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Recyclage de sous-espaces de Krylov et réduction de sous-espaces de déflation pour résoudre une séquence de systèmes linéaires

Hussam Al Daas, Laura Grigori, Pascal Hénon, Philippe Ricoux

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Recyclage de sous-espaces de Krylov et réduction de sous-espaces de déflation pour résoudre une séquence de systèmes linéaires

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Résumé : Ce papier présente plusieures stratégies de déflation liées aux méthodes de recyclage de sous-espaces de Krylov pour résoudre une séquence de systèmes linéaires. À côté de stratégies de déflation très connues qui sont basées sur la déflation des vecteurs de Ritz et Ritz harmonique, on introduit une technique de déflation basée sur la décomposition en valeurs singulières. On considère deux contextes du recyclage, le recyclage de l'espace de Krylov entre les cycles de resart et le recylcage de l'espaces de déflation quand la matrice change dans la séquence. L'efficacité de la méthode proposée est étudiée sur des séquence de systèmes linéaires issues de la modélisation de réservoirs.

Mots-clés : espaces de Krylov, solveur linéaires itératifs, recyclage, déflation, réduction, svd

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Recycling Krylov subspaces and reducing deflation subspaces for solving sequence of linear systems

Abstract: This paper presents deflation strategies related to recycling Krylov subspace methods for solving one or a sequence of linear systems of equations. Besides well-known strategies of deflation, Ritz- and harmonic Ritz-based deflation, we introduce an SVD-based deflation technique. We consider the recycling in two contexts, recycling the Krylov subspace between the cycles of restarts and recycling a deflation subspace when the matrix changes in a sequence of linear systems. Numerical experiments on real-life reservoir simulations demonstrate the impact of our proposed strategy.

Key-words: Krylov subspaces, iterative linear solvers, recyling, deflation, reduction, svd

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1 Introduction

We consider the system

$$Ax = b \tag{1}$$

where $A \in \mathbb{C}^{n \times n}$ is non-singular and diagonalizable, $b \in \mathbb{C}^n$ is a right-hand side vector and $x \in \mathbb{C}^n$ is the vector of unknowns. We are interested in solving (1) by using a Krylov iterative method. GMRES [13], CG [5] and BiCG [9] are widely used Krylov subspace methods. Several works have studied the impact of the spectrum of the matrix A on the convergence rate of Krylov methods [9, 12, 8, 7, 17]. It was noticed that the small and nearly isolated eigenvalues induce a slow rate of convergence. In order to have robustness and fast convergence, a deflation subspace correction can annihilate the impact of this part of the spectrum. Different types of deflation techniques depending on the Hermitian structure of the matrix A are described in the literature [8, 3, 7]. These strategies construct a deflation subspace during the iterative solver. Usually, deflation subspaces in recycling methods are approximations of eigenspaces. In [3], the authors propose an approximation by using the Ritz vectors. To preserve the structure of the Arnoldi procedure, Morgan presented the deflation of harmonic Ritz vectors [8]. Few papers discuss the impact of singular vectors on the convergence of Krylov methods. In [15], Simoncini shows the impact of very small singular values on the convergence of several restarted variants of Krylov methods when the right-hand side has a large coefficient in the direction of the left singular vector. This work inspired us to think about deflation of approximate singular vectors rather than eigenvectors.

In several applications, we have to solve successive linear systems $A_i x = b_i$ where $A_i \approx A_{i+1}$, e.g., large eigenvalue problems or the Newton's method for solving nonlinear systems. Furthermore, the Newton's method combined with a Constrained Pressure Residual (CPR [18, 14]) solver have both cases, $A_i \approx A_{i+1}$ and $A_i = A_{i+1}$. Thus, the deflation subspace that is built during the solution of $A_i x = b_i$ can be useful to solve $A_{i+1} x = b_{i+1}$. In [10], the authors present the method GCRO-DR that benefits from previous solutions when solving a sequence of linear systems. In the Hermetian case, Vuik et al. [17] prove that for any deflation subspace the effective condition number of the preconditioned matrix without deflation. However, this is not true for the non-symmetric case. Moreover, we will see in the numerical experiments 7 how the deflation subspace, in several cases might deteriorate the convergence of the method. Thus, when considering a new matrix in the sequence of linear systems, it is important to check whether the deflation subspace is appropriate or not. For this aim, we introduce different strategies to reduce the deflation subspace based on the approximation type of the deflation.

We introduce the deflation based on singular vectors. We generalize the formulation of exact deflation for a symmetric linear system given in [17] to general linear systems. Then we derive the approximation strategy for deflating the singular vectors. We compare the combination of the reduction of the deflation subspace and the deflation of singular vectors to the existing method GCRO - DR on a sequence of linear systems arising from reservoirs simulations. Since we are interested in solving a sequence of linear systems, we follow the presentation of Parks et al. in [10].

The paper is organized as follows. In the following section, we give a brief review about the Krylov subspace and the Arnoldi procedure, and then we present the deflation subspace correction. Afterwards, we review the method GCRO-DR [10] with an abstract deflation subspace related to the previous search subspace. In Section 3, we introduce the deflation of singular vectors. We generalize the formulation of deflation of exact eigenspaces in the Hermitian case [17] to the non-symmetric case. Section 4 presents strategies of recovering deflation vectors for GCRO-DR. At first, we recall the Ritz pairs and the harmonic Ritz pairs, and then we introduce a strategy of deflating approximate singular vectors. In Section 5, we present the reduction of the deflation subspace in recycling Krylov subspace methods. For each type of deflated vectors we propose a reduction strategy. Section 6 details the parallel implementation of the method and analyzes its computation and communication cost. In Section 7, we compare the three variants of deflation techniques combined with reduction of deflation subspace on a sequence of linear systems arising from reservoirs simulations.

2 Background

In this section, we review the Krylov subspaces, the Arnoldi procedure, and the exact deflation of eigenvectors. We review the GCRO-DR algorithm to which we propose a modification in the following sections.

Notation

For an integer m > 0 we refer to the $m \times m$ identity matrix as I_m . Let the integers m > 0, p > 0, and let $h_{ij} \in \mathbb{C}$, for $i = 1, \ldots, m, j = 1, \ldots, p$, we refer to the matrix whose (i, j) coefficient is $h_{i,j}$ for $i = 1, \ldots, m, j = 1, \ldots, p$ as $(h_{ij})_{\substack{1 \leq i \leq m \\ 1 \leq j \leq p}}$. If for some pair (i, j) such that $1 \leq i \leq m$ and $1 \leq j \leq p$ the element $h_{i,j}$ is not defined, it is set to zero. Let $H = (h_{i,j})_{\substack{1 \leq i \leq m \\ 1 \leq j \leq p}} \in \mathbb{C}^{m \times p}$, we refer to the matrix $(h_{i,j})_{\substack{1 \leq i \leq m-1 \\ 1 \leq j \leq p}} \in \mathbb{C}^{(m-1) \times p}$ as \overline{H} . The matrix $(h_{i,j})_{\substack{2 \leq i \leq m \\ 1 \leq j \leq p}} \in \mathbb{C}^{(m-1) \times p}$ as \underline{H} .

2.1 Krylov subspaces and Arnoldi procedure

The Krylov subspace of dimension k > 0 associated to A and b is defined as

$$K_k(A,b) = \operatorname{span}\left\{b, Ab, \dots, A^{k-1}b\right\}.$$
(2)

Let

$$k_{A,b} = \min_{\substack{P \in \mathcal{P} \\ P(0) = 1}} \{ deg(P) \mid P(A)b = 0 \},\$$

where \mathcal{P} represents the set of polynomials with complex coefficients. The vectors $b, Ab, \ldots, A^{k-1}b$ for $k \leq k_{A,b}$ form a basis of the Krylov subspace. Due to round-off errors, an orthogonalization procedure is necessary to construct a basis vectors for the Krylov subspace. To this aim the Arnoldi procedure is widely used, see Algorithm 1.

