{(u)} = P_{j,i}^{(v)}$
         11: else
            12: $P_{j,i}^{(u)} = P_{j,i}^{(0)}$
         13: end if
      14: end for
      15: if $f(P_{j,i}^{(u)}) < f(P_{j,i}^{(0)})$ then
         16: $P_{j,i}^{(1)} = P_{j,i}^{(u)}$
      17: else
         18: $P_{j,i}^{(1)} = P_{j,i}^{(0)}$
      19: end if
   20: end for
21: end for

Algorithm 1 describes the general outline of our DE implementation. Initially, the algorithm randomly generates and evaluates $n_p$ candidate solutions (2:). Each solution is made of $B - 1$ buckets thresholds. Next, for a predefined number of generations, $n_G$, the following steps are repeated. For each current element of the population a new candidate solution is generated through differential mutation (7:) and uniform crossover with the current element (9:-14:). Differential mutation generates a new solution by multiplying
the difference between two randomly selected solution vectors by the scaled factor $F$ and adding the result to a third vector. Then, a uniform crossover is applied. During crossover, the algorithm recombines the initial solution with the new candidates by replacing each component $P_{j,i}^{(0)}$ with mutant ones resulting from the differential mutation step $P_{j,i}^{(v)}$ with a probability $CR$. The resulting new trial solution is denoted by $P_{j,i}^{(u)}$. Then, the objective function $f$ is calculated and the new candidate solution replaces the current one only if it has better fitness value. Otherwise, the current solution is carried over to the next generation. The algorithm terminates after a predefined number of generations.

The tuning parameters of the DE implementation are the scaling factor $F$ and the crossover probability $CR$. These settings might affect the quality of results depending on the properties of the problem. To determine the parameter values that result in the best objective function values, we run the algorithm 30 times for different combinations of $F$ and $CR$, ranging between 0.5 and 0.9. Thereby, the objective function (6) was considered. The distribution of the results indicate that, for the specific problem instance, tuning the technical parameters does not affect the solution quality for values of $CR$ larger than 0.6, while the choice of $F$ within the given interval appears to be irrelevant.
3.2 Threshold Accepting

The idea of TA is to iteratively compare the objective function values of two candidate solutions that belong to the same neighborhood and to select one of them for further refinement. Thereby, the current candidate solution is replaced by a new one

- if this results in an improvement of the objective function value, and
- if a deterioration of the objective function value does not exceed a threshold as defined by a threshold sequence.

Due to the second feature, TA may overcome local optima.

TA requires to set an initial candidate solution and a criterion that terminates the search process. It turns out to be best to determine an initial candidate solution completely at random. Moreover, the search is stopped after a predetermined number of iterations. A nice feature of this stopping criterion is that the computation time can be controlled quite effectively.

In TA the current candidate solution is compared with a neighboring solution. Thus, the implementation requires to define a neighborhood structure. It is adjuvant to define neighborhoods quite large at the beginning of the search but small towards its end. The idea underlying this procedure is to put more emphasis on exploring wide areas of the search space first but emphasizing a narrow search and refinement of a supposedly good candidate solution towards the end of the search.
Suppose the TA algorithm has generated for 7 bucket the starting solution $g_c = (3\%, 6\%, 10\%, 12\%, 17\%, 21\%)$ and PDs in our dataset are bound by the interval $[0.2\%; 24\%]$. Suppose further that the second bucket threshold is randomly selected for modification. The new candidate solution will be a neighbor to the old one if the second bucket threshold is determined randomly from all PDs in the interval $]3\%; 10\%[$. The intervals for the remaining bucket thresholds can be found accordingly. The procedure is illustrated in Figure 1.

![Bucket intervals](image)

Figure 1: Bucket intervals.

As the search proceeds, these intervals shrink linearly in the current number of iterations relative to the total number of iterations. I.e., the contraction factor takes the form $[(I + 1) - i]/I$. Consequently, after performing for example 20\% of the iterations the second bucket threshold would be determined from the interval $]6\% - 0.8 \cdot (6\% - 3\%); 6\% + 0.8 \cdot (10\% - 6\%)[$.

New candidate solutions are generated from old ones by first determining randomly a bucket threshold of the current candidate solution and then replacing it with a random element from the above interval. This procedure is advantageous in at least two aspects. First, the objective function value of the new candidate solution $g_n$ differs from the objective function value of the
current candidate solution \( g_c \) only in the contribution of the two buckets that are affected by the alteration. Thus, fast updating of the objective function is feasible. Moreover, computation time becomes vastly independent of the number of buckets. This is due to the fact that for any number of buckets \( \text{TA} \) only has to compute the fitness of two buckets per iteration. On the contrary, in \( \text{DE} \), as implemented in (Krink et al., 2007), the fitness for all buckets is computed in every iteration. This results in a higher computation time. This disadvantage of \( \text{DE} \) becomes more pronounced for higher numbers of buckets.\(^3\) Second, since in \( \text{TA} \) new bucket thresholds are chosen from the \( PDs \) in the dataset, each new candidate solution constitutes a different partition and, consequently, a different value of the objective function which is not the case for our \( \text{DE} \) implementation on a continuous search space.

A final crucial element of any \( \text{TA} \) implementation is its threshold sequence since it determines \( \text{TA}'s \) ability to overcome local optima. Basically, the idea is to accept \( g_n \) if its objective function value is better or if it is not much worse than that of \( g_c \) where \textit{not much worse} means the deterioration may not exceed some threshold \( T \) defined by the threshold sequence.

We propose a threshold sequence that is based on the differences in the fitness of candidate solutions that are found in a certain area of the search space. Instead of using an ex ante simulation of local differences of the fitness function as proposed by Winker and Fang (1997), the local differences\(^3\) It has to be left for future research to consider updating rules for \( \text{DE} \) similar to the ones employed for \( \text{TA} \) in the present application.
actually calculated during the optimization run are considered. By using a moving average, a smooth threshold sequence is obtained. Algorithm 2 provides the pseudocode for the TA implementation with the data driven generation of the threshold sequence.

