A NUMERICAL COMPARISON OF SOME HEURISTIC STOPPING RULES FOR NONLINEAR LANDWEBER ITERATION*

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Abstract. The choice of a suitable regularization parameter is an important part of most regularization methods for inverse problems. In the absence of reliable estimates of the noise level, heuristic parameter choice rules can be used to accomplish this task. While they are already fairly well understood and tested in the case of linear problems, not much is known about their behaviour for nonlinear problems and even less in the respective case of iterative regularization. Hence, in this paper, we numerically study the performance of some of these rules when used to determine a stopping index for Landweber iteration for various nonlinear inverse problems. These are chosen from different practically relevant fields such as integral equations, parameter estimation, and tomography.

Key words. heuristic parameter choice rules, Landweber iteration, inverse and ill-posed problems, nonlinear operator equations, numerical comparison

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1. Introduction. In this paper, we consider nonlinear inverse problems of the form

\[(1.1) \quad F(x) = y,\]

where \(F : D(F) \subset X \rightarrow Y\) is a continuously Fréchet-differentiable nonlinear operator between real Hilbert spaces \(X\) and \(Y\). Furthermore, we assume that, instead of exact data \(y\), we are only given noisy data \(y^\delta\) that satisfy

\[(1.2) \quad \|y - y^\delta\| \leq \delta,\]

where \(\delta\) denotes the noise level. Typically, inverse problems are also ill-posed, which means that they may have no or even multiple solutions and, in particular, that a solution does not necessarily depend continuously on the data. This entails a number of difficulties due to which one usually has to regularize the problem.

During the last decades, a large number of different regularization approaches have been developed; see, e.g., \([22, 49, 81]\) and references therein. Two of the most popular methods, which also serve as the bases for a wide variety of other regularization approaches, are Tikhonov regularization \([83, 84]\) and Landweber iteration \([62]\). In its most basic form, Tikhonov regularization determines a stable approximation \(x_\alpha^\delta\) to the solution of \((1.1)\) as the minimizer of the Tikhonov functional

\[(1.3) \quad T_\alpha^\delta(x) := \|F(x) - y^\delta\|^2_Y + \alpha \|x - x_0\|^2_X,\]

where \(\alpha \geq 0\) is a regularization parameter and \(x_0\) is an initial guess. In order to obtain convergence of \(x_\alpha^\delta\) to a solution \(x_*\) of \((1.1)\), the regularization parameter \(\alpha\) has to be suitably chosen. If the noise level \(\delta\) from \((1.2)\) is known, then one can use either a priori parameter

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choice rules such as $\alpha \sim \delta$ or a posteriori rules like the discrepancy principle, which determines $\alpha$ as the solution of the nonlinear equation with some $\tau > 1$:

\[(1.4) \quad \|F(x^\delta_\alpha) - y^\delta\| = \tau \delta.\]

Unfortunately, in many practical applications, estimates of the noise level $\delta$ are either unavailable or unreliable, which renders the above parameter choice rules impracticable. Hence, a number of so-called heuristic parameter choice rules have been developed over the years; see, e.g., [35, 36, 64, 65, 80, 85, 86] and references therein. Most of them determine a regularization parameter $\alpha^*$ via

\[(1.5) \quad \alpha^* \in \arg\min_{\alpha \geq 0} \psi(\alpha, y^\delta),\]

where $\psi : \mathbb{R}_0^+ \times \mathbb{Y} \to \mathbb{R} \cup \{\infty\}$ is some lower semi-continuous functional. For example, the following popular choices in turn define the heuristic discrepancy (HD) principle [35], the Hanke–Raus (HR) rule [35, 79], the quasi-optimality (QO) rule [84], and the simple L (LS) rule [59]:

\[
\psi_{\text{HD}}(\alpha, y^\delta) := \frac{1}{\sqrt{\alpha}}\|F(x^\delta_\alpha) - y^\delta\|, \\
\psi_{\text{HR}}(\alpha, y^\delta) := \frac{1}{\alpha^2}(F(x^\delta_{\alpha,2}) - y^\delta, F(x^\delta_\alpha) - y^\delta), \\
\psi_{\text{QO}}(\alpha, y^\delta) := \|x^\delta_{\alpha,2} - x^\delta_\alpha\|, \\
\psi_{\text{LS}}(\alpha, y^\delta) := \langle x^\delta_\alpha - x^\delta_{\alpha,2}, x^\delta_\alpha \rangle,
\]

where $x^\delta_{\alpha,2}$ denotes the so-called second Tikhonov iterate [33], which is defined by

\[x^\delta_{\alpha,2} := \arg\min_{x \in \mathbb{X}} \{\|F(x) - y^\delta\|^2_\mathbb{Y} + \alpha\|x - x^\delta_\alpha\|^2_\mathbb{X}\}.\]

As is usually the case, rules of the form (1.5) are not restricted to Tikhonov regularization but can be used (possibly in a different form) in conjunction with various other regularization methods as well. Owing to their practical success in the treatment of linear inverse problems, heuristic parameter choice rules have received extended theoretical attention in recent years; see, e.g., [30, 31, 51, 52, 56, 79] and Section 2.2 for an overview.

A potential drawback of using Tikhonov regularization is the need to minimize the Tikhonov functional (1.3). Although for linear operators this reduces to solving a linear operator equation, the case of nonlinear operators is much more involved since one typically has to apply iterative optimization algorithms for the minimization of (1.3). Moreover, in order to employ either the discrepancy principle (1.4) or a heuristic rule of the form (1.5), this minimization usually has to be done repeatedly for many different values of $\alpha$, which can render it infeasible for many practical applications.

Hence, a popular alternative in order to circumvent these issues is to directly use so-called iterative regularization methods. As noted above, perhaps the most popular of these methods is Landweber iteration [22, 49], which is defined by

\[x^\delta_{k+1} = x^\delta_k + \omega F'(x^\delta_k)^* (y^\delta - F(x^\delta_k)),\]

where $\omega > 0$ is a step-size parameter. In order to obtain a convergent regularization method, Landweber iteration has to be combined with a suitable stopping rule such as the discrepancy principle, which now determines the stopping index $k_*$ by

\[(1.6) \quad k_* := \min\{k \in \mathbb{N} \mid \|F(x^\delta_k) - y^\delta\| \leq \tau \delta\},\]
for some parameter $\tau \geq 1$. Note that in contrast to (1.4), for the choice of $\alpha$ in Tikhonov regularization, the discrepancy principle for Landweber iteration can be verified directly during the iteration and does not require the iteration to be run more than once.

Now, analogously to (1.5), most heuristic parameter choice (stopping) rules for Landweber iteration determine a stopping index $k^*$ via

$$
\psi(k, y^\delta) := \psi^\text{HD}(k, y^\delta), \quad \psi(k, y^\delta) := \psi^\text{HR}(k, y^\delta), \quad \psi(k, y^\delta) := \psi^\text{QO}(k, y^\delta), \quad \psi(k, y^\delta) := \psi^\text{LS}(k, y^\delta),
$$

with $\psi : \mathbb{N} \times Y \to \mathbb{R} \cup \{\infty\}$ again being some lower semi-continuous functional [35, 51, 52, 56, 75]. At least conceptually, the regularization parameter $\alpha^*$ in Tikhonov regularization and the stopping index $k^*$ for Landweber iteration play inversely proportional roles, i.e., $\alpha \sim 1/k$.