According to the notations in Algorithm 1 the following relation holds at each iteration k

$$AV_k = V_{k+1}H_k,\tag{3}$$

where $V_k = \{v_1, \ldots, v_k\}, V_{k+1} = [V_k, v_{k+1}], \text{ and } H_k = (h_{i,j})_{\substack{1 \le i \le k+1 \\ 1 \le j \le k}}$

2.2 Deflation of eigenvectors

Let S be an A-invariant subspace of dimension $k \ge 0$ related to the smallest eigenvalues of the matrix A. Let $Z \in \mathbb{C}^{n \times k}$ be a matrix whose vectors form an orthonormal basis of the subspace S. Let T be the projection of A on the subspace S, i.e., $T = Z^H A Z$. We note that the following relation holds

$$AZ = ZT$$

Algorithm 1 Arnoldi (A, v_1, m)

Require: normal vector $v_1 \in \mathbb{C}^n$, matrix $A \in \mathbb{C}^{n \times n}$, number of iterations m **Ensure:** orthonormal basis vectors V_m , Hessenberg matrix $H_m \in \mathbb{C}^{(m+1) \times m}$ 1: for j = 1 : m do 2: $w = Av_i$ 3: 4: end for 5: $w = w - \sum_{i=1}^{j} h_{ij} v_i$ $h_{j+1,j} = \|w\|_2$ 6: 7: if $h_{j+1,j} = 0$ then 8: Set $m = j, v_{m+1} = 0$ 9: 10: else $v_{j+1} = \frac{w}{h_{j+1,j}}$ 11: 12:end if $V_m = [v_1, \cdots, v_m], V_{m+1} = [V_m, v_{m+1}], H_m = (h_{ij})_{i,j}$ 13: 14: **end for**

Let $\{\lambda_1, \ldots, \lambda_n\}$ be the eigenvalues of A ordered increasingly by the magnitude of their absolute value, i.e., $|\lambda_1| \leq \ldots \leq |\lambda_n|$. Consider the preconditioned system

$$(I + Z(T^{-1} - I)Z^H)Ax = (I + Z(T^{-1} - I)Z^H)b$$
(4)

The following Theorem 1 can be found in [3]. It describes how to deflate the eigenvalues of the matrix A by adding a low-rank correction.

Theorem 1 The eigenvalues of the matrix $A + Z(T^{-1} - I)Z^H A$ are $\{1, \ldots, 1, \lambda_{k+1}, \ldots, \lambda_n\}$.

The equation (4) can be rewritten as :

$$\left(\left(A - ZZ^{H}A\right) + ZT^{-1}Z^{H}A\right)x = \left(\left(b - ZZ^{H}b\right) + ZT^{-1}Z^{H}b\right)$$

In other words, to solve (1), we look for the solution in two subspaces, the first is inside S, the second is outside S. The first finds the part of the solution that is generated by the basis vectors V. This is done by solving the deflation subspace problem with the matrix T. The second finds the part that is orthogonal to S. This part is left to the Krylov method. Since the eigenvectors of the matrix are not orthogonal in general, we need at each iteration of the Krylov method to correct the new computed basis vector. In [17], the authors present how to avoid doing this correction at each iteration when the considered matrix is Hermitian Positive Definite (HPD).

2.3 Ritz pairs

In this paragraph we review the definition of Ritz pairs of a matrix related to a subspace.

Definition 1 Let $B \in \mathbb{C}^{n \times n}$ and let S be a subspace of dimension k < n. We say that the pair $(u, \theta) \in \mathbb{C}^n \times \mathbb{C}$ is a Ritz pair of B related to the subspace S if

$$\begin{cases} u \in \mathcal{S}, \\ y^H(Bu - \theta u) = 0, \forall y \in S. \end{cases}$$
(5)

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In the case where the subspace $S = K_k(A, b)$, the Ritz pairs of the matrix A correspond to

$$\mathcal{R} = \{ (V_k u, \theta), \overline{H}_k u = \theta u \},\$$

see Algorithm 1. If we consider the subspace $S = AK_k(A, b)$, the Ritz pairs of the matrix A^{-1} correspond to

$$\mathcal{R} = \{ (V_k u, \theta), \overline{H}_k^H u = \theta H_k^H H_k u \}$$

Lemma 1 Let $B \in \mathbb{C}^{m \times m}$ and let S be a subspace of dimension k < m that contains an eigenpair (λ, u) of the matrix B. Let $V \in \mathbb{C}^{m \times k}$ be a matrix whose columns form an orthonormal basis of S. Then, there exists a vector $w \in \mathbb{C}^k$ such that (λ, w) is an eigenpair of $V^H BV$, the restriction of B to the subspace S. Furthermore, $w = V^H u$.

Proof 1 The proof is simple and direct.

In Section 4, we will review in detail the derivation of these Ritz pairs. Furthermore, we derive the Ritz pairs relative to the singular vectors and values of A.

In the following, we review the GCRO-DR method and the construction of an abstract deflation subspace related to it.

2.4 GCRO-DR

We recall that CGRO-DR is an iterative method for solving a sequence of linear systems which takes advantage of the history of previous search subspaces by performing recycling. In [10], the authors present the algorithm with a deflation subspace related to the harmonic Ritz vectors. We explain briefly the method. Given the sequence of linear systems $A_i x = b_i$, for $i = 1, \ldots, p$, where p > 1, the method GCRO-DR(m, k) works as the following. To solve the linear system $A_1 x = b_1$, the first cycle is performed as a usual GMRES cycle. The subspace relation that holds at the end of the cycle is the usual Arnoldi relation :

$$A_1 V_m = V_{m+1} H_m.$$

The solution x_1 and the residual r_1 are computed as usual. Based on a given strategy, a column matrix $P_k \in \mathbb{C}^{m \times k}$ is computed (Section 4 details the computation of this matrix following different strategies). This matrix is expanded by V_m to form the basis of the deflation subspace $\tilde{Y}_k = V_m P_k$. The image of the deflation subspace by A_1 is computed

$$A_1 V_m P_k = V_{m+1} H_m P_k,$$

then orthonormalized,

$$C_k = V_{m+1}Q,$$

where Q stands for the unitary factor of the QR factorization of $H_m P_k = QR$. The relation that holds at this stage is :

$$A_1 Y_k R^{-1} = C_k.$$

We note $U_k = \tilde{Y}_k R^{-1}$. In order to avoid round-off errors, the vectors of U_k are normalized. We note D_k the diagonal matrix such that the following holds

$$\tilde{U}_k = U_k D_k^{-1},$$
$$A_1 \tilde{U}_k = C_k D_k.$$

At that moment, the method is ready to continue to the next cycle, m - k Arnoldi iterations are performed with $(I - C_k C_k^H) A_1$ as an operator and the normalized residual as the starting vector. The following relation holds

$$A_1[U_k, V_{m-k}] = [C_k, V_{m-k+1}]G_m,$$

where V_{m-k+1} is the resulting orthonormal basis constructed by the Arnoldi procedure, $G_m = \begin{pmatrix} D_k & B_{m-k} \\ 0 & H_{m-k} \end{pmatrix}$, $B_{m-k} = C_k^H A_1 V_{m-k}$ and H_{m-k} is the resulting Hessenberg matrix of the Arnoldi procedure. The solution x and the residual r are computed following the constraint $||r_2||_2 = ||r_1 - A_1 x||_2$ has a minimal norm over the subspace spanned by the columns of $[\tilde{U}_k, V_{m-k}]$. The extraction of the deflation subspace is performed as before by replacing the matrix H_m by G_m . This cycle is repeated until the convergence is achieved. In order to solve the second linear system, all what is necessary is to update the image of the deflation subspace, i.e., C_k , by the image of the deflation subspace by the new matrix A_2 . Algorithm 2 presents the GCRO-DR method with an abstract deflation subspace. We will discuss the deflation variants in Section 4.