**Algorithm 2** Pseudocode for TA with data driven generation of threshold sequence.

1: Initialize $I$, $Ls = (0, \ldots, 0)$ of length 100
2: Generate at random an initial solution $g_c$, set $T = f(g_c)$
3: for $i = 1$ to $I$ do
4:  Generate at random $g_n \in N(g_c)$
5:  Delete first element of $Ls$
6:  if $f(g_n) - f(g_c) < 0$ then
7:    add $|f(g_n) - f(g_c)| \cdot (i/I)$ as last element to $Ls$
8:  else
9:    add $|f(g_n) - f(g_c)| \cdot (1 - i/I)$ as last element to $Ls$
10: end if
11: $T = Ls \cdot (1 - i/I)$
12: if $f(g_n) + T \leq f(g_c)$ then
13:  $g_c = g_n$
14: end if
15: end for

The threshold sequence is calculated during the run time of the algorithm and exhibits the following properties. First, it adapts to the region of the search space to which the current solution belongs. Second, it takes into account the current definition of the neighborhood. Third, and most importantly, it adapts to the objective function used. As a result, this data driven threshold sequence is readily available for use with any objective function, constraint handling technique or neighborhood structure and does not require any fine-tuning.
The current value of the threshold $T$ is defined as the weighted mean $\overline{Ls}$ over the last 100 fitness differences (11:). A general requirement in TA is that thresholds should be larger at the beginning of the search in order to overcome local optima and decrease to zero at the end in order to reach at least a local, if not the global optimum. In order to satisfy this requirement, the weighted mean $\overline{Ls}$ is multiplied with a scaling factor decreasing linearly from one to zero with the number of iterations (11:).

Apart from this global weights, each fitness difference entering the vector $Ls$ obtains a particular weight. At the beginning of the search process, one might expect many fitness improvements. For not being too generous in accepting deteriorations of the objective function, objective fitness differences corresponding to improvements are downweighted by the factor $i/I$, i.e., the share of iterations already done (7:). In contrast, towards the end of the search procedure, one has to expect that most trials result in a deterioration of the objective function. To avoid too generous thresholds, the corresponding elements of $Ls$ are downweighted by the factor $(1-i/I)$ decreasing to zero with the number of iterations (9:). It is obvious that this threshold sequence adapts to the local structure of the search space. If the algorithm moves candidate solutions towards an optimum, fitness improvements are likely to become smaller the closer the algorithm approaches this optimum. Then, $T$ declines and forces the algorithm not to deviate from its track towards the optimum. Once a (local) optimum is found only fitness deteriorations will be observed resulting which makes $T$ to increase and eventually allows the
algorithm to depart from that optimum and examine another part of the search space. By using a moving average, a smooth threshold sequence is obtained (11:).

3.3 Constraint Handling

When running the optimization heuristics TA and DE, the constraints described in Section 2 have to be taken into account. To this end, two alternative methods can be considered: rewriting the definition of domination, such that it includes the constraint handling (Deb et al., 2002) or imposing a penalty on infeasible solutions.

The first possibility has been described for the current application in Krink et al. (2007). The intuitive idea of this constraint handling technique is to leave the infeasible area of the search space as quickly as possible and never return. For minimization problems, the procedure can be described as follows within Algorithm 2:

1. If the new candidate solution $g_n$ and the current candidate solution $g_c$ satisfy the constraints, $g_n$ replaces $g_c$ if its fitness $f(g_n)$ satisfies the condition $f(g_n) + T \leq f(g_c)$. In TA $T$ represents the threshold as defined by the threshold sequence. In DE, we set $T = 0$.

2. If only one candidate solution is feasible, select the feasible one

3. If both solutions violate constraints, . . .
(a) ... select the one that violates fewer constraints.

(b) ... if both solutions violate the same number of constraints, \( g_n \) replaces \( g_c \) if its fitness \( f(g_n) \) satisfies the condition \( f(g_n) + T \leq f(g_c) \). Again, \( T \) either takes a value as defined by the threshold sequence or we set \( T = 0 \) in DE.

In contrast, the penalty technique allows infeasible candidate solutions while running the algorithm as a stepping stone to get closer to promising regions of the search space. In this case, the objective function is multiplied by a penalty term. Solutions should be penalized the stronger the more they violate the constraints. Moreover, in order to guarantee a feasible solution at the end, the penalty should increase over the runtime of the algorithm. Equation (10) states that the objective function value \( f_u \) of a candidate solution is increased by some penalty factor \( A \in [1; 2] \) that puts more weight on penalties the more the current iteration \( i \) approaches the overall number of iterations \( I \). The exponent \( a \) may take values in the interval \([0; 1]\). No penalty is placed on \( f_u \) if no constraint is violated so that \( a = 0 \). However, if the constraints are violated most strongly, i.e., all borrowers are concentrated in one bucket leaving the remaining buckets empty, the exponent takes the value \( a = 1 \). A more formal description of this penalty technique is given in the appendix.

\[
f_c(g) = f_u(g) \cdot A = f_u(g) \cdot \left(1 + \sqrt{\frac{i}{I}}\right)^a \tag{10}
\]
For the current application, DE is only implemented with the constraint-dominated handling technique, while for TA both methods are implemented. Generally, the constraint-dominated handling technique performs well while taking comparatively little computation time. However, depending on the kind of objective function used the penalty technique may improve the reliability of TA, i.e., reduce the variance of the results obtained.

4 Results and Relative Performance

For our empirical application we consider the dataset comprising 11,995 defaulted and non-defaulted borrowers of a major Italian bank already analyzed by Krink et al. (2007). The $PDus$ range between 0.21% and 23.94%. Moreover, conditional probability of default ($PDc$) was computed using equation (2). The $PDcs$ range between 4.52% and 64.88%.

All algorithms are implemented from scratch in Matlab 7.6 and run on a PC with Intel Duo Core processors operating at 2.40 GHz and running Windows XP.

