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BACKWARD ERROR ANALYSIS OF THE SHIFT-AND-INVERT ARNOLDI ALGORITHM

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ABSTRACT. We perform a backward error analysis of the inexact shift-andinvert Arnoldi algorithm. We consider inexactness in the solution of the arising linear systems, as well as in the orthonormalization steps, and take the nonorthonormality of the computed Krylov basis into account. We show that the computed basis and Hessenberg matrix satisfy an exact shift-and-invert Krylov relation for a perturbed matrix, and we give bounds for the perturbation. We show that the shift-and-invert Arnoldi algorithm is backward stable if the condition number of the small Hessenberg matrix is not too large. This condition is then relaxed using implicit restarts. Moreover we give notes on the Hermitian case, considering Hermitian backward errors, and finally, we use our analysis to derive a sensible breakdown condition.

1. INTRODUCTION

Consider an implementation of the Arnoldi algorithm [4, 25]. Not much meaning can be given to the computed quantities if they deviate too much from the recurrence that underpins the algorithm in exact arithmetic:

$$AV_k = V_{k+1}\underline{H}_k, \quad \underline{H}_k = H(1:k+1,1:k).$$

Luckily, good implementations, where in particular the orthogonalization is done with care, can be shown to be backward stable [3, 8, 10, 20] in the sense that the computed quantities V_{k+1} and \underline{H}_k satisfy an exact recurrence with a slightly perturbed matrix:

(1)
$$(A + \Delta A)V_k = V_{k+1}\underline{H}_k$$

This means that we can compute a basis of an exact Krylov subspace corresponding to a nearby matrix. Since the basis will in general not be perfectly orthonormal, so $V_{k+1}^H V_{k+1} \neq I$, we use the term "Krylov recurrence" instead of "Arnoldi recurrence" when referring to recurrences like (1). If A is Hermitian, then it can be shown that the computed basis spans a Krylov subspace associated with a perturbed *Hermitian* matrix $A + \Delta A$ [14]. There is a catch in this case, though: the small $(k+1) \times k$ matrix

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Algorithm 1 The Shift-and-invert Arnoldi algorithm

Input: A, σ, b, k **Output:** $V_{k+1} := [v_1, \dots, v_{k+1}], \underline{H}_k = [h_{ij}]_{i=1:k+1,j=1:k}$ $v_1 = b/||b||$ **for** $j = 1, 2, \dots, k$ $w_j = (A - \sigma I)^{-1}v_j$ $[w'_j, h_{1:j}] = \text{orthogonalization}(w_j, V_j)$ $h_{j+1 j} = ||w'_j||$ **if** $h_{j+1 j} = 0$ **break** $v_{j+1} = w'_j/h_{j+1 j}$ **end for**

associated with this Krylov subspace is in general not the computed Hessenberg matrix.

In this paper we perform a similar backward error analysis of the shift-and-invert Arnoldi algorithm. For example, we show that an implementation of the Arnoldi algorithm applied to A^{-1} , yields computed matrices V_{k+1} and \underline{H}_k such that

$$(A + \Delta A)^{-1}V_k = V_{k+1}\underline{H}_k,$$

and we give an upper bound for $\|\Delta A\|_2$. Perturbed versions of the shift-and-invert Arnoldi algorithm have been considered in the literature as a part of the theory of *inexact methods*, see [15, 18]. However, these results neglect that the orthonormalization is not performed exactly, and furthermore, assume bounds on linear system residuals that may be unattainable (more on this in Section 2). We consider more general linear system residuals and take the error from the orthonormalization into account. Our analysis of how the orthonormalization errors propagates into the shift-and-invert Krylov recurrence highlight the importance of *columnwise* backward error bounds for QR factorizations, and is thus of a different flavor than the corresponding analysis for standard Arnoldi, done in, for example [8].

We also use our error analysis to motivate when "breakdown" should be declared, that is, when $h_{j+1,j}$ may be considered to be "numerically zero."

The algorithm we study can be divided into two main subproblems: solving linear systems and orthonormalizing vectors. We state our backward error results in such a way that they are independent of how these subproblems are being solved, but we also discuss relevant and commonly used approaches for solving these two tasks.

1.1. Technical outline. We study floating point implementations of Algorithm 1, where A is assumed to be of size $n \times n$, σ is the shift, b the starting vector, and k is the maximum number of steps we perform. Throughout the paper $\|\cdot\|$ refers to the 2-norm. In exact arithmetic, we have

orthogonalization $(w_j, V_j) := [w_j - V_j(V_j^H w_j), V_j^H w_j],$

which corresponds to classical Gram-Schmidt if implemented as it stands. In floating point arithmetic, orthogonalization routines with better numerical properties, such as modified Gram-Schmidt, are usually employed.

In the *j*th iteration in Algorithm 1, a new vector w_j is computed and decomposed into a linear combination of v_1, \ldots, v_j and a new component that will be

the definition of v_{j+1} . In exact arithmetic, this can be described by the Arnoldi recurrence

$$(A - \sigma I)^{-1}v_j = V_k h_{1:j,j} + h_{j+1,j}v_{j+1}.$$

When the corresponding thing is done in practice, however, errors are present in all steps of the computation. First, we need to solve a linear system. If we use a direct solver the matrix $A - \sigma I$ needs to be formed. We consider the rounding error in this step as part of the residual from the linear system. This does not affect the norm of the residual significantly, because the rounding error is very small,

$$\|\operatorname{float}(A - \sigma I) - (A - \sigma I)\| < \max_{1 \le i \le n} |a_{ii} - \sigma| u \le u \|A - \sigma I\|.$$

Here float $(A - \sigma I)$ refers to the computed shifted matrix and u is the unit roundoff. Let r_j be the said residual from the linear system, so

(2)
$$(A - \sigma I)w_j = v_j + r_j$$

is the actual linear system that has been solved. Then we have the following equality for the computed quantities:

$$(A - \sigma I)^{-1}(v_j + r_j) = w_j = V_{j+1}h_{1:j+1,j} + g_j,$$

where g_j is an error coming from the orthonormalization process. Defining

$$f_j = r_j - (A - \sigma I)g_j$$

and $F_k = [f_1, f_2, \dots, f_k]$ yields a perturbed recurrence

$$(A - \sigma I)^{-1}(V_k + F_k) = V_{j+1}\underline{H}_k.$$

We discuss the residual r_j and the error g_j in Section 2 and Section 3, respectively, and provide bounds for both quantities. In Section 4, we use these bounds in order to bound F_k , and subsequently the backward error for the shift-and-invert Arnoldi recurrence. In Section 5, we explain how the idea of implicit restarting can be used to gain further insight into the backward error. We also discuss in what sense we have Hermitian backward errors if the method is applied to a Hermitian matrix A. Finally, we talk about breakdown conditions: in floating point arithmetic, the test if $h_{j+1,j} = 0$ in Algorithm 1 is rarely done. Instead one usually checks whether $h_{j+1,j}$ is "small enough." This case is referred to as *breakdown*. A sensible definition of "small enough" is when the quantity is dominated by errors. We discuss this in more detail and derive backward error bounds for this case.

1.2. Notation. The scalar σ refers to a shift while $\sigma_{\min}(X)$ refers to the smallest singular value of X. The dagger notation X^{\dagger} refers to the Moore-Penrose pseudoinverse of X. The lower letter u is reserved to denote the unit roundoff if real arithmetic is used, and $\sqrt{5}$ times the unit roundoff if complex arithmetic is used [7], respectively. When the matrix size is understood from the context, we denote zero matrices and identity matrices as 0 and I, respectively. Similarly, the vector e_i denotes the *i*th column of the identity matrix whose size is understood from the context. For a matrix X, the lower case x_i refers to the *i*th column of X and X_k to $[x_1, x_2, \ldots, x_k]$, that is, the first k columns of X.

2. Errors from linear systems

In this section we discuss bounds on the residual r_i from (2).