Considering the notations in Algorithm 2, if the matrix \tilde{Y}_k is not defined, then (3) holds. Otherwise, the following relation holds

$$A[\tilde{U}_k, V_{m-k}] = [C_k, V_{m-k+1}] \begin{pmatrix} D_k & B_{m-k} \\ 0 & H_{m-k} \end{pmatrix}.$$
 (6)

The matrix $[C_k, V_{m-k+1}]$ is unitary, but this is not true, in general, for the matrix $[\tilde{U}_k, V_{m-k}]$. In this case the subspace generated by the columns of \tilde{U}_k, V_{m-k} is not a Krylov subspace.

In the following section we discuss the deflation of exact singular vectors.

3 Deflation based on singular vectors

As previously mentioned, the work of Simoncini on the convergence of several restarted Krylov methods and its relation with the singular vectors associated to the smallest singular values [15] motivated us to investigate the deflation based on singular vectors approximation. This section introduces the deflation of singular vectors. Given a set of singular vectors (right or left), we show how to deflate these vectors during the Krylov method.

Theorem 2 Let x_* be the exact solution of (1). Let

$$A = U\Sigma V^H$$

be the singular value decomposition of A such that the singular values are ordered increasingly. Let k_{τ} be the number of singular values smaller than a given threshold $\tau > 0$. Consider the splitting of the SVD decomposition

$$A = U_1 \Sigma_1 V_1^H + U_2 \Sigma_2 V_2^H,$$

where $\Sigma_2 \in \mathbb{C}^{k_\tau \times k_\tau}$ is a diagonal matrix whose diagonal elements are the singular values smaller than τ . Consider \tilde{x} an approximate solution of the following linear system of equations

$$(I - U_2 U_2^H) A x = (I - U_2 U_2^H) b, (7)$$

such that $\|\hat{x} - \tilde{x}\|_2 \leq \varepsilon$, where \hat{x} is an exact solution of (7) and $\varepsilon > 0$. Then, the following holds

$$||x_* - (I - V_2 V_2^H)\tilde{x} - V_2 \Sigma_2^{-1} U_2^H b||_2 \le \varepsilon.$$

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Algorithm 2 GCRO-DR

Require: the maximum dimension of the search subspace m, the dimension of the deflation subspace k, the convergence tolerance ε , the initial guess x_0 .

Ensure: approximate solution \tilde{x} of Ax = b. 1: initialization $r_0 = b - Ax_0, i = 1$ 2: if Y_k is defined (from solving a previous linear system) then let [Q, R] be the QR factorization of AY_k 3: $\begin{aligned} C_k &= Q\\ U_k &= \tilde{Y}_k R^{-1} \end{aligned}$ 4: 5:6: 7:8: else $v_1 = r_1 / ||r_1||$ 9: $c = ||r_0||_2 e_1$ 10: perform m steps of GMRES, solving min $||c - H_m y||_2$ for y and generating V_{m+1} and H_m 11: 12: $x_1 = x_0 + V_m y$ $r_1 = V_{m+1}(c - H_m y)$ 13:compute k deflation vectors P_k 14: $Y_k = V_m P_k$ 15:let [Q, R] be the QR factorization of $H_m P_k$ 16: $C_k = V_{m+1}Q$ 17: $U_k = \tilde{Y}_k R^{-1}$ 18:19: end if 20: while $||r_i||_2 > \varepsilon$ do i = i + 121: perform m - k Arnoldi steps with the linear operator $(I - C_k C_k^H)A$, letting $v_1 =$ 22: $|r_{i-1}/||r_{i-1}||_2$ and generating V_{m-k+1}, H_{m-k} , and $B_{m-k} = C_k^H A V_{m-k}$ let D_k be a diagonal scaling matrix such that $\tilde{U}_k = U_k D_k$, where the columns of \tilde{U}_k have 23: a unit norm $\hat{V}_m = [\hat{U}_k, V_{m-k}], \, \hat{W}_{m+1} = [C_k, V_{m-k+1}]$ 24: $G_m = \begin{pmatrix} D_k & B_{m-k} \\ 0 & H_{m-k} \end{pmatrix}$ solve $min \|G_m y - \hat{W}_{m+1}^H r_{i-1}\|_2$ for y25:26: $x_i = x_{i-1} + V_m y$ 27: $r_i = r_{i-1} - \hat{W}_{m+1}G_m y$ 28:compute k deflation vectors P_k 29: $\tilde{Y}_k = \hat{V}_m P_k$ 30: let [Q, R] be the QR factorization of $G_m P_k$ 31: $C_k = \hat{W}_{m+1}Q$ $U_k = \tilde{Y}_k R^{-1}$ 32: 33: 34: end while

35: let $\tilde{Y}_k = U_k$

Proof 2 First, we remark that x_* is a solution of (7) and the set of solutions of (7) can be written as $S = \{x = x_* + V_2u, u \in \mathbb{C}^{k_\tau}\}$. Indeed, let x be a solution of (7). Then,

$$Ax = (I - U_2 U_2^H)b + U_2 U_2^H Ax, Ax = Ax_* - U_2 U_2^H (b - Ax).$$

We multiply by A^{-1} both sides. We obtain

$$x = x_* + V_2 \Sigma_2^{-1} U_2^H (Ax - b)$$

= $x_* + V_2 \Sigma_2^{-1} (\Sigma_2 V_2^H x - U_2^H b)$
= $x_* + V_2 V_2^H x - V_2 \Sigma_2^{-1} U_2^H b.$

On the other side, each element of S is a solution of (7). We have

$$V_2 V_2^H x_* = V_2 V_2^H A^{-1} b$$

= $V_2 \Sigma_2^{-1} U_2^H b.$

Hence, we can write

$$\begin{aligned} x_* &= (I - V_2 V_2^H) x_* + V_2 V_2^H x_*, \\ &= (I - V_2 V_2^H) \hat{x} + V_2 \Sigma_2^{-1} U_2^H b, \end{aligned}$$

thus,

$$x_* - (I - V_2 V_2^H)\tilde{x} - V_2 \Sigma_2^{-1} U_2^H b = (I - V_2 V_2^H)(\hat{x} - \tilde{x}).$$

Finally, we obtain

$$\begin{aligned} \|x_* - (I - V_2 V_2^H) \tilde{x} - V_2 \Sigma_2^{-1} U_2^H b \|_2 &\leq \|(I - V_2 V_2^H) (\hat{x} - \tilde{x})\|_2, \\ &\leq \|\hat{x} - \tilde{x}\|_2, \\ &\leq \varepsilon. \end{aligned}$$

Note that the matrix $(I - U_2 U_2^H)A$ is singular. However, solving the linear system of equations (7) with a Krylov method is possible. The approximate solution of (1) can be recovered as

$$x = (I - V_2 V_2^H)\tilde{x} + V_2 x_2,$$

where $x_2 = \sum_{i=1}^{n-1} U_2^H b$ and \tilde{x} is the approximate solution given by the Krylov method.

The previous discussion is a generalization of the theory of deflation of eigenvectors for an SPD matrix to a general invertible matrix developed in [17]. In the next section, we discuss different approaches for recovering deflation information from a Krylov subspace. Two approaches are based on approximating eigenvectors of the matrix, and these approaches can be found in literature [8, 10]. A third approache is introduced here and it is based on approximating the right singular vectors of the matrix.