4.1 Results for Fixed Number of Buckets

Tables 1 to 3 report the empirical results of the two heuristic algorithms for 7, 10, and 15 buckets, the two different constraint handling techniques and the three objective functions described above. Both algorithms were restarted 30 times on each problem instance to control for the stochasticity
of heuristic optimization techniques. For the comparison of the two methods, we report the best value, the median, the worst value, the variance, the 80% percentile, the 90% percentile, and the frequency the best value occurs in all 30 repetitions.

Table 1: Objective function (5) in EUR

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<th>Best</th>
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<th>Worst</th>
<th>s.d.</th>
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<th>q90%</th>
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\(B = 7\)

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<tr>
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\(B = 10\)

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<td>3.9190</td>
<td>3.9366</td>
<td>0.0080</td>
<td>3.9155</td>
<td>3.9366</td>
<td>19/30</td>
</tr>
<tr>
<td>(DE)</td>
<td>3.9155</td>
<td>9.9319</td>
<td>4.1663</td>
<td>0.0496</td>
<td>3.9195</td>
<td>3.9527</td>
<td>2/30</td>
</tr>
</tbody>
</table>

\(B = 15\)

<p>| | | | | | | | |</p>
<table>
<thead>
<tr>
<th></th>
<th></th>
<th></th>
<th></th>
<th></th>
<th></th>
<th></th>
<th></th>
</tr>
</thead>
<tbody>
<tr>
<td>(TA^a)</td>
<td>2.8842</td>
<td>2.8848</td>
<td>2.8929</td>
<td>0.0016</td>
<td>2.8855</td>
<td>2.8855</td>
<td>6/30</td>
</tr>
<tr>
<td>(TA^b)</td>
<td>2.8842</td>
<td>2.8874</td>
<td>2.9064</td>
<td>0.0053</td>
<td>2.8929</td>
<td>2.8933</td>
<td>7/30</td>
</tr>
<tr>
<td>(DE)</td>
<td>2.8964</td>
<td>2.9761</td>
<td>3.0199</td>
<td>0.0428</td>
<td>3.0083</td>
<td>3.0140</td>
<td>1/30</td>
</tr>
</tbody>
</table>

\(^a\)Rejection based constraint handling technique  
\(^b\)Penalty technique

For DE the initial parameter settings were population size \(n_p = 100\) and number of generations \(n_G = 1000\), while the scaling factor \(F\) and the crossover rate \(CR\) were kept constant at 0.5 and 0.8, respectively.\(^4\) In the case of TA, the algorithm was run for \(I = 100 000\) iterations, in order to attain analogy with DE’s population size and number of generations.\(^5\)

\(\text{extensive parameter tuning on DE suggests that DE is rather insensitive to the choice of } F \text{ and } CR\). Results are available upon request.\(^4\)  
\(\text{It should be noted that due to the local updating method in TA, the 100 000 iterations}\)
Table 2: Objective function (6) in EUR

<table>
<thead>
<tr>
<th></th>
<th>Best</th>
<th>Mean</th>
<th>Worst</th>
<th>s.d.</th>
<th>q80%</th>
<th>q90%</th>
<th>Freq</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>4,582.86</td>
<td>4,582.86</td>
<td>4,582.86</td>
<td>0.0000</td>
<td>4,582.86</td>
<td>4,582.86</td>
<td>30/30</td>
</tr>
<tr>
<td></td>
<td>4,582.86</td>
<td>4,582.86</td>
<td>4,582.86</td>
<td>0.0000</td>
<td>4,582.86</td>
<td>4,582.86</td>
<td>30/30</td>
</tr>
<tr>
<td></td>
<td>4,582.86</td>
<td>4,587.35</td>
<td>4,671.38</td>
<td>16.5569</td>
<td>4,583.52</td>
<td>4,585.09</td>
<td>2/30</td>
</tr>
<tr>
<td>$B = 10$</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>3,471.68</td>
<td>3,479.97</td>
<td>3,483.92</td>
<td>1.8592</td>
<td>3,480.01</td>
<td>3,480.21</td>
<td>1/30</td>
</tr>
<tr>
<td></td>
<td>3,471.68</td>
<td>3,480.33</td>
<td>3,483.92</td>
<td>2.8705</td>
<td>3,483.66</td>
<td>3,483.92</td>
<td>2/30</td>
</tr>
<tr>
<td></td>
<td>3,471.51</td>
<td>3,475.47</td>
<td>3,498.96</td>
<td>5.4891</td>
<td>3,479.96</td>
<td>3,480.18</td>
<td>4/30</td>
</tr>
<tr>
<td></td>
<td>2,821.00</td>
<td>2,833.55</td>
<td>2,865.98</td>
<td>14.1177</td>
<td>2,844.24</td>
<td>2,856.92</td>
<td>8/30</td>
</tr>
<tr>
<td></td>
<td>2,821.00</td>
<td>2,830.55</td>
<td>2,860.02</td>
<td>11.5971</td>
<td>2,840.99</td>
<td>2,844.24</td>
<td>3/30</td>
</tr>
<tr>
<td></td>
<td>2,866.23</td>
<td>2,943.09</td>
<td>3,122.39</td>
<td>57.9240</td>
<td>2,958.48</td>
<td>3,005.19</td>
<td>1/30</td>
</tr>
</tbody>
</table>

$^a$Rejection based constraint handling technique

$^b$Penalty technique

Table 1 presents a statistical summary of the results using objective function (5). The TA algorithm was run using both the rejection based constraint handling technique and the penalty technique. The results are affected by the choice of the constraint handling technique, as the best value is obtained with an equal or higher frequency when using the penalty technique. However, these results cannot be generalized for the alternative objective functions (6) and (7) (see Tables 2 and 3), especially for a larger number of buckets, i.e., $B = 10, 15$. In general, the performance of the TA implementation is excellent for the case of seven buckets and still gives good results with low variance of TA will require less computing time than the corresponding run of DE. The relative merits of both methods in terms of computational load and result quality are reported in Section 4.2.
for the larger problem instances.