Hence, the heuristic rules from (1.5) now correspond to

$$
\psi_{\text{HD}}(k, y^\delta) := \sqrt{k} \|F(x^\delta_k) - y^\delta\|,
\psi_{\text{HR}}(k, y^\delta) := k(y^\delta - F(x^{2k}_k), y^\delta - F(x^\delta_k)),
\psi_{\text{QO}}(k, y^\delta) := \|x^{\delta}_{2k} - x^\delta_k\|,
\psi_{\text{LS}}(k, y^\delta) := \langle x^{\delta}_k, x^{\delta}_{2k} - x^\delta_k \rangle.
$$

Note that the analogue of the second Tikhonov iteration $x^{\delta}_{\alpha,2}$ in the Landweber method corresponds to applying $k$ steps of a Landweber iteration with initial guess $x^\delta_k$, which is easily seen to be identical to simply doubling the iteration number, i.e., calculating $x^\delta_{2k}$.

Several variants and modifications of the rules in (1.8) are possible: The functionals $\psi(k, \delta)$ can be divided by $\|x^\delta_k\|$, leading to the so-called ratio methods; see, e.g., [59] for the simple L rule. Furthermore, the factor $k$ or $\sqrt{k}$ for the first two instances can be adapted in case of variable step sizes (see [48]) or it can be replaced by a shifted version $k + c$ or $\sqrt{k + c}$, respectively; see, e.g., [35, 48]. The latter only changes the functional values by a factor $O(1)$ and seems to yield no significant advantages at the cost of an additional constant $c$ to be chosen.

Similarly to the case of the discrepancy principle (1.6), these functionals can be evaluated during a single run of Landweber iteration. Hence, in particular in the nonlinear case, the combination of heuristic parameter choice rules together with Landweber iteration (or other iterative regularization methods) suggests itself.

Heuristic parameter choice rules are already fairly well understood in the case of linear problems. Despite a number of obstacles such as the Bakushinskii veto [6], many theoretical results on both the convergence and other aspects of these rules are already available; see, e.g., [51] and references therein. Furthermore, numerical tests have been carried out for various different linear test problems [14, 31, 75]. In contrast, for the case of nonlinear problems, not much is known with respect to convergence theory nor is there a unifying framework for heuristic rules, and, furthermore, there are no numerical performance studies for heuristic rules.

Hence, the main motivation of this article is to close this gap. In particular, we consider the behaviour of the four rules given in (1.8), i.e., the heuristic discrepancy principle, the Hanke–Raus rule, the quasi-optimality rule, and the simple L rule, using Landweber iteration and for different nonlinear test problems from practically relevant fields such as integral equations, parameter estimation, and tomography. The aim of this paper is two-fold: On the one hand, we want to demonstrate that heuristic parameter choice rules can indeed be used successfully not only for linear but also for nonlinear inverse problems. On the other hand, we want to provide some useful insight into potential difficulties and pitfalls that one might encounter, and what can be done about them. Since we also compare the heuristics with the
discrepancy principle, the results in this article additionally provide a numerical study of the performance of the latter in the nonlinear case.

The outline of this paper is as follows: In Section 2, we provide some general background on heuristic parameter choice rules and nonlinear Landweber iteration. In Section 3, we then introduce the various problems on which we want to test the different heuristic parameter choice rules. The corresponding results are then presented in Section 4, which is followed by a short conclusion in Section 5.

2. Landweber iteration and stopping rules. In this section, we recall some basic results on Landweber iteration and heuristic parameter choice rules. Since in this paper we are mainly interested in a numerical comparison of the different rules, we only provide a general overview of some of the main results from the literature, focusing in particular on those aspects which are relevant to understand the numerical results presented below.

2.1. Landweber iteration for nonlinear problems. One of the main differences from the linear case is that for nonlinear problems only local convergence can be established. For this, one needs to impose certain restrictions on the nonlinearity of the operator $F$, such as the tangential cone condition [34, 49], which is given by

\begin{equation}
\|F(x) - F(\tilde{x}) - F'(x)(x - \tilde{x})\| \leq \eta \|F(x) - F(\tilde{x})\|, \quad \eta < 1/2.
\end{equation}

This condition has to hold locally in a neighbourhood of a solution $x^*$ of the problem, and the initial guess $x_0$ of the iteration has to be contained inside of it. Furthermore, the parameter $\tau$ in the discrepancy principle (1.6) has to satisfy

\begin{equation}
\tau > 2 \frac{1 + \eta}{1 - 2\eta} \geq 2,
\end{equation}

where the factor 2 can be slightly improved by an expression depending on $\eta$ which tends to 1 as $\eta \to 0$, thereby recovering the optimal bound in the linear case [32].

If in addition the step size $\omega$ is chosen small enough such that locally

\begin{equation}
\omega \|F'(x)\|^2 \leq 1
\end{equation}

holds, then Landweber iteration combined with the discrepancy principle (1.6) converges to $x^*$ as the noise level $\delta$ goes to 0 [34, 49]. Furthermore, convergence to the minimum-norm solution $x^\dagger$ can be established given that, in a sufficiently large neighbourhood, it holds that $N(F'(x^\dagger)) \subset N(F'(x))$. In order to prove convergence rates, in addition to source conditions, further restrictions on the nonlinearity of $F$ are necessary [22, 49].

Even though the tangential cone condition (2.1) holds for a number of different applications (see, e.g., [50] and references therein) and even though attempts have been made to replace it by more general conditions [53], these can still be difficult to prove for specific applications (cf., e.g., [54] for an analysis of the electrical impedance tomography (EIT) problem and [42] for an analysis of a parameter estimation problem in linear elastography). Furthermore, even if the tangential cone condition can be proven, the exact value of $\eta$ typically remains unknown. Since this also renders condition (1.2) impracticable, the parameter $\tau$ in the discrepancy principle then has to be chosen manually; popular choices include $\tau = 1.1$ or $\tau = 2$. These work well in many situations but are also known to fail in others (compare with Section 4 below). In any case, this shows that, for practical applications involving nonlinear operators, informed “heuristic” parameter choices remain necessary even if the noise level $\delta$ is known.
2.2. Heuristic stopping rules. As mentioned in the introduction, in many practical situations one does not have knowledge of the noise level. Thus, applying it with unreliable estimates of the noise level is rarely fruitful. The remedy is the use of heuristic (also known as data-driven or error-free) rules, where the iteration is terminated at \( k_* := \hat{k}(y^\delta) \), which depends only on the measured data and not on the noise level.

The analysis of the so-called heuristic rules for Tikhonov regularization has been examined extensively for linear problems [51, 52, 56] but less so beyond the linear case. Some results for convex problems and nonlinear problems in Banach spaces can be found in [46, 47, 48]. The numerical performance of these rules in the linear case has also been studied [14, 31, 75] and for convex Tikhonov regularization in [55]. However, for Landweber iteration for nonlinear problems, neither has an analysis been given nor has the numerical performance of heuristic stopping rules been investigated so far.