4 Recovering deflation vectors

In this section, we derive three different types of deflation subspaces. These subspaces are considered in the context of GCRO-DR [10]. In Section 2, we reviewed the method GCRO-DR for

an abstract deflation subspace issued from the last Krylov subspace. Even though any deflation vectors can be chosen for GCRO-DR, this choice has an impact on the convergence of the method. Two types of deflation vectors have been presented in previous works [8, 3, 10]. They are based on the smallest and largest Ritz values of the matrices A, A^{-1} respectively. During the GCRO-DR method, one of the relations (3) and (6) holds. In the following, we present how to recover the deflation vectors mentioned above. These strategies were introduced in previous works [8, 3, 10]. We present them for the sake of completeness. Afterwards, we introduce the deflation vectors that are related to the right singular vectors.

We consider the following environment. Let K be a search subspace (Krylov subspace or else) of dimension m associated to the matrix A, let $V \in \mathbb{C}^{n \times m}$ be a matrix whose columns form a basis of K. Let $W \in \mathbb{C}^{n \times (m+1)}$ be a unitary matrix such that the following relation holds

$$AV = WG,\tag{8}$$

where $G \in \mathbb{C}^{(m+1) \times m}$ is a Hessenberg matrix.

We will derive the eigenvalue problem that is solved in order to compute the deflation vectors. For each approximation strategy, we formulate the computation following the notations in Algorithm 2. Thus, we will distinguish two cases in each strategy, depending on whether the deflation subspace already exists or not.

4.1 Approximation based on the smallest Ritz values of A

In this case, the Ritz pairs (u, θ) verify

$$\begin{cases} u \in K, \\ y^H(Au - \theta u) = 0, \forall y \in K \end{cases}$$

Since the columns of V span the subspace K, we have

$$V^H(AVw - \theta Vw) = 0,$$

where u = Vw. This can be written as

$$V^H W G w = \theta V^H V w.$$

By adapting the notations in Algorithm 2, the deflation vectors can be recovered by solving the following eigenvalue problem

$$\begin{cases} \overline{H}_{m}w = \theta w, & \text{if } \tilde{Y}_{k} \text{ is not defined} \\ \begin{pmatrix} \tilde{U}_{k}^{H}C_{k} & \tilde{U}_{k}^{H}V_{m-k+1} \\ 0 & (I_{m-k}, 0) \end{pmatrix} G_{m}w = \theta \begin{pmatrix} \tilde{U}_{k}^{H}\tilde{U}_{k} & \tilde{U}_{k}^{H}V_{m-k} \\ V_{m-k}^{H}\tilde{U}_{k} & I_{m-k} \end{pmatrix} w, & \text{if } \tilde{Y}_{k} \text{ is defined.} \end{cases}$$
(9)

4.2 Approximation based on the largest Ritz values of A^{-1}

To approximate the eigenvectors of A^{-1} we look for the approximate vectors in the subspace AK. We write

$$\begin{cases} u \in AK, \\ y \perp (A^{-1}u - \theta u), \forall y \in AK, \end{cases}$$

The columns of AV span the subspace AK. Hence, we have

$$V^H A^H (Vw - \theta A Vw) = 0,$$

this can be written as

$$G^H W^H V w = \theta G^H G w.$$

By using the notations of Algorithm 2, this requires solving the following generalized eigenvalue problem

$$\begin{cases} \overline{H}_{m}^{H}w = \theta H_{m}^{H}H_{m}w, & \text{if } \tilde{Y}_{k} \text{ is not defined} \\ C_{k}^{H}\tilde{U}_{k} & 0 \\ V_{m-k+1}^{H}\tilde{U}_{k} & \begin{pmatrix} I_{m-k} \\ 0 \end{pmatrix} \end{pmatrix} w = \theta G_{m}^{H}G_{m}w, & \text{if } \tilde{Y}_{k} \text{ is defined.} \end{cases}$$
(10)

We note that Morgan, in his paper [8], refers to the values θ^{-1} as the harmonic Ritz values.

4.3 Approximation based on the smallest Ritz values of $A^{H}A$

We follow the definition of the Ritz pairs given in Definition 1. In order to approximate the left singular vectors of the matrix A, we consider an approximation of the eigenvectors of the matrix $A^H A$. Thus, we are looking for the pair (u, θ) such that

$$\begin{aligned} & u \in K, \\ & y \perp (A^H A u - \theta u), \forall y \in K \end{aligned}$$

The columns of V span the subspace K, the previous relation can be written as

$$V^H(A^HAVw - \theta Vw) = 0,$$

where u = Vw. This can be written as

$$G^H G w = \theta V^H V w.$$

The formulation of this eigenvalue problem in the context of Algorithm 2 is as follows

$$\begin{cases} H_m^H H_m w = \theta w, & \text{if } \tilde{Y}_k \text{ is not defined} \\ G_m^H G_m w = \theta \begin{pmatrix} \tilde{U}_k^H \tilde{U}_k & 0 \\ 0 & I_{m-k} \end{pmatrix} w, & \text{if } \tilde{Y}_k \text{ is defined.} \end{cases}$$
(11)

5 Deflation subspace reduction

In this section we introduce a strategy to reduce the dimension of the deflation subspace. Consider a sequence of linear systems solved by GCRO-DR. This method builds or updates a deflation subspace during the solution of each linear system in the sequence. This deflation subspace is used in the solution of the following system. Since the matrices in the sequence can differ, updating the deflation subspace, which is constructed during the previous system, may not be useful anymore. When the matrices are Hermitian, the worst case that can happen is that the deflation subspace does not improve the convergence; see [17]. However, this is not the case for non-Hermitian matrices, where an improper choice of the deflation subspace may even deteriorate the convergence. Thus, one needs a strategy to reduce the dimension of the deflation subspace judiciously if necessary. Such strategy should keep in the reduced deflation subspace a vector which well approximates an eigenvector of $(A, A^H A, \text{ or } A^{-1})$ depending on the deflation strategy. The motivation for this constraint is the following. Let B be the matrix related to the deflation strategy $(B = A, A^H A, \text{ or } A^{-1})$. Suppose that the deflation subspace contains an exact eigenvector u of B associated to the eigenvalue θ that is smaller than $(\tau, \tau^2, \text{ or } \frac{1}{\tau}, \text{ respectively})$, where $\tau > 0$ is a predefined value. Lemma 1 proves that the projection of the vector u to the deflation subspace is an eigenvector of the projected matrix to the deflation subspace. The threshold τ can be chosen to be relative either with respect to the largest approximated eigenvalue or the largest approximated singular value depending on the deflation strategy. Hence, for each strategy of recycling, presented in Section 4, the reduction strategy is associated to it, i.e., if the recycling strategy deflates the Ritz vectors associated to the small (resp. large) Ritz values (corresponding to the search subspace), the reduction strategy disregards the Ritz vectors associated to the large (resp. small) Ritz values (corresponding to the subspace). The reduced subspace is a subset of the original deflation subspace.

In the following, we derive the generalized eigenvalue problems that are necessary to perform the reduction of the deflation subspace. We suppose that the columns of the unitary matrix Zspan the deflation subspace S. Table 1 presents the eigenvalue problems to solve in order to reduce the deflation subspace corresponding to each strategy of deflation.

Reference matrix	Eigenvalue problem	Disregarded part
A	$Z^H A Z w = \theta w$	largest magnitude
A^{-1}	$Z^H A^H Z w = \theta Z^H A^H A Z w$	smallest magnitude
$A^H A$	$Z^H A^H A Z w = \theta w$	largest magnitude

TABLE 1 – Generalized eigenvalue problems reducing the deflation subspace, *Reference matrix* stands for the matrix whose eigenvectors are approximated, *Eigenvalue problem* refers to the problem to solve in order to reduce the deflation subspace, *Disregarded part* refers to the part of the spectrum to which are associated the disregarded vectors.

We note that for the case of an approximation of singular vectors it is more robust to perform an SVD factorization of AZ rather than solving the eigenvalue problem $Z^H A^H A Z w = \theta w$.