2.1. Backward error bounds. The normwise backward error associated with a computed solution y of a linear Ax = b is defined as

$$\eta_{A,b}(y) := \min\{\epsilon : (A + \Delta A)y = b + \Delta b, \|\Delta A\| \le \epsilon \|A\|, \|\Delta b\| \le \epsilon \|b\|\},\$$

and given by the formula

(3)
$$\eta_{A,b}(y) = \|r\|/(\|A\|\|y\| + \|b\|)$$

where r = Ay - b [19]. See also [11, p. 120]. This result is true for any vector norm $\|\cdot\|$ and its subordinate matrix norm. Thus, if we solve the linear systems in Algorithm 1, up to a backward error $\epsilon_{\rm bw}$, then it holds that

(4)
$$||r_j|| \le (||A - \sigma I|| ||w_j|| + ||v_j||)\epsilon_{\text{bw}}$$

where r_j is defined in (2). If the linear systems are solved by a backward stable direct method, we have $\epsilon_{\rm bw} \leq \phi(n)u$, where $\phi(n)$ is an algorithm dependent constant. If we are interested in the smallest possible $\epsilon_{\rm bw}$ such that (4) holds, then we need to compute $||r_j||/(||A - \sigma I||||w_j|| + ||v_j||)$. However, this may not be feasible for the 2-norm, due to the term $||A - \sigma I||$. In these cases we can replace $||A - \sigma I||$ by a lower bound (the tighter the better), and thus obtain an upper bound for $\epsilon_{\rm bw}$. We can for instance do a few iterations of the power method applied to $(A - \sigma I)^H (A - \sigma I)$. MATLAB's **normest** function does exactly this. This would lead to a lower bound of $||A - \sigma I||$, since convergence is always from below. Another possibility is to use the (lower) bound in [12]. We can also bound the matrix 2-norm in terms of the corresponding infinity-norm or 1-norm. The following proposition shows that such bounds can be satisfactory for many sparse matrices, in particular those which can be permuted to banded form.

Proposition 2.1. Let k_{row} and k_{col} denote the maximum number of nonzero entries in a row and column of A, respectively. Then the following two upper and lower bounds hold.

$$\frac{1}{\sqrt{k_{\rm col}}} \|A\|_2 \leq \|A\|_{\infty} \leq \sqrt{k_{\rm row}} \|A\|_2,$$

$$\frac{1}{\sqrt{k_{\rm row}}} \|A\|_2 \leq \|A\|_1 \leq \sqrt{k_{\rm col}} \|A\|_2.$$

Proof. We have $||A||_{\infty} = ||Ax||_{\infty}$ for some x with $||x||_{\infty} = 1$ and at most k_{row} nonzeros. We get

$$||A||_{\infty} \le ||Ax||_{\infty} \le ||Ax||_{2} \le ||A||_{2} ||x||_{2} \le \sqrt{k_{\text{row}}} ||A||_{2},$$

which is the desired upper bound for $||A||_{\infty}$. Further, we have

$$||A||_1 = ||A^T||_{\infty} \le \sqrt{k_{\text{col}}} ||A^T||_2 = \sqrt{k_{\text{col}}} ||A||_2,$$

which is the desired upper bound for $||A||_1$.

The lower bounds follow from [24, Theorem 4.2].

The inequality (4) can also be used as a stopping criterion for iterative linear system solvers [2]. In this case, $\epsilon_{\rm bw}$ denotes the desired backward error, which is given prior to execution. If we replace $||A - \sigma I||$ with a lower bound, then we get a more stringent stopping criterion.

2.2. Residual reduction bounds. An alternative to (4) is to use the bound

(5)
$$||r_j|| \le ||v_j||\epsilon_{\text{tol}}$$

This bound is commonly used as a stopping condition when the linear systems are solved by iterative methods. Unfortunately, as a stopping condition, (5) "may be very stringent, and possibly unsatisfiable" [11, p. 336]. See also [9, pp. 72–73] for a 2×2 example that illustrates the pitfall of comparing the norm of the residual with the norm of the right hand side. However, since (5) is de facto commonly used in computer codes it is still worth to study it under the assumption that the stopping criterion is met.

2.3. Auxiliary residual bounds. In order to treat both (4) and (5) in a unified way, we consider the following auxiliary bound

(6)
$$||r_j|| \le ||v_j||\epsilon_1 + ||A - \sigma I|| ||w_j||\epsilon_2.$$

Clearly, the substitutions $(\epsilon_1, \epsilon_2) \leftarrow (\epsilon_{\text{bw}}, \epsilon_{\text{bw}})$ and $(\epsilon_1, \epsilon_2) \leftarrow (\epsilon_{\text{tol}}, 0)$ give back (4) and (5), respectively. We can simplify the bound in (6) in cases when $A - \sigma I$ is not too ill-conditioned with respect to ϵ_2 . To see this we need the following lemma.

Lemma 2.2. If $\kappa(A - \sigma I)\epsilon_2 < 1$ and (6) hold, then

$$\|r_j\| \le \frac{\epsilon_1 + \kappa (A - \sigma I)\epsilon_2}{1 - \kappa (A - \sigma I)\epsilon_2} \|v_j\|.$$

Proof. We have

$$||r_j|| \le ||A - \sigma I|| ||(A - \sigma I)^{-1} (v_j + r_j)||\epsilon_2 + ||v_j||\epsilon_1$$

$$\le \kappa (A - \sigma I) ||v_j + r_j||\epsilon_2 + ||v_j||\epsilon_1$$

$$\le \kappa (A - \sigma I) (||v_j|| + ||r_j||)\epsilon_2 + ||v_j||\epsilon_1.$$

Reordering gives the result.

The following result yields a family of new residual bounds independent of $||v_i||$.

Proposition 2.3. Let $(A - \sigma I)^{-1}(v_j + r_j) = w_j$ and assume (6) hold. If

(7)
$$0 < \frac{\epsilon_1 + \kappa (A - \sigma I)\epsilon_2}{1 - \kappa (A - \sigma I)\epsilon_2} \le \gamma < 1,$$

then

$$||r_j|| \le \left(\epsilon_2 + \frac{\epsilon_1}{1-\gamma}\right) ||A - \sigma I|| ||w_j||.$$

Proof. From (6) we have

$$||r_j|| \le \left(\epsilon_2 + \epsilon_1 \frac{||v_j||}{||A - \sigma I|| ||w_j||}\right) ||A - \sigma I|| ||w_j||.$$

Thus we need to show $||v_j||/(||A - \sigma I||||w_j||) \le 1/(1 - \gamma)$. We have

$$\frac{\|v_j\|}{\|A - \sigma I\| \|w_j\|} = \frac{\|v_j\|}{\|A - \sigma I\| \|(A - \sigma I)^{-1}(v_j + r_j)\|} \le \frac{\|v_j\|}{\|v_j + r_j\|},$$

and from the reverse triangle inequality,

$$\frac{\|v_j\|}{\|v_j + r_j\|} \le \frac{\|v_j\|}{\||v_j\| - \|r_j\||}.$$

Now, by Lemma 2.2 and assumption (7), we have

$$\|r_j\| \le \frac{\epsilon_1 + \kappa (A - \sigma I)\epsilon_2}{1 - \kappa (A - \sigma I)\epsilon_2} \|v_j\| \le \gamma \|v_j\|.$$

Putting everything together yields

$$\frac{\|v_j\|}{\|A - \sigma I\| \|w_j\|} \le \frac{\|v_j\|}{\|\|v_j\| - \|r_j\|\|} \le \frac{1}{1 - \gamma}.$$

In particular if $\kappa(A - \sigma I) \leq (1 - 2\epsilon_1)/(3\epsilon_2)$, then we have $\kappa(A - \sigma I)\epsilon_2 < 1$ and can take $\gamma = 1/2$ in Proposition 2.3, to obtain

(8)
$$||r_j|| \le (2\epsilon_1 + \epsilon_2)||A - \sigma I||||w_j||.$$

This is the same bound as we get from (6) if we replace (ϵ_1, ϵ_2) with $(0, 2\epsilon_1 + \epsilon_2)$. In particular, if the linear systems are solved in a backward stable manner so that (4) holds, and $\kappa(A - \sigma I) \leq (1 - 2\epsilon_{\rm bw})/(3\epsilon_{\rm bw})$, then (8) holds with $2\epsilon_1 + \epsilon_2 = 3\epsilon_{\rm bw}$.