Let us briefly illustrate the rationale behind the heuristic minimization-based rules (1.7) (or (1.5)). It is well known that the total error between the regularized solution \( x_k^\delta \) and the exact solution \( x^\dagger \) can be split into approximation and stability error:

\[
\|x_k^\delta - x^\dagger\| \leq \|x_k - x^\dagger\| + \|x_k - x_k^\delta\|. \tag{2.4}
\]

Here, \( x_k \) denotes the regularized solution when exact data \( (\delta = 0) \) would be given. An ideal optimal parameter choice would be one that minimizes over \( k \) the total error or the upper bound in (2.4). However, this “oracle” parameter choice is not possible, as in (2.4) only the element \( x_k^\delta \) is at hand, and neither the exact solution nor the exact data are available. The idea of minimization-based heuristic rules is to construct a computable functional \( \psi \) that estimates the total error sufficiently well, i.e., \( \psi(k, y^\delta) \sim \|x_k^\delta - x^\dagger\| \), such that a minimization over \( k \) is expected to yield a reasonable parameter choice of \( k_* \) as well. In analogy to the above splitting of approximation and stability, we may use a similar splitting for \( \psi \), i.e., \( \psi(k, y^\delta) \leq \psi_a(k, y^\delta) + \psi_d(k, y^\delta) \), and search for conditions such that the respective parts estimate the approximation and stability error separately:

\[
\psi_a(k, y^\delta) \sim \|x_k - x^\dagger\|, \tag{2.5}
\]
\[
\psi_d(k, y^\delta) \sim \|x_k - x_k^\delta\|. \tag{2.6}
\]

As briefly mentioned earlier, the pitfall for heuristic stopping rules manifests itself in the form of the so-called Bakushinskii veto, the consequence of which is that a heuristic stopping rule cannot yield a convergent regularization scheme in the worst-case scenario [6], i.e., for all possible noise elements \( y = y^\delta \). A direct consequence is that there cannot exist a \( \psi \) with the error-estimating capabilities as mentioned above, in particular, such that (2.6) holds! Essentially, the main reason for the worst-case failure of heuristic rules is that they cannot discriminate between exact data and noisy data with “smooth noise”; cf., e.g., [58].

However, there is a way to overcome the negative result of Bakushinskii by restricting the class of permissible noise elements \( y = y^\delta \). In this way, one can prove convergence in a restricted noise case scenario. (Of course, for this approach to be meaningful, the restrictions should be such that “realistic” noise is always permitted.) Some noise restriction were used, e.g., in [27, 35], although the restrictions there were implicit and very hard to interpret. (For instance, in [27], essentially, condition (2.6) was postulated rather than derived from more lucid conditions.)

A major step towards an understanding of heuristic rules was made in [51, 56], when a full convergence analysis in the linear case with explicit and interpretable restrictions was given. These conditions, which were proven to imply (2.6), take the form of a Muckenhoupt-type inequality: Let \( (\sigma_i, u_i, v_i) \) be the singular system of the forward operator. Then, for
\( p \in \{1, 2\} \), the \( p \)-Muckenhoupt inequality holds if there is a constant \( C \) and a \( t_0 \) such that, for all admissible noise \( y - y^\delta \), it holds that

\[
\sum_{\sigma_i^2 \geq t} \frac{t}{\sigma_i^2} |(y - y^\delta, v_i)|^2 \leq C \sum_{\sigma_i^2 < t} \left( \frac{\sigma_i^2}{t} \right)^{p-1} |(y - y^\delta, v_i)|^2 \quad \forall \ t \in (0, t_0).
\]

It has been shown in [51, 56, 59] that, for most of the classical regularization schemes and for the four above-mentioned rules, this condition suffices for being able to estimate the stability error (2.6), and thus convergence can be proven. The inequality (2.7) has the interpretation of an “irregularity” condition for the noise vector \( y - y^\delta \); by postulating (2.7), the noise must be distinguishable from a smooth data error (which never satisfies (2.7)). However, this anyway agrees with the common idea of noise.

**Remark 2.1.** The above heuristic rules require slightly different Muckenhoupt conditions, which lead to two groups of rules: For HD and HR, (2.7) with \( p = 1 \) suffices; while for QO and LS, the condition with \( p = 2 \) (which is a slightly stronger requirement) has to be postulated [51, 56, 59]. Thus the former might be successful even if the latter fail. However, it has to be kept in mind that the error analysis shows that, as long as they can be successfully applied, QO and LS in general lead to smaller errors.

Indeed, it has been shown in [56] that the above-mentioned Muckenhoupt inequality is satisfied in typical situations, and in [57] it was shown in a stochastic setting that it is also satisfied for coloured Gaussian noise almost surely for many cases. Below, we discuss cases where the Muckenhoup inequality might not be satisfied and when heuristic rules may fail.

The above-mentioned noise restrictions are heavily rooted in the linear theory and in particular make use of spectral theory and the functional calculus of operators. In the case of a nonlinear operator, we are no longer afforded the luxury of having these tools available. Some alternative noise conditions in the nonlinear case have been established in [46] for convex variational Tikhonov regularization, in [89] for Bregman iteration in Banach spaces, and in [66] for general variational regularization. However, as yet, these conditions could not be deciphered into a palatable explanation as to when a rule will work or otherwise. An attempt was made in [58] to formulate an analogous Muckenhoupt-type inequality for convex Tikhonov regularization in a somewhat restrictive setting of a diagonal operator over \( \ell^p \) spaces with \( p \in (1, \infty) \).

However, to the knowledge of the authors, neither a convergence analysis nor a numerical investigation of heuristic rules for nonlinear Landweber iteration seems to be available in the literature. In light of all of this, we thus have great incentive to investigate heuristic stopping rules for nonlinear Landweber iteration numerically and to compare them with the more tried and tested a posteriori rules.

The Muckenhoupt inequality covers the convergence theory. When it comes to the practical capabilities of heuristic rules, also convergence rates are important. For this, efficient estimates of the approximation error as in (2.5) are vital, and, in the linear case, sufficient conditions for this have been established, this time in the form of conditions for the exact solution \( x^\dagger \). For instance, using the singular system \((\sigma_i, u_i, v_i)\), the regularity condition (for \( p = 1 \) or \( p = 2 \), depending on the rule)

\[
\sum_{\sigma_i^2 \geq t} |(x^\dagger, u_i)|^2 \leq C \sum_{\sigma_i^2 > t} \left( \frac{\sigma_i^2}{t} \right)^{p-1} r(\sigma_i^2, \alpha)^2 |(x^\dagger, u_i)|^2 \quad \forall \ t \in (0, t_0),
\]

is sufficient for (2.5) and, together with smoothness conditions, yields (optimal-order) convergence rates in many situations [51]. Here, \( r(\lambda, \alpha) \) is the residual spectral filter function.
of the regularization method, i.e., \( r(\lambda, \alpha) = \alpha/(\alpha + \lambda) \) for Tikhonov regularization and \( r(\lambda, \alpha) = (1 - \lambda)^{1/\alpha} \) with \( \alpha = k^{-1} \) for Landweber iteration. Note that the regularity condition depends strongly on the regularization method via \( r(\lambda, \alpha) \), in contrast to the Muckenhoupt condition. The rough interpretation of (2.8) is that the exact solution has coefficients \((x^\dagger, u_i)\) that do not deviate too much from a given decay (that is encoded in a smoothness condition).

2.3. Challenges and practical issues for heuristic rules. Next, let us point out possible sources of failure for heuristic rules and some peculiarities for the case of nonlinear Landweber regularization.

**Failure of Muckenhoupt condition.** A general problem for heuristic rules, in both the linear and nonlinear cases, is when the Muckenhoupt condition (2.7) is not satisfied. This can happen for standard noise for super-exponential ill-posed problems, for instance, for the backward heat equation (see [51, 52]). Less obvious but practically important is the case that the Muckenhoupt inequality might also fail to hold for standard noise if the problem is nearly well-posed, i.e., when the singular values decay quite slowly (e.g., as \( \sigma_i \sim i^{-\beta} \) with, say, \( \beta < 1 \)). In this case, exact data can be quite irregular as the operator is only little smoothing, and it is hard to distinguish between exact data and noise, and this is indeed a relevant possible source of failure.