Algorithm 3 Deflation subspace reduction

Require: matrix A, deflation basis vectors \tilde{Y} , threshold τ , def the matrix to which the deflation subspace is related

Ensure: reduced deflation basis vectors \tilde{Y} 1: compute the QR factorization of $\tilde{Y} = ZR$ if def = A then 2: solve $Z^H A Z w = \theta w$ 3: 4: else if $def = A^{-1}$ then solve $Z^{H}A^{H}AZw = \theta Z^{H}A^{H}Zw$ 5: else if $def = A^H A$ then 6: solve $Z^H A^H A Z w = \theta^2 w$ $7 \cdot$ end if 8: 9: form Y whose columns are Zw where w is an eigenvector associated $\theta < \tau$

Here, we propose the method Generalized Minimal Residual with Modified Deflated Restarting (GMRES-MDR), Algorithm 4. This method is based on the GCRO-DR method. The



FIGURE 1 – Deflation and search subspace basis vectors storage for GMRES-MDR method

main difference is that an estimation of the previous deflation subspace (when the matrix in the sequence changes) is performed in order to keep only necessary vectors.

6 Parallel design of GMRES-MDR

In this section, we describe the parallel design of the method GMRES-MDR. At first we describe the distribution of data over processors. Then, we analyse the cost of the method and compare the three variations related to the deflation that we use in our implementation.

6.1 Data distribution

Matrices and vectors in our implementation can be separated into two types, distributed and redundant. The matrix A and the search subspace (Krylov and deflation) vectors are distributed in block row wise. All matrices and vectors whose both dimensions are less than or equal to the dimension of the search subspace are stored redundantly on each processor. Using this distribution, the application of the sparse operator requires communication between neighbour processors. Only dot product operations require global communication.

In the following, we describe how data is stored in memory.

6.1.1 Memory management

Dense matrices in iterative solvers are usually stored in one of two ways, either in column major or in row major. We use column major format due to constraints in the reservoir simulator. The vectors of the search subspace are stored in *full storage scheme* since they do not have a special structure. We refer the reader to [6] for more details on storage scheme structures. In our implementation, all dense matrices are stored in full storage scheme. Figure 1 shows how the vectors of the basis are stored in memory. This order $\tilde{U}_k, C_k, V_{m-k}$, in column ordering, has the following advantage. The orthogonalization steps in the Arnoldi procedure and the deflation correction can be done simultaneously. Sparse matrices are stored in the Compressed Sparse Row (CSR) storage scheme.

6.2 Parallel interaction environment and implementation language

We use message passing interface MPI [16] for communication between processors and we use OpenMP for multithreading. These two libraries are compatible with the language C that we chose to use.

Algorithm 4 GMRES-MDR

Require: the maximum dimension of the search subspace m, the maximum dimension of the deflation subspace k_{max} , the convergence tolerance ε , the initial guess x_0 , the matrix to which the deflation subspace is related def, τ threshold of reduction of the deflation subspace. **Ensure:** approximate solution \tilde{x} of Ax = b. 1: initialization $r_0 = b - Ax_0, i = 1, k = k_{max}$ 2: if Y_k is defined (from solving a previous linear system) then call Algorithms 3 and get Y_k , where k is the number of columns in Y_k 3: 4: let [Q, R] be the QR factorization of AY_k $\begin{array}{l} C_k = Q \\ U_k = \tilde{Y}_k R^{-1} \end{array}$ 5: 6: $x_1 = x_0 + U_k C_k^H r_0$ 7: $r_1 = r_0 - C_k C_k^{\tilde{H}} r_0$ 8: 9: else 10: $v_1 = r_1 / ||r_1||$ 11: $c = ||r_0||_2 e_1$ perform m steps of GMRES, solving min $||c - H_m y||_2$ for y and generating V_{m+1} and H_m 12: $x_1 = x_0 + V_m y$ 13: $r_1 = V_{m+1}(c - H_m y)$ 14:compute k deflation vectors P_k following def and Section 4 15:16: $Y_k = V_m P_k$ let [Q, R] be the QR factorization of $H_m P_k$ 17: $C_k = V_{m+1}Q$ 18: $U_k = \tilde{Y}_k R^{-1}$ 19: 20: end if 21: while $||r_i||_2 > \varepsilon$ do 22:i = i + 1perform m - k Arnoldi steps with the linear operator $(I - C_k C_k^H)A$, letting $v_1 =$ 23: $|r_{i-1}/||r_{i-1}||_2$ and generating V_{m-k+1}, H_{m-k} , and $B_{m-k} = C_k^H A V_{m-k}$ let D_k be a diagonal scaling matrix such that $\tilde{U}_k = U_k D_k$, where the columns of \tilde{U}_k have 24:a unit norm $\hat{V}_m = [\tilde{U}_k, V_{m-k}], \ \hat{W}_{m+1} = [C_k, V_{m-k+1}] \\ G_m = \begin{pmatrix} D_k & B_{m-k} \\ 0 & H_{m-k} \end{pmatrix}$ 25:26:solve $\min \|G_m y - \hat{W}_{m+1}^H r_{i-1}\|_2$ for y 27: $x_i = x_{i-1} + V_m y$ 28: $r_i = r_{i-1} - \hat{W}_{m+1}G_m y$ 29: set $k = k_{max}$ compute k deflation vectors P_k following def and Section 4 30: $\hat{Y}_k = V_m P_k$ 31:let [Q, R] be the QR factorization of $G_m P_k$ 32: $C_k = \hat{W}_{m+1}Q$ 33: $U_k = \tilde{Y}_k R^-$ 34: 35: end while 36: let $Y_k = U_k$

Operations	cost (flops)
Preparation of the residual, the solution, and the starting vector	n(3k+4)
Arnoldi orthonormalization	$n(m^2 + 4(m-k) - k^2)$
Hessenberg and residual updates	$2m^2 - 2mk + k^2 + m - k$
Solve the least squares problem and expand the solution	$nm + \frac{1}{2}(m^2 + 3m)$
Prepare matrices and solve GEVP (Ritz A)	$n(mk + k^2 + k) + 68m^3 + 2m^2$
Prepare matrices and solve GEVP (Ritz A^{-1})	$nmk + 68m^3 + 2m^2$
Prepare matrices and solve GEVP (Ritz $A^H A$)	$nk^2 + 15m^3 + m^2$
Prepare C_k and \tilde{U}_k	$nk(2m+3) + m^{2}k + m(\frac{5k^{2}+3k}{2}) - \frac{2}{3}k^{3}$

TABLE 2 – Detailed computation cost of one cycle of GMRES-MDR supposing that k deflation vector exist. The maximum dimension of the search subspace is m. GEVP stands for generalized eigenvalue problem. Ritz A, Ritz A^{-1} , and Ritz A^HA refer to the deflation strategy.

Since GMRES-MDR, does not need an explicit coefficient matrix, we use a reverse communication interface (RCI) [2]. Algorithms 5, 6 present the schemes of the RCI loop and the RCI routine, respectively.