Table 3: Objective function (7) in EUR

<table>
<thead>
<tr>
<th></th>
<th>Best</th>
<th>Mean</th>
<th>Worst</th>
<th>s.d.</th>
<th>q80%</th>
<th>q90%</th>
<th>Freq</th>
</tr>
</thead>
<tbody>
<tr>
<td>B = 7</td>
<td>TA&lt;sup&gt;a&lt;/sup&gt; 45,791.49 45,793.18 45,825.62</td>
<td>6.4870</td>
<td>45,791.49 45,791.72</td>
<td>26/30</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>TA&lt;sup&gt;b&lt;/sup&gt; 45,791.49 45,794.11 45,826.57</td>
<td>7.7787</td>
<td>45,791.49 45,801.65</td>
<td>25/30</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>B = 10</td>
<td>TA&lt;sup&gt;a&lt;/sup&gt; 31,942.19 31,951.30 31,996.89</td>
<td>19.9770</td>
<td>31,946.42 31,994.72</td>
<td>21/30</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>TA&lt;sup&gt;b&lt;/sup&gt; 31,942.19 31,951.25 31,994.73</td>
<td>18.8050</td>
<td>31,946.42 31,992.74</td>
<td>18/30</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>DE 31,995.28 32,166.86 32,299.00</td>
<td>119.9837</td>
<td>32,299.00 32,299.00</td>
<td>1/30</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>B = 15</td>
<td>TA&lt;sup&gt;a&lt;/sup&gt; 20,711.93 20,729.26 20,973.53</td>
<td>62.6797</td>
<td>20,713.36 20,714.88</td>
<td>10/30</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>TA&lt;sup&gt;b&lt;/sup&gt; 20,711.93 20,725.99 20,951.01</td>
<td>49.5889</td>
<td>20,714.88 20,715.61</td>
<td>1/30</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>DE 20,970.37 24,916.30 35,003.82</td>
<td>4875.7503</td>
<td>31,215.84 33,676.88</td>
<td>1/30</td>
<td></td>
<td></td>
<td></td>
</tr>
</tbody>
</table>

<sup>a</sup>rejection based constraint handling technique
<sup>b</sup>penalty technique

Considering the performance of the DE implementation, we observe that the best value is obtained for $B = 7$ at a frequency of 9 out of 30 restarts. While, for a higher number of buckets, the best value does not deviate much from the optimum, the efficiency worsens. The same pattern is observed for all three objective functions.

We conclude that the TA implementation is superior for most problem instances in terms of mean solution quality and variance for all objective functions considered. The clustering of credit risk is a problem on a discrete search space. In contrast to the DE algorithm, the TA implementation takes this discrete feature of the search space into account. This might explain its
superior performance.

4.2 Relative Performance of DE and TA

Section 4.1 provides evidence of the good performance of both algorithms for the credit risk bucketing problem. Given that TA exploits the discrete structure of the search space and uses a local updating procedure, it is significantly faster than DE for a given number of function evaluations.

Therefore, in order to obtain a fair comparison of both algorithms, we consider two settings. First, we analyze the distribution of results obtained from both algorithms when running them for the same time. Second, we fix a quality goal, e.g., not to deviate by more than 1% from the best solution documented above. Then, both algorithms are run using increasing computational time until at least 50% of the restarts meet the quality goal.

For the first approach, the following setup is used. We run the DE algorithm with the same parameters as above, i.e., population size $n_p = 100$ and number of generations $n_G = 1000$, and – to have a comparison for a small amount of computational resources – with $n_p = 40$ and $n_G = 50$. Then, we estimate the number of iterations $I$ which can be performed by our TA algorithm using the same computational time. In fact, this number of iterations will depend on the objective function used and on the number of buckets $B$, as the advantage of updating becomes more pronounced for larger $B$.

Table 4 summarizes the findings. The first four columns report respectively the objective function, the number of buckets $B$, the population size
$n_P$ and the number of generations $n_G$ of DE. Column (5) displays the computing time for a single restart of our implementations. Column (6) reports the number of iterations $I$ in TA that equalizes computation time for DE and TA. Columns (7) and (8) displays the difference in the mean and standard deviation between TA and DE. Thereby, negative values indicate an advantage of the TA implementation. Finally, column (9) reports the number of times, TA outperforms DE.

**Table 4: Relative performance of DE and TA for given computing time**

<table>
<thead>
<tr>
<th>Obj.</th>
<th>$B$</th>
<th>$n_P$</th>
<th>$n_G$</th>
<th>CPU time</th>
<th>$I$</th>
<th>$\Delta$</th>
<th>$\Delta$</th>
<th>better result for TA</th>
<th>for DE</th>
</tr>
</thead>
<tbody>
<tr>
<td>(7)</td>
<td>7</td>
<td>100</td>
<td>1000</td>
<td>32.4m</td>
<td>786600</td>
<td>-18.4</td>
<td>-38.3</td>
<td>25/30</td>
<td>0/30</td>
</tr>
<tr>
<td></td>
<td>10</td>
<td>100</td>
<td>1000</td>
<td>36.0m</td>
<td>1050000</td>
<td>-220.7</td>
<td>-108.0</td>
<td>21/30</td>
<td>0/30</td>
</tr>
<tr>
<td></td>
<td>15</td>
<td>100</td>
<td>1000</td>
<td>192.7m</td>
<td>6760511</td>
<td>-4202.8</td>
<td>-4873.9</td>
<td>5/30</td>
<td>0/30</td>
</tr>
<tr>
<td>(7)</td>
<td>7</td>
<td>40</td>
<td>50</td>
<td>42.3s</td>
<td>17000</td>
<td>-13145</td>
<td>-10210</td>
<td>6/30</td>
<td>0/30</td>
</tr>
<tr>
<td></td>
<td>10</td>
<td>40</td>
<td>50</td>
<td>51.7s</td>
<td>25000</td>
<td>-16355</td>
<td>-8630</td>
<td>6/30</td>
<td>0/30</td>
</tr>
<tr>
<td></td>
<td>15</td>
<td>40</td>
<td>50</td>
<td>66.7s</td>
<td>38000</td>
<td>-23441</td>
<td>-64880</td>
<td>1/30</td>
<td>0/30</td>
</tr>
</tbody>
</table>