**Spurious first local minimum for Landweber iteration.** Recall that the effective performance of heuristic rules depends also on efficient estimates of the approximation errors and in particular on the regularity condition to hold. As pointed out before, this strongly depends on the regularization method. We have extensive numerical evidence that, for linear and nonlinear Landweber iteration, the approximation error (i.e., (2.5)) is often only badly estimated by \( \psi(k, y^\beta) \) for the first few iterations (i.e., \( k \) small) and that (2.8) holds only with a bad constant for large \( t = k^{-1} \). In practical computations, this has the consequence that \( \psi(k, y^\beta) \) typically has an outstanding local minimum for small \( k \). However, this local minimum is rarely the global minimum, which usually appears much later for larger \( k \), and inexperienced users are often tempted to mistakenly take this local minimum for the global one to save having to compute later iterations. This happens quite often in the linear and the nonlinear cases for Landweber iteration, but a similar problem for Tikhonov regularization is rarely observed. The deeper reason for this discrepancy is the different shape of the residual filter function for the two methods, which makes the regularity condition (2.8) more restrictive for Landweber iteration and for large \( t \).

**Discretization cut-off.** It is known that, due to discretization, the theoretically global minimum of \( \psi(k, y^\beta) \) for finite-dimensional problems is at \( k = \infty \), which does not provide a correct stopping index. Thus, in practical computations, we have to restrict the search space by fixing an upper bound for \( k \) (or, for the continuous regularization method, by a lower bound for \( \alpha \)). Some rules as to how to do this together with an accompanying analysis in the linear case are given in [52]; however, for nonlinear problems, no such investigation exists. This issue is relevant for very small noise levels or for coarse discretizations, and in practice one takes a pragmatic approach and assumes a reasonable upper bound for the iteration index and looks for interior minima rather than global minima at the boundary of the search space.

**Only local convergence in the nonlinear case.** The established convergence theory in the nonlinear case is a local one. One can only prove convergence when the initial guess is sufficiently close to the exact solution \( x^\dagger \), and, in the case when noise is present, the iteration usually diverges out of the neighborhood of \( x^\dagger \) as \( k \to \infty \). In particular, it is possible that \( x^\dagger_k \) “falls” out of the domain of the forward operator. As a consequence, it might happen that the functionals \( \psi(k, y^\beta) \) in (1.7) are not defined for very large \( k \). By definition, however, one would have to compute a minimizer over all \( k \), which is then not practically possible. (This is different from Tikhonov regularization, whose solution is always well defined for any \( \alpha \).)
In practice, as a remedy, one would introduce an upper limit for the number of iterations up to which the functional $\psi(k, y^\delta)$ is computed. Additionally, one could monitor the distance to the initial guess $\|x_k^\delta - x_0\|$ and terminate if this becomes too large.

In this section, we have stated some practical aspects of heuristic rules; a deeper mathematical analysis (especially in the nonlinear case) is outside the scope of this article. For further aspects on heuristic stopping rules from both theoretical and practical viewpoints, we refer the reader to [76] and references therein.

3. Test problems. In this section, we introduce a number of test problems on which we evaluate the performance of the heuristic stopping rules described above. These nonlinear inverse problems belong to a variety of different problem classes, including integral equations, tomography, and parameter estimation. For each of them, we briefly review their background and describe their precise mathematical setting and relevant theoretical results below.

3.1. Nonlinear Hammerstein operator. A commonly used nonlinear inverse problem [34, 40, 41, 70, 71, 72] for testing, in particular, the behaviour of iterative regularization methods is based on so-called nonlinear Hammerstein operators of the form

$$F : H^1[0, 1] \to L^2[0, 1], \quad F(x)(s) := \int_0^1 k(s, t)\gamma(x(t)) \, dt,$$

with some given function $\gamma : \mathbb{R} \to \mathbb{R}$. Here, we look at a special instance of this operator, namely

$$F(x)(s) := \int_0^s x(t)^3 \, dt,$$

for which the tangential cone condition (2.1) holds locally around a solution $x^\dagger$, given that it is bounded away from zero (see, e.g., [72]). Furthermore, the Fréchet derivative and its adjoint, which are required for the implementation of Landweber iteration, can be computed explicitly.

3.2. Diffusion coefficient estimation. Another classic test problem [49] in inverse problems is the estimation of the diffusion coefficient $a$ in the partial differential equation (PDE)

$$-\text{div}(a \nabla u) = f$$

from measurements of $u$ and given knowledge of the source term $f$ and (Dirichlet) boundary conditions on $u$. For this test problem, we focus on the one-dimensional version

$$-(a(s)u(s))_s = f(s), \quad s \in (0, 1),$$

$$u(0) = u(1) = 0,$$

which leads to an inverse problem of the form (1.1) with the nonlinear operator

$$F : D(F) := \{ a \in H^1[0, 1] \mid a(s) \geq a > 0 \} \to L^2[0, 1],$$

$$a \mapsto F(a) := u(a),$$

where $u(a)$ is the solution of (3.2) above. The computation of the Fréchet derivative and its adjoint of $F$ now requires solving PDEs of the form (3.2). Furthermore, it was shown (see, e.g., [49]) that the tangential cone condition (2.1) holds locally around a solution $a^\dagger \geq c > 0$. 
3.3. Linear elastography. A practically relevant nonlinear inverse problem of parameter estimation type appears in linear elastography. Therein, one tries to estimate the Lamé parameters $\lambda$ and $\mu$ from measurements of the displacement field $u$ inside an object (obtained, for example, through optical coherence tomography measurements) that has been subjected to a deformation \[8, 10, 11, 12, 13, 20, 21, 23, 28, 37, 44, 45, 60, 63, 67, 73, 74, 87\]. Under the assumption of a small deformation and an elastic material, the following PDE can be used as a model:

$$-\text{div}(\sigma(u)) = f, \quad \text{in } \Omega,$$

$$u|_{\Gamma_D} = g_D,$$

$$\sigma(u)\vec{n}|_{\Gamma_T} = g_T.$$  

Thereby, the body forces $f$, the displacement $g_D$, and the surface traction $g_T$ are assumed to be known, and the boundary $\partial\Omega$ with outward unit normal vector $\vec{n}$ of the domain $\Omega$ is split into a Dirichlet part $\Gamma_D$ and a Neumann part $\Gamma_N$, with the stress tensor $\sigma$ encoding the stress–strain relation in $\Omega$ being defined by

$$\sigma(u) := \lambda \text{div}(u)I + 2\mu \mathcal{E}u, \quad \mathcal{E}u := \frac{1}{2}(\nabla u + \nabla u^T),$$

where $I$ denotes the identity matrix and $\mathcal{E}$ is the so-called strain tensor. Again, it is possible to define a nonlinear operator $F$ such that this problem can be written in the form (1.1) and to compute its Fréchet derivative and the adjoint thereof. Furthermore, it was shown that the tangential cone condition (2.1) holds locally around a solution $(\lambda^\dagger, \mu^\dagger) \geq c > 0$ given that its values are known in an arbitrarily small region around the boundary (for details, we refer the reader to [42]), which is reasonable in many practical cases.

3.4. Acousto-electrical tomography. A PDE parameter estimation problem, this time from the field of tomography, is the hybrid imaging modality of acousto-electrical tomography (AET) \[4, 25, 61, 88\]. Based on a modulation of EIT by ultrasound waves, AET aims at reconstructing the spatially varying electrical conductivity distribution inside an object from electrostatic measurements of voltages and the corresponding current fluxes on its surface. Compared, for example, to EIT, reconstructions of high contrast and high resolution may be obtained. Mathematically, the problem amounts to reconstructing the spatially varying conductivity $\sigma$ from measurements of the power densities

$$E_j(\sigma) := \sigma|u_j(\sigma)|^2,$$

where the interior voltage potentials $u_j(\sigma)$ are the solution of the elliptic PDEs

$$\text{div}(\sigma \nabla u_j) = 0, \quad \text{in } \Omega,$$

$$\sigma \nabla u_j \cdot \vec{n}|_{\partial\Omega} = g_j,$$  

(3.4)

where $\Omega \subset \mathbb{R}^N$, with $N = 2, 3$, is a bounded and smooth domain and $g_j$ models the current flux on the boundary $\partial\Omega$ in the outward unit normal direction $\vec{n}$.