Algorithm 5 RCI loop		
1: $ido = 0$		
2: while $ido ! = 99 \operatorname{do}$		
3: call the reverse communication interface routine 6 and get the value of <i>ido</i>		
4: if $ido == 1$ then		
5: apply the preconditioned operator		
6: end if		
7: if $ido == 2$ then		
8: apply the preconditioner		
9: end if		
10: end while		

6.3 Cost analysis

Here, we compare the cost of one cycle of presented variations of GMRES-MDR to the cost of the classical GMRES method. Since the cost of the sparse matrix vector multiplication depends on the sparsity of the matrix A, we do not take it into account in our analysis. In our analysis of the cost of GMRES-MDR, we suppose that the deflation vectors exist i.e., starting from the second cycle. The computation costs consider the real-valued matrices and vectors. The computation costs of the generalized eigenvalue problems can be found in [4]. In Table 2, we detail the computation cost of one cycle of GMRES-MDR(m, k). Table 3 shows a comparison between the computation costs of the variants of GMRES-MDR and the classical GMRES. In Table 4, we detail the communication cost of one cycle between the variants of GMRES-MDR (m, k). Table 5 presents a comparison of communication costs of one cycle between the variants of GMRES-MDR and the classical GMRES. It is clear that the computation cost of the SVD variation is the smallest compared to the two other deflation strategies. This cost can be written as :

$$C(n,m,k) = n(m^2 + (2k+5)m + 2k + 4) + m^2(k + \frac{7}{2}) + m\frac{5k^2 - k + 5}{2} - \frac{2}{3}k^3 + k^2 - k.$$

Inria

Algorithm 6 RCI routine

Require: *ido* the interface indicator, parameters of the solver 1: static variable status = 02: if ido == 0 then set up the solver using the parameters 3: if deflation vectors exist then 4: 5: set ido = 1, status = -1return to apply the preconditioned operator on the deflation vectors 6: 7:else 8: status = 1end if 9: 10: end if 11: if status = -1 then prepare the deflation vectors and matrices 12:set ido = 10, status = 113:14: return 15: end if 16: if status = 1 then prepare for the Arnoldi procedure 17:set ido = 1, status = 2 and return 18: 19: end if 20: if status = 2 then 21: orthogonalization of new basis vector and Hessenberg operations and stopping criterion if convergence and max interior iteration are not achieved then 22:ido = 1, status = 223:24:return to apply the preconditioned operator on the last basis vector else 25:solve the least squares problem and recover the solution 26:compute the deflation subspace 27:if convergence or the max outer iteration is achieved then 28:29:ido = 2, status = 99return to recover the unpreconditioned solution 30: 31: else if max interior iteration is achieved then ido = 1, status = 332: return to apply the preconditioned operator on the solution 33: end if 34:35: end if 36: end if 37: if status = 3 then compute the residual 38: set ido = 10, status = 1 and return 39: 40: end if 41: if status = 99 then set ido = 99, status = 0 and return 42: 43: end if

Method	cost(flops)
$\mathrm{GMRES}(m)$	$C(n,m,0) - 2n - m^2$
GMRES-MDR(m,k) (Ritz A)	$C(n,m,k) + nk(m+1) + m^2$
GMRES-MDR (m, k) (Ritz A^{-1})	$C(n,m,k) + nk(m-k) + m^2$
GMRES-MDR (m, k) (Ritz $A^H A$)	C(n,m,k)

TABLE 3 – Comparison of computation cost of one cycle of GMRES-MDR (supposing that k deflation vector exist) and classical GMRES. The maximum dimension of the search subspace is m. GEVP stands for generalized eigenvalue problem. Ritz A, Ritz A^{-1} , and Ritz A^HA refer to the deflation strategy. $C(n, m, k) = n(m^2 + (2k+5)m + 2k+4) + m^2(k + \frac{7}{2}) + m\frac{5k^2 - k + 5}{2} - \frac{2}{3}k^3 + k^2 - k$.

Operations	messages	words
Preparation of the residual, the solution, and the starting vector	2	(k, 1)
Arnoldi orthonormalization at iteration j	2	(j + k, 1)
Prepare matrices for GEVP (Ritz A)	1	k(m+k+1)
Prepare matrices for GEVP (Ritz A^{-1})	1	k(m+1)
Prepare matrices for GEVP (Ritz $A^H A$)	1	k^2
Prepare $ ilde{U}_k$	1	k

TABLE 4 – Detailed communication costs of one cycle of GMRES-MDR supposing that k deflation vectors exist. The maximum dimension of the search subspace is m. GEVP stands for generalized eigenvalue problem. Ritz A, Ritz A^{-1} , and Ritz A^HA refer to the deflation strategy. When the operation counts 2 messages the number of words is separated into two quantities corresponding to each message. The Arnoldi procedure uses the classical Gram-Schmidt variation.

Method	messages
$\mathrm{GMRES}(m)$	2m + 1
GMRES-MDR(m, k)	2m - 2k + 4

TABLE 5 – Cost analysis of one cycle of GMRES-MDR (supposing that k deflation vectors exist) and classical GMRES. The maximum dimension of the search subspace is m. GEVP stands for generalized eigenvalue problem. Ritz A, Ritz A^{-1} , and Ritz A^HA refer to the deflation strategy. $C(n, m, k) = n(m^2 + (2k + 5)m + 2k + 4) + m^2(k + \frac{7}{2}) + m\frac{5k^2 - k + 5}{2} - \frac{2}{3}k^3 + k^2 - k.$

To give an illustration of the difference between the computation cost of the SVD variant of GMRES-MDR and the classical GMRES, we can set m = 2k. Then the difference is

$$C(n, 2k, k) - C(n, 2k, 0) = n(k^2 + 2k + 2) + (9 - \frac{2}{3})k^3 + 4k^2 - 3k.$$

However, it is sufficient that k > 1 to make the communication cost of GMRES-MDR smaller than the communication cost of the classical GMRES.

7 Numerical experiments

In this section we illustrate the impact of the proposed method on the convergence of sequences of linear systems arising from reservoirs simulations. We present sequential experiments performed in MATLAB as well as parallel experiments. Our set of test cases are obtained from the in-house prototype code at Total, which simulates a complex enhanced oil recovery (EOR) mechanism. This simulator relies on a finite volume discretization and a two-point flux approximation. We have two sequences of linear systems denoted with initials BIGCO24 and BIGP1. BIGCO24 corresponds to a simulation of water and gaz injection using a compositional model (8 hydrocarbon components). The permeability field is heterogeneous. The grid has 83587 active cells. BIGP1 comes from the simulation of water injection using a black-oil model. The permeability field is heterogeneous (sector model from a real field case). The grid has 42332 active cells.

7.1 MATLAB numerical experiments



FIGURE 2 – Convergence history, singular vectors deflated GMRES against eigenvectors deflated GMRES. The convergence history of non-deflated GMRES is also plotted, results for the matrix BIGP1 in MATLAB 2017. The 20 computed approximated vectors (singular vectors and eigenvectors) are approximated up to a tolerance of 10^{-6} .

Figure 2 shows the impact of deflating the left singular vectors related to the smallest singular values. In each variant 20 vectors are computed approximately to a tolerance of 10^{-6} . The



FIGURE 3 – Comparison between three variants of deflation subspaces. SVD, REV, and HEV stand for the deflation methods related to the matrices A^HA , A and A^{-1} , respectively. The maximum dimension of the search subspace is 30. The maximum dimension of the deflation subspace is 5. Fifteen linear systems in the sequence BIGP1 are solved, three matrices each one 5 times, with different right-hand sides (for index *i* the three methods share the same right-hand side). On the left no reduction of the deflation subspace is performed, on the right a reduction of the deflation subspace is performed with a threshold 10^{-3} .

deflation of the eigenvectors corresponds to (4) that we rewrite here,

$$(I + Z(T^{-1} - I)Z^H)Ax = (I + Z(T^{-1} - I)Z^H)b,$$

where Z is a unitary matrix whose columns span the approximated eigenspace. The deflation of the singular vectors follows the Theorem 2. We write the preconditioned linear system

$$(I - U_2 U_2^H)Ax = (I - U_2 U_2^H)b,$$

where $U_2 = AV_2\Sigma_2^{-1}$ stands for the approximated left singular vectors. The solution is recovered as $x = (I - V_2V_2^H)\bar{x} + V_2x_2$, where \bar{x} is the solution obtained by GMRES and $x_2 = \Sigma_2^{-1}U_2^H b$. We remark that the speed of convergence for both deflated variants are approximately the same with a small advantage for the SVD deflation. Furthermore, in the first iterations the SVD deflated GMRES has a sharp convergence curve while the curve corresponding to the eigenvectors deflated GMRES stagnates for a number of iterations. This phenomenon is important for our application. In each Newton iteration during the simulation of a reservoir the linearized system is solved by using the CPR solver. An iteration of the latter solves the linear system associated to the pressure variable. This matrix does not change during the CPR solve, only the right-hand side changes. In practice, a large threshold of convergence for the pressure level is sufficient. Thus, having stagnation at the beginning of the method, even for a small number of iterations, can impact the global number of iterations over the entire simulation. The relative tolerance of convergence in the experiments related to the sequences of linear systems is fixed to 10^{-1} .