3. Errors from orthonormalization

In this section we are concerned with the orthonormalization error

$$g_j = w_j - V_{j+1}h_{1:j+1,j}$$

Up to signs, this error can be viewed as the backward error in the (i + 1)st column of a perturbed QR factorization

(9)
$$[v_1 \ w_1 \ w_2 \ \cdots \ w_k] = V_{k+1}[e_1 \ \underline{H}_k] + [0 \ g_1 \ g_2 \ \cdots \ g_k].$$

Thus, we are interested in *columnwise* backward error bounds for QR factorizations. The next theorem shows how such bounds can be obtained from normwise backward error bounds given in the 2-norm or the Frobenius norm. It applies to floating point algorithms $qr(\cdot)$ that are unaffected by power-of-two column scalings, in the sense that if [Q, R] = qr(A), then [Q, RD] = qr(AD) for any $D = diag(d_1, d_2, \ldots, d_k)$ where the d_i are powers of 2. Barring underflow and overflow, this covers commonly used QR algorithms such as classical and modified Gram-Schmidt with and without (possibly partial) reorthogonalization, Householder QR and Givens QR.

Theorem 3.1. Let qr(A) denote an algorithm that computes an approximate QR factorization of an $n \times k$ matrix A in floating point arithmetic. Suppose further that [Q, RD] = qr(AD) for any $D = diag(d_1, d_2, \dots, d_k)$ where the d_i are powers of 2. If Q and R denotes the computed factors, $\Delta A = A - QR$ and $\|\Delta A\|_* \leq \gamma \|A\|_* u$, where $\|\cdot\|_*$ denotes the 2-norm or the Frobenius norm, then $\|\Delta a_i\| \leq 2\gamma \sqrt{k} \|a_i\| u$ for i = 1:k.

Proof. For i = 1:k, we define

$$d_i = 2^{-\lfloor \log_2 \|a_i\| \rfloor}.$$

so $1 \leq ||a_i|| d_i < 2$. Since ΔAD is the backward error from qr(AD) we have

 $d_i \|\Delta a_i\| = \|\Delta ADe_i\| < \|\Delta AD\|_* < \gamma \|AD\|_* u < 2\gamma \sqrt{k} \|ADe_i\| u = d_i 2\gamma \sqrt{k} \|a_i\| u,$ for i = 1: k, from which the theorem follows.

The constant γ in Theorem 3.1 is obviously algorithm dependent and many bounds exist in the literature. Some of them contain both n and k [21], and others only k [5, 1], [11, Theorem 19.13]. In [11, p. 361] a columnwise bound depending on n and k is given. For Krylov methods we usually have $n \gg k$, so bounds independent from n should certainly be favored. We shall assume that

(10)
$$\|g_j\| \le \eta(n,k) \|w_j\| u$$

holds for some function $\eta(n, k)$.

3.1. Columnwise backward errors for modified Gram-Schmidt. Our next theorem shows that for modified Gram-Schmidt (MGS), with and without one round of reorthogonalization, η in (10) does not depend on n and is given by

$$\eta(n,k) = \zeta k$$

where ζ is a modest constant. We need the following forward error result for _axpy operations.

Lemma 3.2. Let α be a scalar and x and y vectors. If

$$s = \text{float}(\alpha x + y) - (\alpha x + y)$$
 then $||s|| \le 2(||\alpha x|| + ||y||)u$.

Proof. The *i*th component of $\alpha x + y$ can be viewed as the inner product $[x_i \ y_i] [\alpha \ 1]^T$. Thus the componentwise forward error is bounded by $|s| \leq 2u(|\alpha x| + |y|)$ [13]. We get

$$||s|| \le ||2u(|\alpha x| + |y|)|| \le 2(||\alpha x|| + ||y||)u.$$

The next theorem gives columnwise backward error bounds for MGS with and without one round of reorthogonalization.

Theorem 3.3. Let Q and R denote the computed factors in the QR decomposition of an $n \times k$ matrix A, which was obtained by a floating point implementation of modified Gram-Schmidt with or without one round of reorthogonalization. Assume

- (i) $||q_j|| = 1$ for j = 1:k, and (ii) $(1 + (n+3)u)^k < 1 + \delta$ for some $\delta > 0$.

Then there exists a ΔA such that $A + \Delta A = QR$ with $\|\Delta a_j\| \leq cj \|a_j\| u$, where $c = 4(1 + \delta)$ if no reorthogonalization was done and $c = 10(1 + \delta)^2$ if one round of reorthogonalization was done.

Let us pause for a while and discuss the assumptions before proceed with the proof. Assumption (i) is imposed to keep our analysis cleaner; it does not affect our final bounds in any significant way. Assumption (ii) is needed for the following reason: if we compute $y = \text{float}(x - q_j(q_j^H x))$ for some $1 \le j \le k$, then, assuming (i), the quantity $1 + (n+3)u = 1 + ||q_j||^2 (n+3)u$ is an upper bound for ||y|| / ||x|| [11, Lemma 3.9]. Thus, (ii) guarantees that we can apply a sequence of k elementary "floating point" projections of the form $I - q_i q_i^H$ to any vector x, and the resulting vector will be bounded in norm by $(1 + \delta) ||x||$.

Proof of Theorem 3.3. Let $R^{(1)}$ and $R^{(2)}$ denote the strictly upper triangular matrices containing the orthogonalization coefficients corresponding to the first and

second round of orthogonalization, respectively. We define $R^{(2)} \equiv 0$, if no reorthogonalization is done. Assume for a while that $R^{(1)}$ and $R^{(2)}$ are given, and suppose we want to compute

$$a_j - \sum_{i=1}^{j-1} r_{ij}^{(1)} q_i - \sum_{i=1}^{j-1} r_{ij}^{(2)} q_i$$

This can be viewed as 2(j-1) saypy operations. We define $a_j^{(0)} = a_j$ and

$$a_{j}^{(i)} = \begin{cases} \text{float}(a_{j}^{(i-1)} - r_{ij}^{(1)}q_{i}) & \text{for } i = 1:j-1, \\ \text{float}(a_{j}^{(i-1)} - r_{(i-j+1)j}^{(2)}q_{i-j+1}) & \text{for } i = j:2(j-1). \end{cases}$$

Using Lemma 3.2 yields

$$a_{j}^{(i)} = \begin{cases} a_{j}^{(i-1)} - r_{ij}^{(1)}q_{i} + s_{i} & \text{for } i = 1:j-1, \\ a_{j}^{(i-1)} - r_{(i-j+1)j}^{(2)}q_{i-j+1} + s_{i} & \text{for } i = j:2(j-1), \end{cases}$$

where

$$\|s_i\| \le \begin{cases} 2(\|r_{ij}^{(1)}q_i\| + \|a_j^{(i-1)}\|)u & \text{for } i = 1:j-1, \\ 2(\|r_{(i-j+1)j}^{(2)}q_{i-j+1}\| + \|a_j^{(i-1)}\|)u & \text{for } i = j:2(j-1) \end{cases}$$

Now, $a_j^{(i-1)}$ is also the result of applying i-1 elementary floating point projections to a_j , so the discussion prior to the proof gives $||a_j^{(i-1)}|| < (1+(n+3)u)^{i-1}||a_j||$. Further, from (ii) we have $(1+nu)||a_j^{(i-1)}|| < (1+\delta)||a_j||$ for i=1:j-1 and $(1+nu)||a_j^{(i-1)}|| < (1+\delta)^2||a_j||$ for i=j:2(j-1). The forward error of a computed inner product float $(x^H y)$, where x and y are of length n, is bounded by nu||x|||y|| [13]. Thus

$$|r_{ij}^{(1)}| \le |\text{float}(q_i^H a_j^{(i-1)})| \le |q_i^H a_j^{(i-1)}| + nu ||a_j^{(i-1)}|| < (1+\delta) ||a_j||$$

and, similarly, that $|r_{ij}^{(2)}| < (1+\delta)^2 ||a_j||$. Thus s_i is bounded by

$$\|s_i\| \le \begin{cases} 4(1+\delta)\|a_j\|u & \text{ for } i=1:j-1, \\ 4(1+\delta)^2\|a_j\|u & \text{ for } i=j:2(j-1). \end{cases}$$