Once again, this problem can be restated as an operator equation of the form (1.1) with a Fréchet-differentiable nonlinear operator. Its Fréchet derivative and the adjoint thereof can be found, for example, in [39], and their evaluation again requires the solution of PDEs of the form (3.4) for different right-hand sides. Note that it is not known whether the tangential cone condition holds for AET (or EIT). Furthermore, it is in general not possible to uniquely determine the conductivity $\sigma$ from a single power density measurement $E_j(\sigma)$ \[7, 43\]. In addition, if $g_j = 0$ on some part $\Gamma \subset \partial\Omega$ of the boundary, then the problem becomes severely...
ill-posed. On the other hand, if \( g_j \neq 0 \) almost everywhere on \( \partial \Omega \), and given a sufficient amount (depending on the dimension \( N \)) of “different” power density measurements \( E_j(\sigma) \), the conductivity \( \sigma \) can be uniquely reconstructed \([2, 9, 17, 68]\). In this case, the problem behaves numerically close to well-posed, which is reflected in the behaviour of the heuristic parameter choice rules; cf. Section 2.3.

3.5. Single photon emission computed tomography. Next, we look at single photon emission computed tomography (SPECT), which is another example from the large field of tomography \([18, 19, 69, 77, 78]\). In this medical imaging problem, one aims at reconstructing the radioactive distribution \( f \) (activity function) and the attenuation map \( \mu \), which is related to the density of different tissues, from radiation measurements outside the examined body. The connection between these quantities is typically modelled by the attenuated radon transform (ART), which is given by \([69]\)

\[
F(f, \mu)(s, \omega) := \int_{\mathbb{R}} f(s\omega^+ + tw) \exp \left( -\int_t^\infty \mu(s\omega^+ + r\omega) \, dr \right) \, dt,
\]

where \( s \in \mathbb{R} \) and \( \omega \in S^1 \). With this, one again arrives at a problem of the form (1.1), where \( y \) then is the measured sinogram. The well-definedness and differentiability of the operator \( F \) with respect to suitable Sobolev spaces has been studied in detail in \([18, 19]\). However, it is still unknown whether also the tangential cone condition (2.1) holds for this problem.

3.6. Auto-convolution. As a final test example, we now consider the problem of (de-)auto-convolution \([16, 24, 29, 77]\). Among the many inverse problems based on integral operators, auto-convolution is particularly interesting due to its importance in laser optics \([5, 15, 26]\). Mathematically, it amounts to solving an operator equation of the form (1.1) with the operator

\[
F : L^2[0, 1] \to L^2[0, 1], \quad F(x)(s) := (x * x)(s) := \int_0^1 x(s-t)x(t) \, dt,
\]

where the functions in \( L^2[0, 1] \) are interpreted as 1-periodic functions on \( \mathbb{R} \). While deriving the Fréchet-differentiability and its adjoint for \( F \) is straightforward, it is not known whether the tangential cone condition (2.1) holds. However, for small enough noise levels \( \delta \), the residual functional is locally convex around the exact solution \( x^\dagger \), given that it has only finitely many non-zero Fourier coefficients \([41]\).

4. Numerical results. In this section, we present the results of using the four heuristic parameter choice rules defined in (1.8) (as well as the classical discrepancy principle for comparison) to determine a stopping index for Landweber iteration applied to the different nonlinear test problems introduced in Section 3.

For each of these problems, we started from a known solution \( x^\dagger \) in order to define the exact right-hand side \( y \). Random noise corresponding to different noise levels \( \delta \) was added to \( y \) in order to create noisy data \( y^\delta \), and a suitable step size \( \omega \) for Landweber iteration was computed via (2.3) based on numerical estimates of \( \|F'(x^\dagger)\| \). Afterwards, we ran Landweber iteration for a predefined number of iterations \( k_{\text{max}} \), which was chosen manually for each problem via a visual inspection of the error, residual, and heuristic functionals, such that all important features of the parameter choice rules were captured for this comparison.

Following each application of Landweber iteration, we computed the values of the heuristic functionals \( \psi \), as well as their corresponding minimizers \( k_\cdot \). As noted in Section 2, the functional values corresponding to the first few iterations have to be discarded in the search for the minimizers due to the spurious first local minimum (tacitly assuming that the noise
level is small enough such that a good stopping index appears later). For each of the different heuristic rules, we then computed the resulting absolute error

\[ \| x_{k}^\delta - x^\dagger \|, \]

and, for comparison, for each problem we also computed the optimal stopping index

\[ k_{\text{opt}} := \arg\min_{k \in \mathbb{N}} \| x_{k}^\delta - x^\dagger \|, \]

together with the corresponding optimal absolute error. Furthermore, we also computed the stopping index \( k_{\text{DP}} \) determined by the discrepancy principle (1.6), which can also be interpreted as the “first” minimizer of the functional

\[ \psi_{\text{DP}}(k) := \| F(x_{k}^\delta) - y^\delta \| - \tau \delta. \]

As noted in Section 2.1, since the exact value of \( \eta \) in (2.1) is unknown for our test problems, a suitable value for \( \tau \) has to be chosen manually. Depending on the problem, we used one of the two popular choices, \( \tau = 1 \) or \( \tau = 2 \), although, as we are going to see below, these are not necessarily the “optimal” ones. In any case, the corresponding results are useful reference points to the performance of the different heuristic parameter choice rules.

Concerning the discretization and implementation of each of the numerical test problems, we refer to the subsequent sections and the references mentioned therein. All computations were carried out in Matlab on a notebook computer with an Intel® Core™ i7-85650 processor with 1.80 GHz (eight cores) and 16 GB RAM, except for the acousto-electrical tomography problem, which was carried out in Python using the FEniCS library [3] on a notebook computer with an Intel® Core™ i7-4810MQ processor with 2.80 GHz (eight cores) and 15.3 GB RAM.

4.1. Nonlinear Hammerstein operator. First, we consider the nonlinear Hammerstein problem introduced in Section 3.1. In order to discretize this problem, the interval \([0, 1]\) is subdivided into 128 subintervals, and the operator \( F \) itself is discretized as described in [70, 72]; cf. also [38, 40]. For the exact solution, we choose \( x^\dagger(s) = 2 + (s - 0.5)/10 \) and compute the corresponding data \( y \) by an application of the Hammerstein operator (3.1). For the initial guess we choose \( x_0(s) = 1 \), and in the discrepancy principle (1.6) we use \( \tau = 2 \). The absolute error (4.1) corresponding to different parameter choice rules and noise levels \( \delta \) from 0.1% to 2% is depicted in Figure 4.1. Typical plots of the heuristic functionals \( \psi \) as well as the evolution of the absolute error over the iteration are depicted in Figure 4.2. There, the marked points denote the corresponding stopping indices selected via the different rules.