As it was noted in the thesis of Parks [11], the choice of the deflation subspace has an impact on the convergence of sequences of linear systems. Figure 3 does not only confirm this note but also shows how it might be possible that in some configurations the deflation subspace becomes a problem and leads to the stagnation of the residual norm. We solve thirty linear systems related to three newton steps for each deflation strategy. Each newton step is solved by ten iterations of a CPR solver. In order to make a fair comparison and avoid the eventual large components of the right-hand sides on special eigenvectors or singular vectors, we generate random righthand sides with seeds corresponding to the index of the linear system, from one to thirty. In the same figure we show the impact of the reduction of the deflation subspace. Figure 4 shows how the reduction strategy keeps the efficiency of the method when the existing deflation subspace



FIGURE 4 – Comparison between three variants of deflation subspaces. SVD, REV, and HEV stand for the deflation methods related to the matrices $A^H A$, A and A^{-1} , respectively. The maximum dimension of the search subspace is 30. The maximum dimension of the deflation subspace is 5. Thirty linear systems in the sequence BIGCO24 are solved, three matrices each one 10 times, with different right-hand sides (for index *i* the three methods share the same right-hand side). On the left no reduction of the deflation subspace is performed, on the right a reduction of the deflation subspace is performed with a threshold 10^{-3} .

is useful. There might be some unnecessary small increase in the number of iterations due to reducing the deflation subspace. It is possible to apply the reduction strategy after one restart cycle of a new system. This also helps measure the efficiency of the deflation subspace based on the relative residual norm at the end of this cycle. Either we keep it or we reduce it. Our numerical experiments lead to two remarks. The first is that deflating the Ritz pairs related to A is efficient when no deflation subspace exists. The second is that deflating the Ritz pairs related to $A^{H}A$ needs more iterations that the one related to A in order to be efficient. To benefit from both subspaces we propose a simple criterion based on the relative residual during the restart cycle to choose adaptively the subspace of deflation. Figures 5 and 6 present numerical experiments on the same two previous sequences. Figures 7 and 8 present the history of the residual norm corresponding to each strategy of deflation. As we noted previously, we notice that the reduction of the residual norm starting from the first iteration makes the SVD variant has more advantage against other strategies. The total iteration numbers in the BIGCO24 sequence corresponding to the deflation related to A, A^{-1} , and the alternating $(A, A^{H}A)$ is 418, 374, and 163, respectively. We also notice that using the adaptive alternate deflation between the Ritz pairs of A and $A^{H}A$ has the benefit of both strategies, quick useful deflation subspace and quick reduction of residual norm. Moreover in this example we see how the residual norm stagnates starting from the third matrix systems with the harmonic Ritz-based deflation subspace. The total iteration numbers in the BIGP1 sequence corresponding to the deflation strategies related to A, and the alternating $(A, A^H A)$ is 1062 and 686, respectively.

7.2 Parallel numerical experiments

In this section we present the numerical experiments of our implementation of the method GMRES-MDR. All experiments are performed on the supercomputer PANGEA at TOTAL. We had access to a maximum of 1024 nodes. Each node has 16 cores. Our set of test matrices consists of the matrices SPE10 and Grid2D-2000x2000. We solve two sequences, each of nine linear systems in which the matrix is fixed. The right-hand sides are different for each linear system in the sequence. All variants have the same right-hand side for the i^{th} linear system, $i = 1, \ldots, 9$. The matrix is SPE10 $\in \mathbb{R}^{n \times n}$ where n is the number of unknowns n = 1094421, the number of the non-zero elements nnz = 7478141. SPE10 corresponds to a simulation of



FIGURE 5 – Comparison between four variants of deflation subspaces. SVD, REV, and HEV stand for the deflation method related to the matrices A^HA , A, and A^{-1} , respectively. SVD-adap stands for an adaptive strategy that alternates the deflation between SVD and REV based on the relative residual norm during the cycle. Fifteen linear systems in the sequence BIGP1 are solved, three matrices each one 5 times, with different right-hand sides (for index *i* the three methods share the same right-hand side).



FIGURE 6 – Comparison between three variants of deflation subspaces. SVD, REV, and HEV stand for the deflation method related to the matrices $A^H A$, A, and A^{-1} , respectively. SVD-adap stands for an adaptive strategy that alternates the deflation between SVD and REV based on the relative residual norm during the cycle. The maximum dimension of the search subspace is 30. The maximum dimension of the deflation subspace is 5. Thirty linear systems in the sequence BIGCO24 are solved, three matrices each one 10 times, with different right-hand sides (for index *i* the three methods share the same right-hand side).



FIGURE 7 – History of residual norm in a sequence of linear systems (a peak corresponds to the beginning of a new linear system). Comparison between three variants of deflation subspaces. SVD, REV, and HEV stand for the deflation method related to the matrices A^HA , A, and A^{-1} , respectively. The SVD variant uses an adaptive strategy that alternates the deflation between SVD and REV deflation methods based on the relative residual norm during the cycle. The maximum dimension of the search subspace is 30. The maximum dimension of the deflation subspace is 5. Thirty linear systems in the sequence BIGCO24 are solved, three matrices each one 10 times, with different right-hand sides (for index *i* the three methods share the same right-hand side).



FIGURE 8 – History of residual norm in a sequence of linear systems (a peak corresponds to the beginning of a new linear system). Comparison between three variants of deflation subspaces. SVD, REV, and HEV stand for the deflation method related to the matrices A^HA , A, and A^{-1} , respectively. The SVD variant uses an adaptive strategy that alternates the deflation between SVD and REV based on the relative residual norm during the cycle. The maximum dimension of the search subspace is 30. The maximum dimension of the deflation subspace is 5. Fifteen linear systems in the sequence BIGP1 are solved, three matrices each one 5 times, with different right-hand sides (for index *i* the three methods share the same right-hand side).



FIGURE 9 – Comparison between runtime of three variants of deflation subspaces. SVD, REV and HEV stand for subspace method related to the matrices $A^H A$, A and A^{-1} , respectively. The maximum dimension of the subspace is 60. The maximum dimension of the deflation subspace is 20. Nine linear systems sharing the same matrix, Grid2D-2000x2000, with different right-hand sides (for index *i* the three methods share the same right-hand side) are solved.