We have

$$a_j - \sum_{i=1}^{j-1} r_{ij}^{(1)} q_i - \sum_{i=1}^{j-1} r_{ij}^{(2)} q_i = a_j^{(2(j-1))} - \sum_{i=1}^{2(j-1)} s_i$$

If we define $d_i = \text{float}(||a_j^{(2(j-1))}||)$ and $q_j = \text{float}(a_j^{(2(j-1))}/d_j)$ and note that

$$a_j^{(2(j-1))} = q_j d_j + f_j \quad \text{with} \quad \|f_j\| \le \|a_j^{(2(j-1))}\| u < (1+\delta)^2 \|a_j\| u,$$

then we get

$$a_j - \sum_{i=1}^{j-1} (r_{ij}^{(1)} + r_{ij}^{(2)})q_i - d_j q_j = f_j - \sum_{i=1}^{2(j-1)} s_i.$$

Finally, defining $R = \text{float}(R^{(1)} + R^{(2)}) + \text{diag}(d_1, d_2, \dots, d_k)$ yields

$$\Delta a_j := a_j - \sum_{i=1}^j r_{ij} q_i = f_j - \sum_{i=1}^{2(j-1)} s_i - \sum_{i=1}^{j-1} \Delta r_{ij} q_i,$$

where

$$\Delta r_{ij} = r_{ij}^{(1)} + r_{ij}^{(2)} - r_{ij}, \quad \text{so} \quad |\Delta r_{ij}| \le |r_{ij}^{(1)} + r_{ij}^{(2)}|u < 2(1+\delta)^2 ||a_j||u.$$

Using the above bounds for f_j , the s_i and Δr_{ij} gives $\|\Delta a_j\| < 10(1+\delta)^2 j \|a_j\| u$. If no reorthogonalization was done, then we have $s_i = 0$ for i = j : 2(j-1), and $\Delta r_{ij} = 0$, $\|f_j\| \le (1+\delta) \|a_j\| u$ for all j. Taking this into account yields $\|\Delta a_j\| < 4(1+\delta)j\|a_j\|u$.

Remark 3.4. Suppose the perturbed QR factorization (9) was computed using MGS. Then, taking $\delta = 1/10$ and assuming the conditions of Theorem 3.3 hold, imply that $\eta(n,k)$ in (10) is bounded by $\eta(n,k) \leq 5k$ if standard MGS is used, and $\eta(n,k) \leq 13k$ if MGS with one round of reorthogonalization is used.

4. Backward error bounds for the shift-and-invert Arnoldi recurrence

Recall the perturbed Krylov recurrence

(11)
$$(A - \sigma I)^{-1}(V_k + F_k) = V_{j+1}\underline{H}_k$$

where $F_k = [f_1, f_2, \ldots, f_k]$ and f_j , for j = 1:k, is defined by $f_j = r_j - (A - \sigma I)g_j$. We discussed in sections 2 and 3 how to bound r_j and g_j , respectively. By using these bounds, we can now easily bound F_k . Assuming (6) and (10) yields

(12)
$$||f_j|| \le ||v_j||\epsilon_1 + ||A - \sigma I|| ||w_j||(\epsilon_2 + \eta(n, j)u)$$

Further, from (9) we see that

$$|w_j|| = ||V_{j+1}h_{1:j+1,j} + g_j|| \le ||V_{j+1}|| ||h_{1:j+1,j}|| + \eta(n,j)||w_j||u,$$

which in turn implies

$$||w_j|| \le \frac{||V_{j+1}|| ||h_{1:j+1,j}||}{1 - \eta(n,j)u},$$

assuming that $\eta(n, j)u < 1$. We get

$$||f_j|| \le ||v_j||\epsilon_1 + ||A - \sigma I|| ||V_{j+1}|| ||h_{1:j+1,j}||c_{jn}(\epsilon_2)|$$

and further (assuming that $\eta(n, k)$ is monotonically increasing in k)

(13)
$$||F_k|| \le \sqrt{k} ||V_k|| \epsilon_1 + \sqrt{k} ||A - \sigma I|| ||V_{k+1}|| ||\underline{H}_k|| c_{kn}(\epsilon_2),$$

where

(14)
$$c_{kn}(\epsilon_2) := \frac{\epsilon_2 + \eta(n,k)u}{1 - \eta(n,k)u}$$

should be thought of as a tiny factor.

Similarly, if we assume the bound (8) instead of (6), we get

(15)
$$||F_k|| \le \sqrt{k} ||A - \sigma I|| ||V_{k+1}|| ||\underline{H}_k|| c_{kn} (2\epsilon_1 + \epsilon_2).$$

This is the same bound we get from (13) if we replace (ϵ_1, ϵ_2) by $(0, 2\epsilon_1 + \epsilon_2)$.

Having established (13) and (15), we are now ready to reshuffle equation (11) in order to derive backward error bounds for the shift-and-invert Krylov recurrence. We will derive perturbed recurrences of the form

(16)
$$V_k = (A + \Delta A - \sigma I)V_{k+1}\underline{H}_k.$$

If we look at this from a backward error perspective, (16) means that we have taken k steps, without errors, of a shift-and-invert Krylov algorithm applied to a

perturbed pencil, and all linear systems that occurred in the process must have been consistent. However, in order to rewrite (16) as

$$(A + \Delta A - \sigma I)^{-1}V_k = V_{k+1}\underline{H}_k,$$

we need to ensure that $A + \Delta A - \sigma I$ is invertible. We need the following lemma to solve this technicality.

Lemma 4.1. Let A and V be matrices of size $n \times n$ and $n \times k$ respectively, such that rank AV = k. Then for any $\epsilon > 0$, there exists a matrix X with $||X|| < \epsilon$ such that A + X is nonsingular and XV = 0. Furthermore, if A is Hermitian, then we may take X to be Hermitian.

Proof. Find a unitary matrix Q such that

(17)
$$Q^H V = \begin{bmatrix} 0\\ V_2 \end{bmatrix}$$

for some $k \times k$ matrix V_2 , and define $AQ = [A_1 \ A_2]$ where A_2 is of size $n \times k$. From rank AV = k, it follows A_2 has rank k. Define Y so its columns span the orthogonal complement to range of A_2 , and set $Z = [Y - A_1 \ 0]$. We have that $A + ZQ^H = [Y \ A_2]Q^H$ is nonsingular and $ZQ^HV = 0$. In particular, this means that the pencil $A + \lambda ZQ^H$ is regular. If λ is any value outside the spectrum of the pencil such that $|\lambda| < \epsilon/||Z||$, then $X = \lambda ZQ^H$ satisfies the conditions of the theorem.