As can be seen from the left plot in Figure 4.2, the heuristic functionals \( \psi \) generally exhibit the same shape as expected from the theoretical considerations discussed above. For example, each of the functionals exhibits a spurious “first” local minimum within the first few iterations, as already discussed in Section 2.3. Apart from this, each functional \( \psi \) has a well-defined minimum reasonably close to the stopping index \( k_{\text{DP}} \) determined by the discrepancy principle. However, for larger noise levels, this minimum vanishes for the HD and the HR rules, which is reflected in Figure 4.1 by their unsatisfactory constant absolute error (the rules select the spurious minimum in this case). In contrast, the QO and LS rules keep their general shape for all noise levels and thus produce stable stopping indices, which are typically larger than those determined by the discrepancy principle. Since the evolution of the absolute error depicted in the right plot in Figure 4.2 flattens for larger iteration numbers (an effect of the discretization), the error curves for the QO and the LS rules remain rather constant on the logarithmic scale. Curiously, and contrary to theoretical expectations, the absolute error curve for the discrepancy principle in Figure 4.1 exhibits a parabolic shape. This indicates that, on the one hand, the
chosen value of $\tau$ is too small, while, on the other hand, the discretization might be too coarse in comparison with the small noise levels. However, in practice, the discretization is often fixed by practical limitations, while a proper value of $\tau$ satisfying (2.2) is typically impossible to determine, or unreasonably large. Hence, this first test already indicates the usefulness of heuristic parameter choice rules (especially the QO and LS rules) in comparison to the discrepancy principle and shows some typical limitations that we now investigate further in the remaining test problems.
4.2. Diffusion coefficient estimation. Next, we consider the diffusion coefficient estimation problem introduced in Section 3.2. For discretizing the problem, we use a standard projection approach (see, e.g., [22]) onto a finite-dimensional subspace of $H^1[0, 1]$ spanned by piecewise linear finite element method (FEM) hat functions defined on a uniform subdivision of $[0, 1]$ into 50 subintervals. For the exact solution, we choose $x^*(s) = 2 + s(1 - s)$ and compute the corresponding data $y$ by applying the operator $F$ defined in (3.3), using a finer grid in order to avoid an inverse crime. For the initial guess, we use $x_0(s) = 2$.

According to the discrepancy principle (1.6) we choose $\tau = 1.1$. As before, Figure 4.3 depicts the absolute errors (4.1) corresponding to the different parameter choice rules, now for noise levels $\delta$ between $0.1\%$ and $1\%$. A typical plot of the heuristic functionals $\psi$ and the evolution of the absolute error can be seen in Figure 4.4.

![Figure 4.3](image_url)

**FIG. 4.3.** Numerical results for the diffusion coefficient estimation problem introduced in Section 3.2. Absolute error (4.1) at the stopping indices $k_\ast$ determined by the discrepancy principle (1.6), the different heuristic parameter choice rules (1.8), and the optimal stopping index $k_{opt}$ defined in (4.2), each for different relative noise levels $\delta$.

We observe again a slight superiority of the QO and LS rules over the HD and HR methods and even over the discrepancy principle, although the rates of these methods (the slopes of the plots in Figure 4.3) are comparable. The HD and HR methods indicate erratic behaviour for large $\delta$, which is explained by the lack of a clear minimum and the resulting suboptimal choice of the spurious minimum. While the discrepancy principle follows quite nicely a theoretically predicted rate, the QO rule (and less so the LS rule) has a jump in the error curve, which is explained by the occurrence of two local minima in the graph of Figure 4.4; at a certain $\delta$ the global minimum switches from one to the other. As before, we can conclude successful results for the discrepancy principle as well as for the heuristic rules.

4.3. Linear elastography. For the third test, we consider the linear elastography problem introduced in Section 3.3. For the exact description of the problem set-up, discretization, and implementation, we refer to [42]. In short, the unknown Lamé parameters $(\lambda^\dagger, \mu^\dagger)$ are uniform circular inclusions in an otherwise constant background over a rectangular domain $\Omega \subset \mathbb{R}^2$. Constant displacement conditions ($g_D = \text{const}$.) are applied on the top and bottom of the domain, while on the sides we impose vanishing surface traction corresponding to
Heuristic stopping rules for Landweber iteration

Fig. 4.4. Numerical results for the diffusion coefficient estimation problem introduced in Section 3.2. Heuristic functionals (1.8) and discrepancy functional (4.3) for $\delta = 0.2\%$ relative noise (left). Corresponding evolution of the absolute error $\|x_k^\delta - x^\dagger\|$, with marked points indicating the stopping indices chosen by the different rules (right).

$g_T = 0$. Again, Figure 4.5 depicts the absolute errors (4.1) corresponding to the different parameter choice rules, this time for noise levels $\delta$ between 0.1% and 1%. A typical plot of the heuristic functionals and the corresponding evolution of the absolute errors (4.1) can be found in Figure 4.6.

First of all, we observe that, for this problem, the LS rule performs very well over the entire range of noise levels, with a clearly defined minimum in the functionals $\psi$. On the other
hand, both the HD and the HR rules fail entirely, due to their corresponding functionals $\psi$ increasing monotonically. The discrepancy principle, here used with the choice of $\tau = 4$, also does not perform well for larger noise levels, although improving for smaller values of $\delta$. This indicates that our conservative choice of $\tau$ is not optimal, which, on the other hand, indicates the importance of a “proper” parameter choice even for deterministic stopping rules. Finally, note that, for large noise levels $\delta$, also the QO rule fails since the corresponding functionals $\psi$ are monotonically increasing. However, this increase stalls for smaller noise levels, until the functionals $\psi$ begin to exhibit clear local minima from $\delta = 0.5\%$ onwards, which results in an excellent performance of the QO rule for smaller noise levels.

4.4. Acousto-electrical tomography. Next, we consider the AET problem introduced in Section 3.4. For a detailed description of the problem set-up, discretization, and implementation, we refer to [39]. In short, the unknown inclusion $\sigma^\dagger$ consists of three uniform disconnected inclusions (two circular, one crescent-shaped) with values of 1.3, 1.7, and 2, respectively, in an otherwise constant background of value 1 over the circular domain $\Omega := \{(r, \theta) \in [0, 1) \times [0, 2\pi]\} \subset \mathbb{R}^2$. Furthermore, we use the boundary flux functions

$$g_j(r, \theta) := \begin{cases} \sin(2j\pi\theta/\alpha), & (r, \theta) \in \Gamma(\alpha), \\ 0, & \text{otherwise}, \end{cases}$$

for $j = 1, 2, 3$,

where $\Gamma(\alpha) := \{(r, \theta) \in \{1\} \times [0, \alpha]\} \subset \partial\Omega$ for $\alpha \in [0, 2\pi]$. Hence, if $\alpha = 2\pi$, then $g_j \neq 0$ almost everywhere on $\partial\Omega$. In the following, this case will be called 100\% boundary data, while the case $\alpha = 3\pi/2$ is analogously called 75\% boundary data. Note that in the 75\% boundary case, the inverse problem shows only mild instability, while in the 100\% case the problem behaves essentially like a well-posed problem. This can be quantified via the condition number of the discretized Fréchet derivative of the underlying nonlinear operator, which is equal to 385 and 12 in the 75\% and 100\% boundary data cases, respectively. The power density data $E_j(\sigma^\dagger)$ is created via solving the PDE (3.4) for $\sigma = \sigma^\dagger$, and for the initial
guess we use $\sigma_0(r, \theta) = 1.5$. For completeness, note that, on the definition space of the underlying nonlinear operator, we use the same weighted inner product as described in [39].

**Fig. 4.7.** Numerical results for the AET problem introduced in Section 3.4 with 75% boundary data. Absolute error (4.1) at the stopping indices $k_\ast$ determined by the discrepancy principle (1.6), the different heuristic parameter choice rules (1.8), and the optimal stopping index $k_{\rm opt}$ defined in (4.2), each for different relative noise levels $\delta$.

**Fig. 4.8.** Numerical results for the AET problem introduced in Section 3.4 with 75% boundary data. Heuristic functionals (1.8) and discrepancy functional (4.3) for $\delta = 2\%$ relative noise (left). Corresponding evolution of the absolute error $\|x_k - x^\dagger\|$ with marked points indicating the stopping indices chosen by the different rules (right).