Table 4 reports results for objective function (7) and $B = 7, 10, 15$ buckets. When considering the original setting for DE with $n_P = 100$ and $n_G = 1000$, the number of iterations for TA can be increased above the value of 100 000 given the same computation time. This further increase in the number of iterations does not affect the quality of results. However, when considering a smaller amount of available computational time, e.g., $n_P = 40$ and $n_G = 50$, it becomes obvious that TA still outperforms DE when using the same computational time. Despite the mean objective function value
being only slightly better, the standard deviation is drastically reduced for TA.

For the second approach mentioned above, we consider three quality levels, i.e., 10%, 5% and 1% departure from the optimum values reported. Taking into account computation time, we only report findings for objective function (6) and $B = 7$ and $B = 15$ buckets, respectively. For DE we fix $n_p = 100$ and increase the number of generations $n_G$, while for TA the number of iterations $I$ varies. For both algorithms the parameter ($n_G$ or $I$) is increased stepwise until the algorithm finds a solution meeting the quality level in at least 50% out of 30 replications. The results are summarized in Table 5 providing both the parameters actually used for the two algorithms and the corresponding CPU times for a single restart.

<table>
<thead>
<tr>
<th>Precision</th>
<th>$B$</th>
<th>$n_p$</th>
<th>$n_G$</th>
<th>DE time</th>
<th>TA time</th>
</tr>
</thead>
<tbody>
<tr>
<td>10%</td>
<td>7</td>
<td>100</td>
<td>30</td>
<td>44.44s</td>
<td>200</td>
</tr>
<tr>
<td>5%</td>
<td></td>
<td></td>
<td></td>
<td>58.45s</td>
<td>300</td>
</tr>
<tr>
<td>1%</td>
<td></td>
<td></td>
<td></td>
<td>104.24s</td>
<td>800</td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>Precision</th>
<th>$B$</th>
<th>$n_p$</th>
<th>$n_G$</th>
<th>DE time</th>
<th>TA time</th>
</tr>
</thead>
<tbody>
<tr>
<td>10%</td>
<td>15</td>
<td>100</td>
<td>100</td>
<td>5.19m</td>
<td>2000</td>
</tr>
<tr>
<td>5%</td>
<td></td>
<td>1000</td>
<td>45.75m</td>
<td>3000</td>
<td>0.05m</td>
</tr>
<tr>
<td>1%</td>
<td></td>
<td></td>
<td>*</td>
<td>15000</td>
<td>0.21m</td>
</tr>
</tbody>
</table>

*: No solution obtained for $n_G \leq 5000$ generations.

It is evident that a given quality of the solutions can be obtained much faster with TA. The relative advantage becomes even more pronounced for the larger problem instance ($B = 15$). This effect is due to the local updating
used with the TA algorithm. In fact, for $B = 15$, the quality goal of 1% from the best value could not be satisfied in at least 50% of the cases by DE even when using $n_G = 5000$ generations. For this parameter setting, a single run of the DE algorithm takes more than 4 hours of CPU time. By contrast, the same quality goal can be obtained by the TA algorithm in less than a minute.

5 Endogenous Determination of Number of Buckets

The Basel II framework requires banks to have a meaningful credit risk rating system. This does not only refer to the clustering of clients into a given number of $PD$-buckets, but also to the choice of the number of buckets. Thereby, a trade-off has to be faced. On the one hand, the clusters of borrowers should be rather homogenous. Increasing the number of buckets will reduce the loss in precision that comes from replacing individual $PD$s with pooled $PD$s. This effect causes objective function values to decline as the number of buckets is raised, resulting in a larger optimum number of $PD$-buckets.

On the other hand, both banks and regulators will be interested in an ex post validation of the classification system. For example, one may want to evaluate ex post if the observed number of defaults matches the ones predicted by the credit risk rating system. In this context, looking at the number of defaults may be seen as a proxy for evaluating whether the credit
risk rating system will predict unexpected losses correctly, which in turn, results in a statement about the adequacy of banks’ regulatory capital. Alternatively, one might consider directly the precision of the estimates of unexpected losses. A crucial factor driving the precision of any ex post evaluation is the number of borrowers per bucket. Thus, imposing a requirement on the minimum number of borrowers in a bucket based on ideas of ex post validation will result in an optimum (maximum) number of buckets still satisfying this constraint.

In the following, we will analyse both the ex post validation of the actual number of defaults and of the unexpected losses.

5.1 Validation of Actual Number of Defaults

First, we compare the actual number of defaults $D_a^b$ in a given bucket $b$ with the forecast $D_f^b$ based on the mean $\overline{PD}_b$ and the number of borrowers $N_b$:

$$D_f^b = N_b \overline{PD}_b.$$ 

In order to be able to judge whether a deviation of $D_a^b$ from $D_f^b$ should be considered as being significant, i.e., challenging the credit classification system, the distribution of $D_f^b$ has to be analyzed under the null hypothesis that $\overline{PD}_b$ is an unbiased estimator.