For the second part, suppose A is Hermitian and Q is such that (17) holds. Write

$$Q^{H}AQ = \begin{bmatrix} A_{11} & A_{12} \\ A_{12}^{H} & A_{22} \end{bmatrix} \text{ and } W = \begin{bmatrix} \omega I - A_{11} & 0 \\ 0 & 0 \end{bmatrix}, \quad \omega > 0,$$

where A_{11} is of size $(n-k) \times (n-k)$. We have that QWQ^H is Hermitian, $QWQ^HV = 0$, and

$$Q(Q^H A Q + W)Q^H = A + QWQ^H$$

Thus, for the same reason as above, it is enough to find one $\omega > 0$ such that $Q^H A Q + W$ is nonsingular. Let

$$A_{22} = U \begin{bmatrix} D & 0 \\ 0 & 0 \end{bmatrix} U^H$$

be a spectral decomposition so D is of full rank and define $[B_1 B_2] = A_{12}U$ such that B_1 has as many columns as D. We have that $Q^H A Q + W$ is nonsingular if and only if

$$\begin{bmatrix} \omega I & B_1 & \omega B_2 \\ B_1^H & D & 0 \\ \omega B_2^H & 0 & 0 \end{bmatrix}$$

is nonsingular. Further, since $[A_{12}^T \; A_{22}^T]^T$ is of full rank, and

$$\begin{bmatrix} B_1 & B_2 \\ D & 0 \\ 0 & 0 \end{bmatrix} = \begin{bmatrix} I & 0 \\ 0 & U^H \end{bmatrix} \begin{bmatrix} A_{12} \\ A_{22} \end{bmatrix} U,$$

it follows that B_2 is also of full rank. We have

$$\begin{bmatrix} \omega I & B_1 & \omega B_2 \\ B_1^H & D & 0 \\ \omega B_2^H & 0 & 0 \end{bmatrix} \begin{bmatrix} I & -\omega^{-1}B_1 & -\omega^{-1}B_2 \\ 0 & I & 0 \\ 0 & 0 & \omega^{-1}I \end{bmatrix}$$
$$= \begin{bmatrix} \omega I & 0 & 0 \\ B_1^H & D - \omega^{-1}B_1^H B_1 & -\omega^{-1}B_1^H B_2 \\ \omega B_2^H & -B_2^H B_1 & -B_2^H B_2 \end{bmatrix},$$
nich is easily seen to be nonsingular for large enough values of ω . \Box

which is easily seen to be nonsingular for large enough values of ω .

If we use the bound on F_k shown in (13), then we can deduce the following theorem.

Theorem 4.2. Let $(A - \sigma I)^{-1}(V_k + F_k) = V_{k+1}\underline{H}_k$ be of full rank and assume F_k is bounded as in (13) and $\sqrt{k\kappa}(V_k)\epsilon_1 < 1$. Then there is a ΔA of rank at most k such that

$$V_k = (A + \Delta A - \sigma I)V_{k+1}\underline{H}_k,$$

and

$$\|\Delta A\| \le \sqrt{k} \|A - \sigma I\| \frac{\kappa(V_k)\epsilon_1 + \kappa(V_{k+1})\kappa(\underline{H}_k)c_{kn}(\epsilon_2)}{1 - \sqrt{k}\kappa(V_k)\epsilon_1}$$

where $c_{kn}(\epsilon_2)$ is given by (14).

Proof. From $V_k + F_k = (A - \sigma I)V_{k+1}\underline{H}_k$ and $V_k = (A + \Delta A - \sigma I)V_{k+1}\underline{H}_k$ we see that any eligible ΔA has to satisfy $\Delta AV_{k+1}\underline{H}_k = -F_k$. We choose $\Delta A = -F_k(V_{k+1}\underline{H}_k)^{\dagger}$ (which is of rank at most k) implying $\|\Delta A\| \leq \|F_k\|/\sigma_{\min}(V_{k+1}\underline{H}_k)$. Substituting $||F_k||$ by the upper bound given in (13) yields

$$\begin{aligned} \|\Delta A\| &\leq \frac{\sqrt{k} \|V_k\| \epsilon_1 + \sqrt{k} \|A - \sigma I\| \|V_{k+1}\| \|\underline{H}_k\| c_{kn}(\epsilon_2)}{\sigma_{\min}(V_{k+1}\underline{H}_k)} \\ &\leq \frac{\sqrt{k} \|V_k\| \epsilon_1}{\sigma_{\min}(V_{k+1}\underline{H}_k)} + \sqrt{k} \|A - \sigma I\| \kappa(V_{k+1}) \kappa(\underline{H}_k) c_{kn}(\epsilon_2) \end{aligned}$$

For the denominator we get

$$\sigma_{\min}(V_{k+1}\underline{H}_k) \ge \sigma_{\min}((A + \Delta A - \sigma I)V_{k+1}\underline{H}_k)/||A + \Delta A - \sigma I||$$

$$\ge \sigma_{\min}(V_k)/(||A - \sigma I|| + ||\Delta A||),$$

where we used $\sigma_{\min}(XY) \leq ||X|| \sigma_{\min}(Y)$ which holds for any matrices X, Y. Thus

$$\|\Delta A\| \leq \frac{\sqrt{k} \|V_k\| (\|A - \sigma I\| + \|\Delta A\|)\epsilon_1}{\sigma_{\min}(V_k)} + \sqrt{k} \|A - \sigma I\| \kappa(V_{k+1})\kappa(\underline{H}_k)c_{kn}(\epsilon_2)$$

which can be reordered to the claimed bound.

If the linear systems are solved up to a normwise backward error $\epsilon_{\rm bw}$, and (8) and (15) hold for $2\epsilon_1 + \epsilon_2 = 3\epsilon_{\rm bw}$, then we get the following corollary.

Corollary 4.3. Let $(A - \sigma I)^{-1}(V_k + F_k) = V_{k+1}\underline{H}_k$ be of full rank and assume F_k is bounded as in (15) with $2\epsilon_1 + \epsilon_2 = 3\epsilon_{bw}$. Then there is a ΔA of rank at most k such that

$$V_k = (A + \Delta A - \sigma I)V_{k+1}\underline{H}_k,$$

and

$$\|\Delta A\| \le \sqrt{k} \|A - \sigma I\| \kappa(V_{k+1}) \kappa(\underline{H}_k) c_{kn}(3\epsilon_{\mathrm{bw}}),$$

where $c_{kn}(\cdot)$ is given by (14).

A few remarks are in order.

Remark 4.4. If $A + \Delta A - \sigma I$ in Theorem 4.2 and Corollary 4.3 is singular, then we can invoke Lemma 4.1 with $V = V_{k+1}\underline{H}_k$ to obtain a backward error $\Delta \widetilde{A}$, arbitrarily close to ΔA , such that $(A + \Delta \widehat{A} - \sigma I)^{-1}V_k = V_{k+1}\underline{H}_k$. The new backward error $\Delta \widehat{A}$ will in general have rank greater than k, but its numerical rank is still bounded by k. Here the definition of numerical rank can be arbitrarily strict, in the sense that we may define the numerical rank as the number of singular values that greater than $\epsilon > 0$, for an arbitrarily small ϵ .

Remark 4.5. If the orthonormalization is done properly, using, for instance, MGS with reorthogonalization, then $\kappa(V_{k+1}) \approx 1$. In this case we can ignore the factors $\kappa(V_{k+1})$ and $\kappa(V_k)$ when evaluating the bounds in Theorem 4.2 and Corollary 4.3. In particular this means that the bounds can be estimated cheaply as long as $||A - \sigma I||$ (or a good estimate of it) is known.

Remark 4.6. For the standard eigenvalue problem, shifts are used to find interior eigenvalues, so any sensible shift satisfies $|\sigma| \leq ||A||$. Thus, we have $||A - \sigma I|| \leq 2||A||$ in practice.

Remark 4.7. In view of [6], we note that our bounds do not contain the loss-oforthonormality term $||V_{k+1}^H V_{k+1} - I||$. Instead we saw that the condition number of the computed basis V_{k+1} plays a role in the bounds of the backward error. We note, however, that a small value of $||V_{k+1}^H V_{k+1} - I||$ implies that V_{k+1} is well-conditioned:

$$\|V_{k+1}^H V_{k+1} - I\| < \epsilon < 1 \quad \Rightarrow \quad \kappa(V_{k+1}) < \sqrt{\frac{1+\epsilon}{1-\epsilon}}.$$

The next example shows how Theorem 4.2 can be used to derive a simple a posteriori backward error bound.