First, we present results for the 75% boundary data case, for which the resulting absolute errors (4.1) corresponding to the different parameter choice rules can be found in Figure 4.7. As before, characteristic plots of the heuristic functionals $\psi$ and the evolution of the absolute
errors can be found in Figure 4.8. First of all, consider the results of discrepancy principle (1.6), here used with $\tau = 2$. While its corresponding stopping index is not too far off the optimal value, the steep shape of the absolute error curve nevertheless results in an overall large error. While this suggests the use of a smaller $\tau$, note that already in this case for $\delta = 1\%$ the discrepancy principle is not attainable within a reasonable number of iterations. Next, note that the HR rule fails, since its corresponding functional $\psi_{HR}$ is monotonically increasing. From the remaining heuristic parameter choice rules, the LS rule gives the best results overall, determining a stopping index close to the optimal one. In contrast, the HD and QO rules stop the iteration relatively early and late, respectively, and thus lead to suboptimal absolute errors.

Next, we consider the results for the 100\% boundary data case, for which the resulting absolute errors (4.1) corresponding to the different parameter choice rules can be found in Figure 4.9. Characteristic plots of the heuristic functionals $\psi$ and the evolution of the absolute errors can be found in Figure 4.10. Following our previous findings, we now consider the discrepancy principle (1.6) with the choice $\tau = 1.1$. However, apart from the case of $\delta = 2\%$, the resulting stopping index is still far away from the optimal one, with the case of $\delta = 1\%$ being non-attainable, as before. Next, note that the HR, QO, and LS rules all fail, with their corresponding functionals being either monotonically increasing or decreasing. We conjecture that this failure might be related to the fact that the Muckenhoupt condition is not satisfied due to a nearly well-posed situation; cf. Remark 2.1. As noted above, the 100\% boundary case behaves essentially like a well-posed problem, which is reflected in the evolution of the absolute error depicted in Figure 4.10 (right). Consequently a suboptimal stopping index has little effect, and the resulting errors when using the QO and the LS rules are comparable to those of the HD rule, which in this setting is the only heuristic rule producing a well-defined stopping index.

As a further illustration of the results, in Figure 4.11 we provide the calculated approximate solution $x_{k*}$ for this problem with noise level 2\% (corresponding to the second column in Figure 4.7). The exact solution is provided on the left-hand side and is given by (nearly)
Heuristic functionals \((1.8)\) and discrepancy functional \((4.3)\) for \(\delta = 2\%\) relative noise (left). Corresponding evolution of the absolute error \(\|x_\delta^k - x^\dagger\|\), with marked points indicating the stopping indices chosen by the different rules (right).

piecewise constant inclusions of various shapes. Note that this article’s scope is within the realm of Hilbert spaces, and for this problem a standard Sobolev space regularization is used, even though in such situations total variation regularization or iterative regularization in Banach spaces (see, e.g., [1]) might be a preferable choice. However, we would like to point out that the theory of heuristics in such non-Hilbertian cases is still in its infancy and that the behaviour of the heuristic functionals is not well understood. Some investigations in those directions can be found in [46, 47, 55, 58, 66, 89].

Still, despite using only Sobolev regularization, the results for some parameter choice rules are quite satisfactory, an indication that the problem is rather mildly ill-posed. The best (in fact, excellent) results are achieved by the LS rule and the QO rule, which match the result in Figure 4.7. The classical discrepancy rule yields comparably weaker results; again, this might be rooted in a rather conservative choice of the \(\tau\) parameter. The HD and HR rules are disappointing. We note that the artefacts on the right lower part correspond to the lack of data, i.e., the zero-flux condition, there.

4.5. Single photon emission computed tomography. For the fifth test, we consider the nonlinear SPECT problem introduced in Section 3.5. For discretizing the problem, we utilize the same approach as used, for example, in [40, 77, 78], using 79 uniformly spaced angles \(\omega\) in the attenuated Radon transform \((3.5)\). For the exact solution \((f^\dagger, \mu^\dagger)\), we choose the MCAT (mathematical cardiac–torso) phantom [82] (see also [40, 77, 78]), and for the initial guess we use \((f_0, \mu_0) = (0, 0)\). For the discrepancy principle \((1.6)\) we selected the rather large value \(\tau = 10\), which was, however, found to lead to better results than the standard choices \(\tau = 1.1\) or \(\tau = 2\). The absolute error \((4.1)\) corresponding to the different parameter choice rules are depicted in Figure 4.12, in this case for the practically realistic case of noise levels \(\delta\) between 1\% and 10\%. Again, typical plots of the heuristic functionals \(\psi\) as well as the evolution of the absolute error over the iteration are depicted in Figure 4.13.

The plots in these figures show that all parameter choice rules work well for this problem, that the error rates follow the optimal one, and that the heuristic plots in Figure 4.13 have a clear minimum. Furthermore, the HD and HR rules are slightly superior to the other rules in
Reconstructions of electrical conductivity $\sigma$ with the heuristic parameter choice rules. The stopping indices $k_*$ indicate the iteration number.

(a) Exact electrical conductivity $\sigma^\dagger$

(b) DP rule, $k_{DP} = 50$

(c) HD rule, $k_{HD} = 21$

(d) HR rule, $k_{HR} = 1$

(e) QO rule, $k_{QO} = 803$

(f) LS rule, $k_{LS} = 162$

(g) Optimal choice, $k_{opt} = 222$

FIG. 4.11. Numerical results for the AET problem introduced in Section 3.4 with 75% boundary data – the inaccessible boundary part is $[3\pi/2, 2\pi)$: (a) the exact electrical conductivity $\sigma^\dagger$; (b)–(f) its reconstructions with the different heuristic parameter choice rules defined in (1.8); and (g) the optimal stopping index $k_{opt}$ defined in (4.2), for a relative noise level of $\delta = 2\%$.

FIG. 4.12. Numerical results for the nonlinear SPECT problem introduced in Section 3.5. Absolute error (4.1) at the stopping indices $k_*$ determined by the discrepancy principle (1.6), the different heuristic parameter choice rules (1.8), and the optimal stopping index $k_{opt}$ defined in (4.2), each for different relative noise levels $\delta$.

terms of the resulting absolute error. Note that, as for the nonlinear Hammerstein problem, the error graph for the discrepancy principle again exhibits a slight parabolic shape, which could again be related either to a too small choice of $\tau$ or to a coarse discretization of the problem.

4.6. Auto-convolution. For the final test, we consider the auto-convolution problem introduced in Section 3.6. For the discretization of this problem, which is based on standard FEM hat functions on a uniform subdivision of the interval $[0, 1]$ into 60 subintervals, we refer to [41]. For the exact solution, we choose $x^\dagger(s) = 10 + \sqrt{2}\sin(2\pi s)$, from which we compute
FIG. 4.13. Numerical results for the nonlinear SPECT problem introduced in Section 3.5. Heuristic functionals (1.8) and discrepancy functional (4.3) for $\delta = 5\%$ relative noise (left). Corresponding evolution of the absolute error $\|x^\delta_k - x^\dagger\|$, with marked points indicating the stopping indices chosen by the different rules (right).

Since the initial guess is rather close to the exact solution, we now consider noise levels $\delta$ between 0.01% and 0.1%. The corresponding absolute errors (4.1) for the different parameter choice rules are depicted in Figure 4.14, while a typical plot of the heuristic functionals $\psi$ and the evolution of the absolute error can be seen in Figure 4.15.