Given that the actual default for a loan is a binary variable, the number of
actual defaults within a bucket can be modeled by the binomial distribution.\(^6\) Consequently, a \(1 - \alpha\) confidence interval for \(D^a_b\) is defined by:

\[
P_{\text{int}} = P_b(D_{b,min} \leq D^a_b \leq D_{b,max})
= \sum_{k=D_{b,min}}^{D_{b,max}} \binom{N_b}{k} \cdot \overline{PD}^k_b \cdot (1 - \overline{PD}^k_b)^{N_b-k} \geq 1 - \alpha. \quad (11)
\]

The corresponding confidence interval for the default rates, i.e. \(D^f_b / N_b\) will shrink with a growing number of borrowers \(N_b\) in bucket \(b\). Thus, any requirement on the size of the confidence interval will impose a lower bound on \(N_b\). In a first approach, we consider symmetric confidence intervals around \(D^f_b\) of size \(2\varepsilon\) as long as the confidence interval falls in the interval \([0, 1]\), otherwise, the confidence interval is censored, i.e.,

\[
D_{b,min} = N_b \cdot \max(\overline{PD}^k_b - \varepsilon, 0). \quad (12)
\]
\[
D_{b,max} = N_b \cdot \min(\overline{PD}^k_b + \varepsilon, 1). \quad (13)
\]

The choice of an absolute definition of approximation errors rather than imposing a relative error margin is motivated by its economic impact. In fact, any deviation of the actual ex post default rates from the estimated ones by, e.g., one percentage point will have the same effect on actual defaults

\(^6\)Thereby, we made the simplifying assumption that the default risks are independent, which might be a sensible assumption for retail loans, but might be challenged for other segments of the loan market.
independent from the level of the estimated default rate ceteris paribus.

Given that we impose a minimum constraint on the precision, we do not have to solve equation (11) for the number of elements in the bucket $N_b$. Instead, for a given bucket $b$ of size $N_b$, we just have to check whether the constraint $P_{int} \geq 1 - \alpha$ is satisfied. Thus, our requirement on the precision of ex post validation imposes an additional constraint to the optimization problem.\footnote{For the consideration of this additional constraint in the penalty term, see the details provided in the appendix.}

Using this concept, we define a credit classification system as \textit{meaningful} if it allows for an ex post validation at a given level of precision as described by the two parameters $\alpha$ and $\varepsilon$. The sample composition, in particular the total sample size, and bank objectives will affect the choice of $\varepsilon$.

However, it has to be taken into account that not all combinations of $\alpha$ and $\varepsilon$ will be feasible for a given total number of loans to be considered and taking into account the other constraints imposed by the Basel II framework. In fact, a rough back of the envelope calculation shows that for our data with a mean $PD$ of around 20\% in the last bucket, values of $\alpha = 0.95$ and $\varepsilon = 1\%$ would require more than 4 000 observations in the last bucket. This is not feasible given the sample size and the other constraints imposed, in particular the constraint that no more than 35\% of total exposure at default should belong to one bucket. Therefore, we restrain from reporting results for this approach as we would have to use unreasonably low values of $\alpha$ and $\varepsilon$.\footnote{For the consideration of this additional constraint in the penalty term, see the details provided in the appendix.}
almost meaningless high values of $\epsilon$ to obtain feasible results. We leave the application of this method to future research working with a larger dataset.

If feasible, the procedure would work in two steps. First, the optimization is done for different fixed numbers of buckets. Then, the optimum number of buckets is selected to be the one corresponding to the best value of the objective function obtained for the different runs. In principle, one could also think about integrating the number of buckets as an additional parameter directly in the optimization routine. However, given its very discrete nature, the sequential procedure might be more advantageous.\footnote{A similar two-step procedure is suggested by Winker and Maringer (2004) in the context of model selection.}

Constructing buckets as previously described allows to easily validate the accuracy (as determined by the choice of $\epsilon$) of a bank’s credit risk rating system. If we find the actual number of defaults in any bucket $b$ to lie outside the interval $[D_{b,\text{min}}; D_{b,\text{max}}]$ we can state with confidence $1 - \alpha$ that the credit risk rating system is not suitable for predicting defaults in that bucket. This may have at least two reasons. First, the objective function that is used for partitioning the dataset may not be appropriate. One may easily check for this problem by using different objective functions and then assessing which ones yield results that do not cause actual defaults to lie outside the above bounds. Second, the bank employs a statistical default prediction model that does not forecast defaults correctly and thus needs to be improved.
5.2 Validating Unexpected Losses

As an alternative to an ex post validation of predicted default rates, validation may also be based on the correct statement of unexpected losses, respectively regulatory capital since $RC = 1.06 \cdot UL$. Supervisory authorities’ objective is to motivate banks to set aside equity capital equaling at least 8% of their risk-weighted assets in order to ensure the stability of the banking system. On the contrary, the objective of profit maximization requires banks’ to back up their risk-weighted assets with no more than the supervisory authority’s minimum requirements. These objectives can be operationalized by stating that in no bucket $b$ actual unexpected losses in a stress-situation ($UL_{b,a}$) shall be smaller or larger than predicted unexpected losses ($UL_b$) plus or minus some fraction $\varepsilon$ of bucket $b$’s stake in total unexpected losses as measured by the percentage of its borrowers ($N_b$) in the number of all borrowers ($N$) (see equation (14)). Total unexpected losses, unexpected losses of bucket $b$ and pooled conditional probabilities of default are given by equations (15), (16), and (17), respectively.

\[
UL_b - \varepsilon \cdot \left( UL \cdot \frac{N_b}{N} \right) \leq UL_{b,a} \leq UL_b + \varepsilon \cdot \left( UL \cdot \frac{N_b}{N} \right) \tag{14}
\]

\[
UL = 0.45 \cdot \sum (EAD_i \cdot (PD_{c,i} - PD_i)) \tag{15}
\]

\[
UL_b = 0.45 \cdot \sum_{i \in b} (EAD_i \cdot (PD_{c,b} - PD_b)) \tag{16}
\]

\[
PD_{c,b} = \frac{\sum_{i \in b} PD_{c,i}}{N_b}. \tag{17}
\]
Equation (14) is not operational since we do not know the distribution of unexpected losses. Given that we know the distribution of defaults, we can approximate $UL_{b,a}$ by $N_b \cdot UL_b$. Then, dividing equation (14) by $UL_b$ and multiplying it with $PD_b$, we obtain equation (18). This equation can be interpreted meaningfully as well since it says that for condition (14) to hold the actual number of defaults in bucket $b$ ($D_a^b$) must lie within an interval $[D_{b,min}; D_{b,max}]$. The size of this interval is determined by several parameters. Obviously, it increases with the expected probability of default and the number of borrowers in bucket $b$. Moreover, it rises with $\varepsilon$. Finally, the interval becomes larger and thus easier to satisfy if the mean unexpected loss in bucket $b$ ($UL_b$) is smaller than the mean of total unexpected loss ($UL$), i.e., the default of a borrower in this bucket is less likely to endanger the bank’s stability than an average borrower’s default. On the other hand, the interval shrinks and thus becomes harder to satisfy if borrowers are likely to cause an above average unexpected loss.