Example 4.8. Suppose a matrix A and a shift σ with $|\sigma| < ||A||$ are given, and suppose we perform k steps of the shift-and-invert Arnoldi algorithm. To solve the linear systems we use an iterative method that employs (5) as stopping condition, that is, the linear systems are considered "solved" when the residuals are less than some tolerance ϵ_{tol} (we ignore the norm of the right hand side since it is approximately one). We use a rather crude tolerance so $\epsilon_{tol} \gg u$. For the orthogonalization we use MGS with one round of reorthogonalization so $c_{kn}(0) \leq 13ku$ (cf. Remark 3.4). If

(18)
$$\epsilon_{\text{tol}} \ge \kappa(\underline{H}_k) c_{kn}(0),$$

then Theorem 4.2, with $\epsilon_1 = \epsilon_{tol}$ and $\epsilon_2 = 0$, and the following remarks, yield that the computed quantities satisfy

$$(A + \Delta A - \sigma I)^{-1}V_k = V_{k+1}\underline{H},$$

where

(19)
$$\|\Delta A\| \le \frac{4\sqrt{k\kappa(V_{k+1})\epsilon_{\text{tol}}}}{1-\sqrt{k\kappa(V_k)\epsilon_{\text{tol}}}} \|A\|.$$

Here we have used the fact that $\kappa(V_{k+1}) \geq \kappa(V_k)$. Since MGS with reorthogonalization was employed, we expect $\kappa(V_{k+1})$ to be close to one. Thus, (19) tells us that the relative backward error $\|\Delta A\|/\|A\|$ is of roughly the same size as the tolerance we used to solve the linear systems. So, in this setting shift-invert-Arnoldi is backward stable.

We end this section with a numerical experiment. We consider the matrix

$$A = \begin{bmatrix} -2 & 1 & & \\ 1 & -2 & 1 & & \\ & 1 & \ddots & \ddots & \\ & & \ddots & \ddots & 1 \\ & & & 1 & -2 \end{bmatrix},$$

of order n = 1000, and the shift $\sigma = -2$. It is well-known that the spectrum of A is a subset of the interval (-4, 0), and the eigenvalues are given by $-2+2\cos(\pi k/(n+1))$, for k = 1:n. It follows that the shifted matrix $A - \sigma I$ is invertible and has norm $2\cos(\pi/(n+1)) \approx 2$.

We implemented the shift-and-invert Arnoldi algorithm in MATLAB R2013a. For orthonormalization we used MGS with one round of reorthogonalization. The matrix $A - \sigma I$ was stored in sparse format, and the linear systems was solved using MATLAB's "backslash" and 1u routines. We took k = 30 steps with the starting vector $[1, 1, \ldots, 1]^T$, and in each iteration we computed the backward error shown in (3), where the residual was evaluated in extended precision (32 digits) and then rounded to double precision. We did this using the vpa function from the Symbolic Math Toolbox. We also computed the error $F_k = V_k - (A - \sigma I)V_{k+1}\underline{H}_k$ in extended precision and rounded the result to double precision. For each j = 1:k, we computed

$$\mathcal{B}(\|\Delta A^{(j)}\|) := \sqrt{j} \|A - \sigma I\| \kappa(\underline{H}_j) c_{jn}(3\epsilon_{\mathrm{bw}}),$$

where $\epsilon_{\rm bw}$ was set to be the largest backward error of the linear systems that was encountered in the algorithm, and $c_{jn}(3\epsilon_{\rm bw}) := (3\epsilon_{\rm bw} + 13ju)/(1 - 13ju)$ (cf. Remark 3.4). As is mentioned in Remark 4.5, the above quantity is a good estimate of the bound in Corollary 4.3. We also evaluated the expression for the backward error, $\Delta A^{(j)} = -F_k(V_{k+1}\underline{H}_k)^{\dagger}$, given in the proof of Theorem 4.2, and estimated its norm. We did this using the MATLAB routines pinv (for the Moore-Penrose pseudo-inverse) and normest. The quantities $\mathcal{B}(\|\Delta A^{(j)}\|)$ and $\|\Delta A^{(j)}\|$ are shown in Figure 1 for j = 1:30. Even though the (estimated) upper bound $\mathcal{B}(\|\Delta A^{(j)}\|)$ can be seen to be rather pessimistic, it does show that the backward error is less than \sqrt{u} . In other words, by evaluating $\mathcal{B}(\|\Delta A^{(k)}\|)$ (which is cheap), we can deduce that the computation is backward stable up to single precision.

5. Further topics

5.1. **Implicit restarting.** The bounds in Theorem 4.2 and Corollary 4.3 contain the factor $\kappa(\underline{H}_k)$, so if $\kappa(\underline{H}_k) \gg 1$ we cannot guarantee a small backward error. If we recall how Arnoldi locates eigenvalues [23, pp. 257–265], we have, unfortunately, reason to suspect that this is the case. Since Arnoldi does not target the largest eigenvalues, but *any* isolated eigenvalue cluster, $H_k := [I_k \ 0]\underline{H}_k$ is likely to have both large and small eigenvalues, which suggests that H_k may be ill-conditioned. We will now show that the situation can be much better than expected if we restrict our attention to the largest eigenvalues of H_k , that is, the ones corresponding to eigenvalues of A closest to the shift σ . The idea is to do an implicit (thick)

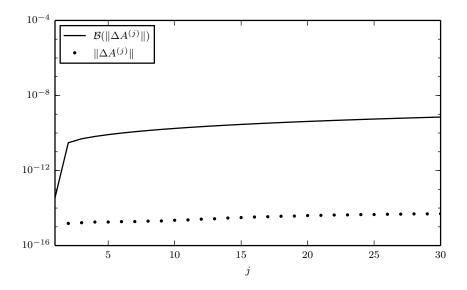


FIGURE 1. Computed backward errors and associated bound.

restart [22], and purge the small eigenvalues of H_k . Since small eigenvalues of H_k correspond to eigenvalues of A far from the shift σ , it is reasonable to assume they are of less interest. Suppose

$$(A - \sigma I)^{-1}(V_k + F_k) = V_{k+1}\underline{H}_k$$

and consider a Schur form $H_k = QTQ^H$ such that t_{ii} , $i = \ell + 1:k$, are the small eigenvalues to be purged. We have

$$(A - \sigma I)^{-1}(U_k + F_k Q) = [U_k \ v_{k+1}] \begin{bmatrix} T \\ h_{k+1,k} e_k^T Q \end{bmatrix}$$

where $U_k = V_k Q$. Throwing away the last $k - \ell$ columns yields

$$(A - \sigma I)^{-1}(U_{\ell} + F_k Q_{\ell}) = [U_{\ell} v_{k+1}] \begin{bmatrix} T_{\ell} \\ h_{k+1,k} e_k^T Q_{\ell} \end{bmatrix},$$

where $Q_{\ell} = Q(:, 1:\ell), U_{\ell} = U(:, 1:\ell)$ and $T_{\ell} = T(1:\ell, 1:\ell)$. Defining $u_{\ell+1} = v_{k+1}$,

$$\underline{T}_{\ell} = \begin{bmatrix} T_{\ell} \\ h_{k+1,k} e_k^T Q_{\ell} \end{bmatrix},$$

and $E_{\ell} = F_k Q_{\ell}$, results in a compact recurrence

(20)
$$(A - \sigma I)^{-1} (U_{\ell} + E_{\ell}) = U_{\ell+1} \underline{T}_{\ell},$$

where $||E_{\ell}|| \leq ||F_k||$. Note that our bound on E_{ℓ} depends on k and not ℓ . We can now repeat the proof of Theorem 4.2, and use the bounds $||E_{\ell}|| \leq ||F_k||$ and $\sigma_{\min}(U_{\ell+1}) \geq \sigma_{\min}(V_{k+1})$, and the recurrence (20) instead of the one assumed in the theorem. We get

$$U_{\ell} = (A + \Delta A - \sigma I)U_{\ell+1}\underline{T}_{\ell},$$

where

(21)
$$\|\Delta A\| \le \|A - \sigma I\| \frac{\sqrt{k\kappa(V_k)\epsilon_1} + \sqrt{k\kappa(V_{k+1})c_{kn}(\epsilon_2)} \|\underline{H}_k\| / \sigma_{\min}(\underline{T}_\ell)}{1 - \sqrt{k\kappa(V_k)\epsilon_1}}$$

Comparing this to the bound in Theorem 4.2 we see that $\kappa(\underline{H}_k)$ has been replaced by $\|\underline{H}_k\|/\sigma_{\min}(\underline{T}_{\ell})$. Further, it holds that

$$\|\underline{H}_k\|/\sigma_{\min}(\underline{T}_\ell) \le \|\underline{H}_k\|/\sigma_{\min}\left(\begin{bmatrix}T\\h_{k+1,k}e_k^TQ\end{bmatrix}\right) = \kappa(\underline{H}_k).$$

It follows that if \underline{H}_k is ill-conditioned due to the small eigenvalues we purged, then $\|\underline{H}_k\|/\sigma_{\min}(\underline{T}_\ell) \ll \kappa(\underline{H}_k)$ and (21) shows that the upper bound for the backward error corresponding to the part of the spectrum we care about is much smaller than the upper bound for the general backward error.