FIG. 4.14. Numerical results for the auto-convolution problem introduced in Section 3.6. Absolute error (4.1) at the stopping indices $k^*_\delta$ determined by the discrepancy principle (1.6), the different heuristic parameter choice rules (1.8), and the optimal stopping index $k_{\text{opt}}$ defined in (4.2), each for different relative noise levels $\delta$. 

The corresponding data $y$ by applying the operator $F$ as defined in (3.6). For the initial guess we use $x_0(s) = 10 + \frac{1}{2}\sqrt{2}\sin(2\pi s)$, and in the discrepancy principle we choose $\tau = 1.1$.
Fig. 4.15. Numerical results for the auto-convolution problem introduced in Section 3.6. Heuristic functionals (1.8) and discrepancy functional (4.3) for $\delta = 0.05\%$ relative noise (left). Corresponding evolution of the absolute error $\|x^*_k - x^\dagger\|$, with marked points indicating the stopping indices chosen by the different rules (right).

We observe that for this problem the HD rule gives the best results overall, yielding stopping indices close to the optimal $k_{opt}$ for all considered noise levels. Furthermore, from Figure 4.15 we can see that also the functional $\psi_{HR}$ exhibits a clearly distinguishable minimum. However, since the absolute error drops steeply between the corresponding stopping index and the optimal one, the resulting absolute error when using the HR rule is significantly higher. In contrast, the functional values $\psi(k, y^\delta)$ for both the QO rule and the LS rule tend towards $-\infty$ as the iteration number increases, leading to no interior minimum. (Note that, in the nonlinear case, non-negativity of $\psi(k, y^\delta)$ is not guaranteed [58].) This coincides with the fact that the absolute error in the iteration stays more or less constant after having reached the minimum value at $k_{opt}$. Consequently, the LS and the QO rule both stop with $k_s = k_{max}$, which in this case by chance leads to generally very good absolute errors. However, the shape of the QO and LS functionals is in contrast to theory, as the graph is expected to diverge for $k \to \infty$.

This observation is a hint that the Muckenhoupt condition is not satisfied here for QO and LS, while it might be for HD and HR; cf. Remark 2.1. This may also be connected to the fact that the residual functional $x \mapsto \|F(x) - y^\delta\|^2$ is locally convex around our chosen solution $x^\dagger$ (see, e.g., [41, Proposition 5.2]), and thus the auto-convolution problem may behave nearly like a well-posed problem; cf. Section 2.3.

5. Summary and conclusion. In the previous sections, we have presented and discussed the results of the different parameter choice rules applied to the test problems introduced in Section 3. As a summary, we now present our findings in Table 5.1, where we classify the results, in a rather informal style, together with comments and suspected background.

In the first column of Table 5.1, we indicate the (suspected) type of ill-posedness of the various test problems and whether the tangential cone condition (TCC) is known to hold. Furthermore, in the second column the performance of the parameter choice rules is classified according to two indicators: The row with “$k_s \sim k_{opt}$” indicates whether the stopping index selected by the respective parameter choice rule is close to the optimal index, and whether the heuristic functionals behave as desired, i.e., have a clear minimum. The row with “Error” indicates how close the resulting absolute error is to the optimal error. These performance
### Table 5.1

Summary of performance for various examples and stopping rules.

| Example                               | Indicator | $\delta$ rule | Heuristic rules |
|---------------------------------------|-----------|----------------|-----------------|
|                                        |           | $\delta$ small: Good | DP  | HD | HR | QO | LS |
| **Hammerstein equation**              | $k_\star \sim k_{opt}$ | Good          | $\delta$ large: Bad | Average |
| Ill-posedness: mild                   |           | $\delta$ small: Good | Good |
| TCC: yes                              |           | $\delta$ large: Bad | Good |
| **Diffusion estimation**              | $k_\star \sim k_{opt}$ | Good          | $\delta$ small: Good | Good |
| Ill-posedness: mild                   |           | $\delta$ large: Bad | Good |
| TCC: yes                              |           | $\delta$ small: Good | Good |
| **Elastography**                     | $k_\star \sim k_{opt}$ | Good          | $\delta$ large: Bad | Good |
| Ill-posedness: mild                   |           | $\delta$ small: Good | Excellent |
| TCC: yes                              |           | $\delta$ large: Bad | Good |
| **Acousto-electrical tomography 75%** | $k_\star \sim k_{opt}$ | Good          | $\delta$ large: Bad | Excellent |
| Ill-posedness: mild                   |           | $\delta$ small: Good | Good |
| TCC: unknown                          |           | $\delta$ large: Bad | Good |
| **Acousto-electrical tomography 100%**| $k_\star \sim k_{opt}$ | Good          | $\delta$ large: Bad | Excellent |
| Near-well-posedness                   |           | $\delta$ small: Good | Good |
| TCC: unknown                          |           | $\delta$ large: Bad | Good |
| **SPECT**                             | $k_\star \sim k_{opt}$ | Good          | $\delta$ large: Bad | Good |
| Ill-posedness: mild                   |           | $\delta$ small: Good | Good |
| TCC: unknown                          |           | $\delta$ large: Bad | Good |
| **Auto-convolution**                 | $k_\star \sim k_{opt}$ | Good          | $\delta$ large: Bad | Good |
| Ill-posedness: mild                   |           | $\delta$ small: Good | Good |
| TCC: unknown                          |           | $\delta$ large: Bad | Excellent |
| **Acousto-electrical tomography 75%** | $k_\star \sim k_{opt}$ | Good          | $\delta$ large: Bad | Excellent |
| Ill-posedness: mild                   |           | $\delta$ small: Good | Good |
| TCC: unknown                          |           | $\delta$ large: Bad | Good |
indicators are stated in a colloquial manner, classified as “Excellent”, “Good”, “Average”, or “Bad”.

Some conclusions can be drawn from these results. First of all, the discrepancy principle works well in all cases, even when the tangential cone condition is not known to hold, as discussed above. However, its reliability strongly correlates with being able to find a proper choice of the parameter $\tau$, which is not at all an easy task in practice. The strange parabolic shape in Figure 4.1 is, for example, attributed to its having been selected too small. Next, we observe that the heuristic rules work well for many cases, but not always. Furthermore, our numerical results show that it is not possible to determine an overall “best” heuristic parameter choice rule for all test problems. This suggests that, in practice, one should always first conduct a series of simulations using multiple different parameter choice rules for any given problem, instead of blindly selecting any single rule among them. From the computational point of view, the authors personally prefer the HD rule, which does not require one to compute and store both iterates $x_k$ and $x_{2k}$ during the iteration. However, also the HD rule fails at times, and the LS rule in particular has been found to be a useful alternative which often showed an excellent performance with respect to the resulting absolute error. One issue with the QO and LS rules is that their corresponding functionals $\psi$ do not exhibit a clear minimum in the case of well-posed or nearly well-posed problems (e.g., AET with 100% boundary data and auto-convolution). However, this can be attributed to the fact that the Muckenhoupt condition is stronger for these rules and hence might not be satisfied; cf. Remark 2.1. On the other hand, even if the QO and LS rules fail in finding a good approximation of $k_{opt}$, the resulting error is then not too dramatic, since the almost well-posedness then only leads to moderate errors.

Finally, our numerical studies also have implications for the further analytic study and development of (novel) heuristic stopping rules for nonlinear Landweber iteration. In particular, our findings indicate that, apart from the relation between the smoothness of the solution and the noise (i.e., noise conditions), also the type of ill-posedness and especially the type of nonlinearity of the problem have to be taken into account. The authors believe that it will in general be impossible to design and analyse a single heuristic parameter choice rule which performs well for all different nonlinear inverse problems. Rather, one may have to consider different heuristic rules depending on the type of nonlinearity and ill-posedness of each specific problem. This then poses an interesting challenge for future research.

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