$$N_b \cdot PD_b \cdot \left[ 1 - \varepsilon \cdot \frac{UL}{UL_b} \right] \leq D_a^b \leq N_b \cdot PD_b \cdot \left[ 1 + \varepsilon \cdot \frac{UL}{UL_b} \right].$$

(18)

The central idea of this approach is to have a sufficient number of borrowers in each bucket so that we can ex ante state with a certain confidence $1 - \alpha$ that the actual number of defaults should lie within the interval $[D_{b,min}; D_{b,max}]$. Since defaults follow a binomial distribution we can express the above idea by equation (11).
In terms of our optimization problem the above condition replaces the constraint to have at least 1% of all borrowers in each bucket. Following the above idea that a high number of buckets reduces the precision error occurring from substituting individual PDs by pooled PDs we then choose as the optimum number of buckets the maximum number of buckets that is consistent with some predefined values for $\alpha$ and $\varepsilon$. In practice, these values have to be chosen meaningfully by regulation authorities and/or banks based on their objectives. As noted for the case of ex post validation of the PDs, not all combinations of $\alpha$ and $\varepsilon$ will be feasible for a given sample size.

It is of some interest how to incorporate this constraint in our optimization algorithm. When constraints are considered based on rejection of infeasible candidate solutions the algorithm described above will not change. However, if the penalty technique is used it is necessary to alter the term that captures the degree of violation of this constraint (see Appendix for details).

5.3 Results for Endogenous Number of Buckets

In this section we evaluate the quality of the UL-constraint proposed in Section 5.2. The results were obtained from running TA 30 times with 200,000 iterations. We evaluate objective functions (5) based on squared differences of PDcs, (6) based on weighted squared differences of PDcs, and (7) based on differences in RC in absolute terms. We choose $\alpha = 10\%$ since it gives
a sufficient level of confidence and allows us to choose $\varepsilon = 30\%$. Thus, if we find ex post the actual number of defaults in all buckets to lie in the interval defined by equation (18) we can state with 90% confidence that actual unexpected losses do not deviate from unexpected losses predicted by the credit risk rating system by more than $\pm 30\%$ of the buckets’ fraction (as measured by the number of borrowers) in total unexpected loss. Taking into account the small size of our sample (11,995 borrowers) we are confident that these values can be improved drastically for larger samples.

Using the UL-constraint gives similar results for objective functions (5) and (6) such that stylized facts on these functions can be presented together.\textsuperscript{10}

1. First results indicate that the best number of buckets is between 10 (for objective function (6)) and 12 (for objective function (5)).

2. When increasing the number of buckets, the algorithm does not always find a feasible solution. In fact, the UL-constraint makes the optimization problem more complex by narrowing the search space even more.

3. For a seven bucket setting an idealized solution-vector of buckets’ mean PDs looks like $g_s = (0.25\%; 0.55\%; 1.5\%; 4\%; 8\%; 14\%; 21\%)$. The UL-\textsuperscript{9}Please note that the higher a value we choose for $\alpha$ the larger will be the risk of a $\beta$-error, i.e. accepting $\overline{PD}_b$ as an unbiased estimator of bucket $b$’s true default rate while it is not.

\textsuperscript{10}Please note that without using the UL-constraint objective functions (5) and (6) produce quite dissimilar results, i.e., objective function (5) places a sizeable amount of borrowers in the first bucket while objective function (6) produces a more evenly distribution of borrowers across buckets.
constraint shapes the solution in a way that we must not reject the validity of the credit risk rating system if we find ex post actual $PD_b$s that deviate from predicted $PD_b$s by less than ± the allowed deviations (in percentage points) given by $d = (0.2\%; 0.25\%; 0.4\%; 1\%; 1.8\%; 3.5\%; 6.5\%).$

(a) We find that the $UL$-constraint imposes constraints on mean $PD$s that are of a reasonable size.

(b) The constraint on the first bucket is quite generous since it contains good borrowers that are unlikely to default.

(c) It is restrictive for mid-range borrowers allowing actual mean $PD$s to only deviate from predicted mean $PD$s by roughly 1/4. This is reasonable since it is highly uncertain whether these borrowers will default and cause a high unexpected risk for the bank.

(d) The $UL$-constraint becomes more generous for the last bucket again, allowing actual mean $PD$s to deviate from predicted mean $PD$s by roughly 1/3. This is reasonable since these borrowers’ default is quite likely such that high provisions have already been recognized. Hence, a smaller portion of their default risk must be backed up with capital requirements.

4. The rejection based constraint handling technique gives us better results (i.e., better objective function values and fewer runs converging to an infeasible solution) than the penalty technique.
Objective function (7) gives slightly different results since it allocates borrowers more evenly and especially puts less borrowers in the first bucket. We find the idealized vectors \( g_s = (0.23\%; 0.3\%; 0.6\%; 0.9\%; 3\%; 7\%; 18\%) \) and \( d = (0.2\%; 0.2\%; 0.25\%; 0.25\%; 1\%; 2\%; 5\%) \) using the terminology introduced before.