5.2. Hermitian backward errors. We now restrict the scope to the Hermitian matrix eigenvalue problem, that is, when $A = A^H$ and σ is real. Let us mention that we still consider the shift-and-invert Arnoldi algorithm, as it is shown in Algorithm 1, and *not* the shift-and-invert Lanczos algorithm with a three-term recurrence. In the Hermitian case, Algorithm 1 is also known as the shift-and-invert Lanczos algorithm with full orthogonalization, and it is used in, e.g., ARPACK[16, routine ssaitr.f] and MATLAB's eigs command.

Is it, for a Hermitian A, possible to find a Hermitian backward error ΔA ? We have seen in the proof of Theorem 4.2 that ΔA has to satisfy $\Delta AV_{k+1}\underline{\mathbf{H}}_k = -F_k$. Unfortunately the following lemma rules out existence of such a Hermitian ΔA in general.

Lemma 5.1. Let $X \in \mathbb{C}^{n \times k}$ and $F \in \mathbb{C}^{n \times k}$. Then there exists a Hermitian E with EX = F if and only if $X^H F$ is Hermitian and $FX^{\dagger}X = F$. In that case, there is such an E with rank $(E) \leq 2k$ and $||E||_* \leq 2||F||_*/\sigma_{\min}(X)$ where $|| \cdot ||_*$ denotes the 2-norm or the Frobenius norm.

Proof. The proof is simple and, for k = 1, is contained in [17]. We give it for completeness. Let E be any matrix such that EX = F. This implies $EXX^{\dagger}X = FX^{\dagger}X$ and (using $XX^{\dagger}X = X$) $EX = FX^{\dagger}X$, contradicting EX = F if $F \neq FX^{\dagger}X$. Thus $F = FX^{\dagger}X$ is necessary for the existence of an E with EX = F. Now, if E is Hermitian, then so is $X^{H}EX = X^{H}F$. Hence, if $X^{H}F$ is not Hermitian, then there is no Hermitian E with EX = F.

On the other hand, if $X^H F$ be Hermitian and $F = F X^{\dagger} X$, then

$$E := FX^{\dagger} + (FX^{\dagger})^H - X^{\dagger H}F^H XX^{\dagger} = FX^{\dagger} + (FX^{\dagger})^H (I - XX^{\dagger})$$

is also Hermitian. Furthermore, $\operatorname{rank}(E) \leq 2k$, EX = F, and (using that $I - XX^{\dagger}$ is an orthogonal projector)

$$||E||_* \le 2||FX^{\dagger}||_* \le 2||F||_* ||X^{\dagger}||_2 = 2||F||_* /\sigma_{\min}(X).$$

The next result shows that one still gets a Hermitian backward error if one replaces the Hessenberg matrix \underline{H}_k by some other $(k+1) \times k$ matrix \underline{G}_k . Before we state the theorem, we should clarify what we mean by "backward error" in this case. If we replace \underline{H}_k by something else, we cannot say that the computed quantities $(V_{k+1} \text{ and } \underline{H}_k)$ satisfy a an exact Krylov recurrence of a perturbed input matrix. We can, however, still say that the *computed subspace* is a Krylov subspace of a perturbed Hermitian input matrix. We refer to this Hermitian perturbation as the backward error.

Theorem 5.2. Let A be Hermitian and $(A - \sigma I)^{-1}(V_k + F_k) = V_{k+1}\underline{H}_k$. Suppose it holds for $\underline{G}_k \in \mathbb{C}^{(k+1)\times k}$ that $V_k^H V_{k+1} \underline{G}_k$ is Hermitian and $V_{k+1} \underline{G}_k$ is of full rank. Then there is a Hermitian ΔA of rank at most 2k such that

$$V_k = (A + \Delta A - \sigma I)V_{k+1}\underline{G}_k$$

and

$$\|\Delta A\| \le 2 \frac{\|(A - \sigma I)\| \|V_{k+1}\| \|\underline{H}_k - \underline{G}_k\| + \|F_k\|}{\sigma_{\min}(V_{k+1}\underline{G}_k)}.$$

Proof. From $V_k = (A + \Delta A - \sigma I)V_{k+1}\underline{G}_k$ and

$$V_k + F_k = (A - \sigma I)V_{k+1}\underline{H}_k = (A - \sigma I)V_{k+1}\underline{G}_k + (A - \sigma I)V_{k+1}(\underline{H}_k - \underline{G}_k)$$

we see that any eligible ΔA has to satisfy

$$\Delta A V_{k+1} \underline{G}_k = (A - \sigma I) V_{k+1} (\underline{H}_k - \underline{G}_k) - F_k = V_k - (A - \sigma I) V_{k+1} \underline{G}_k.$$

Since it is assumed that $V_{k+1}\underline{G}_k$ is of full rank, Lemma 5.1 implies that such a Hermitian ΔA exists if

$$(V_{k+1}\underline{G}_k)^H(V_k - (A - \sigma I)V_{k+1}\underline{G}_k) = (V_{k+1}\underline{G}_k)^H V_k - (V_{k+1}\underline{G}_k)^H (A - \sigma I)V_{k+1}\underline{G}_k$$

is Hermitian. Since the first term on the right hand side is Hermitian by assumption, this is easily seen to be the case. Also by Lemma 5.1, ΔA is bounded by

$$\begin{aligned} \|\Delta A\| &\leq 2\|(A - \sigma I)V_{k+1}(\underline{H}_k - \underline{G}_k) - F_k\|/\sigma_{\min}(V_{k+1}\underline{G}_k) \\ &\leq 2(\|(A - \sigma I)\|_2\|V_{k+1}\|\|\underline{H}_k - \underline{G}_k\| + \|F_k\|)/\sigma_{\min}(V_{k+1}\underline{G}_k), \end{aligned}$$
of rank at most 2k.

and is of rank at most 2k.

Remark 5.3. If $A + \Delta A - \sigma I$ is singular, then we can use the second part of Lemma 4.1 to find a *Hermitian* backward error $\Delta \tilde{A}$ arbitrarily close to ΔA such that $A + \Delta A - \sigma I$ is invertible.