Summarizing, we find that the structure of results when using the UL-constraint is quite reasonable. It puts more emphasis on critical, i.e. mid-range borrowers and yields intervals around mean PDs that reflect the structure of borrowers. Moreover, imposing the UL-constraint somewhat increases the computational burden by narrowing down the search space. As a consequence, for 200 000 iterations the TA optimization heuristic converges towards different solutions in repeated runs. A nice feature of the UL-constraint, even for our small dataset, is to give us feasible solutions for reasonable values of \( \alpha \) and \( \varepsilon \). This enables us to test our validation-hypothesis. Thus, for a larger number of borrowers results may be expected to improve massively.

6 Conclusion

The Basel II capital accord requires banks to group loans according to their creditworthiness and set aside equity in order to self-insure against unexpected losses from borrowers’ defaults that occur under sufficiently negative economic conditions. Previous work has shown that this task can be tackled
as a clustering problem, where the objective is to minimize the loss in precision, which inevitably occurs when borrowers in the same bucket are assigned the same probability of default. Furthermore, real-world constraints can increase the complexity of the optimization problem. Optimization heuristics can then be a reliable and viable tool to use.

In this work, we extend previous research in two directions.

First, we suggest to use the Threshold Accepting algorithm and show that this approach allows to minimize the loss in precision more effectively, more reliably, and more efficiently than Differential Evolution. I.e., TA finds partitions that have a smaller loss in precision than those found by DE. TA converges to better grouping solutions in less computational time and with a smaller number of iterations.

Second, we propose two different approaches for determining the optimal number of buckets. To our knowledge, this topic has not been addressed in the literature before although it is of great importance for practitioners. We aim to tackle the problem by designing a bank’s credit rating system such that its quality may be validated ex-post. The loss in precision by grouping borrowers together rather than treating them as individuals decreases as the number of buckets increases. Moreover, banks and regulatory authorities are concerned with stating regulatory capital (respectively unexpected losses) correctly. Thus, we propose to cluster borrowers such that we may evaluate ex post with a given confidence level whether actual unexpected losses fall within a sufficiently narrow interval around predicted unexpected losses. Then, the
optimal number of buckets is the maximum number of buckets that allows us to support our statement with a given confidence level. Our evaluations of this constraint suggest that it influences the structure of clusters in a reasonable way. Moreover, we find that even for small sample sizes it allows us to use up to eleven buckets for reasonable confidence- and precision-levels.

We show that our approach can provide meaningful insights into the problem of determining the optimal structure of PD buckets. However, we are aware that further research and empirical investigation on larger real-world datasets is required. Moreover, it is of special interest which confidence- and precision-levels may be used for different sample sizes. In this context, also different assumptions about the dependency structure of unexpected losses in a credit portfolio might be considered. Finally, although the constraint imposed on unexpected losses has a strong theoretical support, one might also consider alternative formulations or approximations resulting in a lower computational complexity for calculating the constraints. Thereby, the efficiency of the algorithm could be improved even further.

Appendix

Penalty Term

The exponent $a$ used in the penalty term (10) is defined as follows:
\[
a = \left( 0.5 \cdot \sum_b D_{EAD,b} \cdot \frac{\sum_{i \in b} EAD_{b,i} - 35\% \cdot \sum_b \sum_{i \in b} EAD_{b,i}}{65\% \cdot \sum_b \sum_{i \in b} EAD_{b,i}} \right) + \left( 0.5 \cdot \frac{\sum_b D_{N,b} \cdot \frac{x \cdot N - N_b}{x \cdot N}}{\sum_b D_{N,b}} \right).
\]

The idea of the penalty technique is to allow infeasible candidate solutions while running the algorithm as a stepping stone to get closer to promising regions of the search space. In this case, a penalty is multiplied on the objective function value that depends on the extent of constraint violations. In order to guarantee a feasible solution at the end, this penalty should increase over the runtime of the algorithm. The problem-specific penalty weights used in our application are defined by equations (10) and (19). They state that the objective function value \( f_u \) of a candidate solution is increased by some penalty factor \( A \in [1; 2] \) that puts more weight on penalties the more the current iteration \( i \) approaches the overall number of iterations \( I \). No penalty is placed on \( f_u \) if no constraint is violated so that \( a = 0 \). However, the variable \( a \) may take values up to 1 if the violation of the constraints reaches its maximum value. If the sum of \( EAD \) in some bucket \( b \) exceeds 35\% of total \( EAD D_{EAD,D} \) takes value 1 and 0 otherwise. \( D_{N,b} \) takes value 1 if bucket \( b \) contains less than 1\% of all borrowers. Both binding constraints are equally weighted.
Confidence Interval

Let us define the dummy variable $D_{N,b}$, which takes the value 1 if the constraint is violated and 0 otherwise. When constraints are considered based on rejection of infeasible candidate solutions the algorithms described above will not change. However, if the penalty technique is used it is necessary to alter equation (19) by removing the second summand in (19) and adding a term for the degree of violation of the additional constraint as exhibited by the second term in (20):

$$a = \left( 0.5 \cdot \sum_b \cdots \right) + \left( 0.5 \cdot \frac{\sum_b D_{N,b} \cdot \frac{1-\alpha - P_{int}}{1-\alpha}}{\sum_b D_{N,b}} \right).$$

Results with Ex Post Validation

In the following, the numerical results shall be presented that are discussed and interpreted in Section 5. In this section we evaluate the quality of the $UL$-constraint. The results were obtained from running TA 30 times with 200,000 iterations. We evaluate objective functions (5) based on squared differences of $PDcs$, (6) based on weighted squared differences of $PDcs$, and (7) based on differences in $RC$ in absolute terms. We choose $\alpha = 10\%$ and $\varepsilon = 30\%$. The last column gives the number of runs that converge towards the best solution relative to all runs that produce a solution meeting all constraints. For the problem instances, for which no feasible solution could be found in 30 runs, we report “n.a.” in the corresponding cells of the tables.
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^aRejection based constraint handling technique
^bPenalty technique
Table 7: Objective function (6) in EUR

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\(^a\) Rejection based constraint handling technique

\(^b\) Penalty technique
Table 8: Objective function (7) in EUR

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^aRejection based constraint handling technique
^bPenalty technique
References


45


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