In order to obtain a small Hermitian backward error, we need to find a matrix \underline{G}_k close to $\underline{\mathbf{H}}_k$ such that $V_k^H V_{k+1} \underline{G}_k$ is Hermitian. One possibility is

(22)
$$\underline{G}_k := R_{k+1}^{-1} \begin{bmatrix} T_k \\ h_{k+1,k} e_k^T \end{bmatrix} R_k,$$

where R_k, R_{k+1} are the upper triangular QR factors of V_k, V_{k+1} , respectively, and T_k is the real symmetric tridiagonal matrix with $t_{j+1,j} = t_{j,j+1} = h_{j+1,j}$ and $t_{j,j} = \Re(h_{jj})$. Then <u>G</u>_k is Hessenberg and computing Ritz pairs is particularly easy: we need to find vectors z and scalars μ such that

$$V_k^H (A + \Delta A - \sigma I)^{-1} V_k z = \mu V_k^H V_k z.$$

Here we have used Remark 5.3 in order to ensure that $A + \Delta A - \sigma I$ is invertible. By using the Krylov relation $(A + \Delta A - \sigma I)^{-1}V_k = V_{k+1}\underline{G}_k$ we obtain

$$V_k^H V_{k+1} \underline{G}_k z = \mu V_k^H V_k z.$$

Inserting the QR factorizations $V_j = Q_j R_j$, j = k, k + 1 and the formula for \underline{G}_k shown in (22) yields

$$R_{k}^{H}[I \ 0]R_{k+1}R_{k+1}^{-1}\begin{bmatrix}T_{k}\\h_{k+1,k}e_{k}^{T}\end{bmatrix}R_{k}z = \mu R_{k}^{H}R_{k}z,$$

which simplifies to $T_k \tilde{z} = \mu \tilde{z}$ where $\tilde{z} = R_k z$. So, the Ritz values are just the eigenvalues of T_k (which are real, since T_k is Hermitian). To obtain the Ritz vectors,

we would have to multiply \tilde{z} with R_k^{-1} . However, since R_k is close to the identity matrix if the orthogonalization has been done properly (for instance, by using MGS with reorthogonalization) we can approximate \tilde{z} by z. Thus, (approximations of) Ritz pairs for the choice (22) of \underline{G}_k can be obtained without computing R_k, R_{k+1} . We also note that choosing the eigenpairs of T_k to construct Ritz pairs is what is done in practice.

5.3. Conditions for breakdown. We now discuss how to derive a sensible breakdown criterion based on our error analysis. We saw in Section 1.1 that the computed quantities V_{j+1} and \underline{H}_k satisfy

$$(A - \sigma I)^{-1}(V_j + F_j) = V_{j+1}\underline{H}_j.$$

This recurrence can be rewritten as

$$(A - \sigma I)^{-1}(V_j + \widetilde{F}_j) = V_j H_j$$

where $\widetilde{F}_j = F_j - (A - \sigma I)h_{j+1,j}v_{j+1}e_j^T$. Note that the first j-1 columns of \widetilde{F}_j and F_j are identical. For the last column, we have

$$f_j = r_j - (A - \sigma I)(g_j + h_{j+1,j}v_{j+1}).$$

It is natural to declare breakdown when the error introduced by neglecting $h_{j+1,j}$ is of the same order as the errors that are present in the computation. This leads us to the following breakdown condition:

$$h_{j+1,j} < ||g_j|| + ||r_j|| / ||(A - \sigma I)v_{j+1}||$$

We can simplify this condition by replacing $||g_j||$ with its bound in (10). This yields

(23)
$$h_{j+1,j} < \eta(n,j) \|w_j\| u + \|r_j\| / \|(A - \sigma I)v_{j+1}\|.$$

We now discuss how to evaluate (23) in practice. Surely, if $h_{j+1,j} < \eta(n,j) ||w_j||u$ we can declare breakdown without further work. Otherwise we have to take the second term in (23) into consideration. If an iterative linear system solver that guarantees a residual less than some tolerance is used, then we can substitute $||r_j||$ in (23) by the given tolerance. If, for example, (6) is used as a stopping condition for the linear system solver, then $||r_j||$ is replaced by the right hand side of (6). If the residual, or any good bound for it, is not given, then we need to compute it. This is generally the case when the linear systems are solved by a direct method. Let mbe a constant such that the following forward error bound holds for an arbitrary vector x

$$\|\operatorname{float}((A - \sigma I)x) - (A - \sigma I)x\| \le mu \|A - \sigma I\| \|x\|.$$

If $A - \sigma I$ is given as a dense matrix, we have $m = n^{3/2}$ [11, p. 70]. For sparse matrices, m can be much smaller. The computed residual \hat{r}_i satisfies

$$\|\widehat{r}_{j}\| \leq (1+u)\|\text{float}((A-\sigma I)w_{j}) - v_{j}\| \\ \leq (1+u)(\|r_{j}\| + mu\|A-\sigma I\|\|w_{j}\|).$$

By comparing to (4), we recognize $||A - \sigma I|| ||w_j||mu$ as a part of the norm of a residual associated with a computed solution with corresponding backward error mu. Thus, we can compute a satisfactory \hat{r}_j if we use an extended precision \overline{u} such that $m\overline{u} < u$.

For the computation of the vector $(A - \sigma I)v_{i+1}$, we have

$$\begin{aligned} \|\text{float}((A - \sigma I)v_j) - (A - \sigma I)v_j\| &\leq mu\|A - \sigma I\|\|v_j\| \\ &\leq mu\kappa(A - \sigma I)\|(A - \sigma I)v_j\|, \end{aligned}$$

and, using the reverse triangle inequality, that

$$\|(A - \sigma I)v_j\| (1 - mu\kappa(A - \sigma I)) \le \|\text{float}((A - \sigma I)v_j)\|.$$

Thus the norm of the computed vector is accurate enough as long as $mu\kappa(A-\sigma I) \ll$ 1. If $(A-\sigma I)$ is so ill-conditioned so this is not satisfied, then we can use an extended precision \overline{u} such that $m\overline{u}\kappa(A-\sigma I) \ll 1$.

If (23) and (6) hold, then

$$\|\widetilde{f}_{j}\| \leq 2(\|v_{j}\|\epsilon_{1} + \|A - \sigma I\|\|w_{j}\|(\epsilon_{2} + \eta(n, j)u)).$$

By derivations similar to those leading to (13), we get

(24)
$$\|\widetilde{F}_j\| \leq 2\left(\sqrt{k}\|V_k\|\epsilon_1 + \sqrt{k}\|A - \sigma I\|\|V_{k+1}\|\|\underline{H}_k\|c_{kn}(\epsilon_2)\right).$$

From this we obtain the following "breakdown analogue" of Theorem 4.2.

Theorem 5.4. Let $(A - \sigma I)^{-1}(V_j + \widetilde{F}_j) = V_j H_j$ be of full rank and assume \widetilde{F}_k is bounded as in (24) and $\sqrt{k}\kappa(V_k)\epsilon_1 < 1$. Then there is a ΔA of rank at most k such that

$$V_k = (A + \Delta A - \sigma I)V_kH_k$$

$$\|\Delta A\| \le 2\sqrt{k} \|A - \sigma I\| \frac{\kappa(V_k)\epsilon_1 + \kappa(V_{k+1}) \|\underline{H}_k\| / \sigma_{\min}(H_k)c_{kn}(\epsilon_2)}{1 - \sqrt{k}\kappa(V_k)\epsilon_1}.$$

where $c_{kn}(\cdot)$ is given by (14).

The proof is omitted since it is essentially the same as the proof of Theorem 4.2. In a similar manner, we can get corresponding breakdown analogues to Corollary 4.3 and Theorem 5.2.

6. CONCLUSION

We have shown that a floating point implementation of shift-and-invert Arnoldi, where errors from all steps of the computation are taken into account, yields computed quantities that satisfy an exact shift-and-invert Krylov recurrence of a perturbed matrix. Here, the word "Krylov" is used instead of "Arnoldi" since the computed basis cannot be guaranteed to be perfectly orthogonal. We saw that the condition number of the computed basis V_{k+1} plays a role in the bounds of the backward error. Further, we have seen that the norm of the backward error ΔA depends on $\kappa(\underline{H}_k)$. We have seen that large $\kappa(\underline{H}_k)$ are acceptable if the linear systems are only solved to a loose tolerance (18). Otherwise we argued that even if this condition number is large, the restriction to the most important part of the recurrence (that is, what is left after purging the small eigenvalues of H_k) can have a small backward error.

For Hermitian matrices A, we have shown that there is a Hermitian backward error ΔA such that the computed basis, that is, the columns of V_{k+1} , spans a Krylov subspace associated with $A + \Delta A$. However, as in the case of standard Arnoldi [14], the small $(k+1) \times k$ matrix associated this subspace is generally not the computed Hessenberg matrix. Finally, we noted that our error analysis yields a sensible condition for when to declare breakdown. If this condition is met, we could derive a new set of backward error bounds, which show that an invariant subspace of a perturbed matrix has been found.

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