a black oil with a heterogeneous permeability field [1]. The grid has 1094421 active cells. The matrix Grid2D-2000x2000 $\in \mathbb{R}^{n \times n}$, where n is the number of unknowns $n = 4 \times 10^6$, the number of non-zero elements $nnz = 2 \times 10^7$. This matrix corresponds to a discretization of a pressure system in which the permeability field follows a log-normal distribution with standard deviation $\sigma = 3$ and a mean $\mu = loq(500)$. The discretized domain is the unit square with 2000×2000 grid points. In all experiments, a linear system is solved to warm up the machine before solving each sequence. Figure 9 shows the scalability of the three variants discussed in Section 4. The sequence of linear systems is related to the test matrix Grid2D-2000x2000. The number of nodes (processors) varies from 16 up to 1024 with 2 cores per processor. A block Jacobi preconditioner with LU factorization in each block is used. The number of blocks is equal to the number of processors. In Listing 1 and Listing 2, we present two post-analysis of GMRES-MDR with SVD deflation variant that is associated to Figure 9. Figure 10 shows the scalability of the three variants discussed in Section 4. The sequence of linear systems is related to the test matrix SPE10. The number of nodes (processors) varies from 16 up to 256, 1 core per MPI processor. A block Jacobi preconditioner with ILU(0) factorization in each block is used. The number of blocks is equal to the number of processors. We notice that the three variants scale well with a smaller runtime for the SVD variant. We note that the runtime for the three variants of deflation does not scale on 512 nodes in the test case SPE10. See Figure 11 and the discussion below for the explanation. Figure 11 shows the percentage of the average time (over all processors) spent in the operations during the solution of a representative linear system of the sequence by using the SVD deflation strategy. This runtime corresponds to a solution of the system on 256 nodes. We remark that the time spent in the computation of the generalized eigenvalue problem becomes much more important than all other operations. This explains why the scaling of the runtime stopped for this sequence after 256 nodes.

```
Listing 1 – Post analysis of GMRES-MDR with SVD deflation variant. First linear system in the sequence Grid2D-2000x2000.
```

```
Solver information:
Number of processors: 1024
Number of threads: 2
Preconditioner: LU of block Jacobi
Orthogonalization strategy: Double Block Classical Gram-Schmidt
Deflation strategy: SVD values
Max dimension of search subspace: 60
Max dimension of deflation subspace: 20
```

Post analysis:

Number of iterations: 1395 Number of cycles: 36 Residual norm: 9.96438063135672e-02 Relative residual norm: 9.96438063135672e-02 Relative tolerance of convergence: 1.000000000000000e-01

Times over processors	:	MAX	MIN	AVERAGE
Total	:	3.237501	3.237501	3.237501 s
Preconditioning	:	1.577664	1.209435	1.294308 s
SPMV	:	0.447073	0.081644	0.166228 s
SPMV Diag	:	0.031328	0.025887	0.026805 s
SPMV Off Diag	:	0.068317	0.016273	0.044505 s
SPMV Comm	:	0.322190	0.036086	$0.090636 \ s$
DeflationComputation	:	0.280072	0.275017	0.276312 s
DeflationComm	:	0.021925	0.016195	0.020721 s
Orthogonalization	:	1.539758	1.132159	1.445061 s
Orthogonalization Comm	:	1.094884	0.642479	0.984363 s
HessenbergOperations	:	0.002427	0.002085	0.002221 s



FIGURE 10 – Comparison between runtime of three variants of deflation subspaces. SVD, REV and HEV stand for subspace method related to the matrices $A^H A$, A and A^{-1} , respectively. The maximum dimension of the subspace is 30. The maximum dimension of the deflation subspace is 10. Nine linear systems sharing the same matrix, SPE10, with different right-hand sides (for index *i* the three methods share the same right-hand side) are solved.

Listing 2 – Post analysis of GMRES-MDR with SVD deflation variant. Ninth linear system in the sequence Grid2D-2000x2000.

Solver information: Number of processors: 1024 Number of threads: 2 Preconditioner: LU of block Jacobi Orthogonalization strategy: Double Block Classical Gram-Schmidt Deflation strategy: SVD values Max dimension of search subspace: 60 Max dimension of deflation subspace: 20

Post analysis:

Number of iterations: 39 Number of cycles: 1 Residual norm: 9.99391435927701e-02 Relative residual norm: 9.99391435927701e-02Relative tolerance of convergence: 1.00000000000000e-01 AVERAGE Times over processors MAX MIN 0.069219 0.069219 0.069219 Total \mathbf{s} Preconditioning 0.0382460.0340660.036175 \mathbf{S} SPMV 0.0065810.0021300.004383 \mathbf{s} SPMV Diag 0.0010150.0007340.000759 \mathbf{S} SPMV Off Diag 0.0019260.000461 0.001252 \mathbf{s} SPMV Comm 0.0043820.000921 0.002392 \mathbf{S} DeflationComputation 0.0035000.0032770.003379 \mathbf{S} DeflationComm 0.000360 0.0001730.000268 \mathbf{S} Orthogonalization 0.0256000.0207880.023068 s Orthogonalization Comm 0.0139470.0088240.011384 \mathbf{S} HessenbergOperations0.0000740.000053 $0.000063 \ s$:



FIGURE 11 – Average time (over processors) of different operations in GMRES-MDR as a percentage of the total runtime. The communication and the computation part of the deflation and the orthogonalization are separated. This runtime corresponds to the solution of a representative linear system during the sequence. Number of nodes = 256. Deflation strategy is the SVD.

8 Conclusion

In this chapter we reviewed the deflation concept in Krylov iterative methods. Usual choice of deflation subspaces is related to the approximation of eigenvectors. Based on the study in [15] we introduced the deflation of singular vectors. Since we are interested in solving a sequence of linear systems, we reviewed the method GCRO-DR [10] with an abstract deflation subspace. Two previously introduced subspaces are briefly presented. Afterwards, we proposed a deflation subspace related to an approximation of singular vectors. To validate the proposed strategy, a comparison between the three variants was made on different sequences arising from reservoirs simulations. This validation was performed on a sequential code as well as on a parallel code. The numerical experiments demonstrated the impact of choosing the SVD variant. The gain in the global iteration count is up to 40%. Concerning the parallel implementation, the SVD variant gave better performance than other variants on a short sequence of linear systems. This improvement would be of larger impact when the length of the sequence of linear systems is longer.

Références

- M. A. Christie and M. J. Blunt. Tenth spe comparative solution project : A comparison of upscaling techniques. 2001.
- [2] Jack Dongarra, Victor Eijkhout, and Ajay Kalhan. Reverse communication interface for linear algebra templates for iterative methods, 1995.
- [3] Jocelyne Erhel, Kevin Burrage, and Bert Pohl. Restarted GMRES preconditioned by deflation. Journal of Computational and Applied Mathematics, 69(2):303 – 318, 1996.
- [4] Gene H. Golub and Charles F. Van Loan. Matrix Computations (3rd Ed.). Johns Hopkins University Press, Baltimore, MD, USA, 1996.
- [5] M. R. Hestenes and E. Stiefel. Methods of conjugate gradients for solving linear systems. Journal of research of the National Bureau of Standards., 49 :409–436, 1952.
- [6] Intel(R). Math kernel library.
- [7] Ronald B. Morgan. A restarted GMRES method augmented with eigenvectors. SIAM Journal on Matrix Analysis and Applications, 16(4):1154–1171, 1995.
- [8] Ronald B. Morgan. GMRES with deflated restarting. SIAM Journal on Scientific Computing, 24(1) :20–37, 2002.
- [9] Dianne P. O'Leary. The block conjugate gradient algorithm and related methods. *Linear Algebra and its Applications*, 29:293 322, 1980.
- [10] Michael L. Parks, Eric de Sturler, Greg Mackey, Duane D. Johnson, and Spandan Maiti. Recycling krylov subspaces for sequences of linear systems. SIAM Journal on Scientific Computing, 28(5) :1651–1674, 2006.
- [11] Michael Lawrence Parks. The iterative solution of a sequence of linear systems arising from nonlinear finite element analysis, 2005.
- [12] Y. Saad. Iterative Methods for Sparse Linear Systems. Society for Industrial and Applied Mathematics, Philadelphia, PA, USA, 2nd edition, 2003.
- [13] Youcef Saad and Martin H. Schultz. GMRES : A generalized minimal residual algorithm for solving nonsymmetric linear systems. SIAM Journal on Scientific and Statistical Computing, 7(3):856–869, 1986.

- [14] H. Cao SPE Schlumberger, H.A. Tchelepi SPE Stanford U., J. Wallis Wallis Consulting Inc., and H. Yardumian SPE Chevron. Parallel scalable unstructured CPR-Type Linear Solver for Resevoir Simulation. SPE 96809, 2005.
- [15] V. Simoncini. On the convergence of restarted krylov subspace methods. SIAM Journal on Matrix Analysis and Applications, 22(2):430–452, 2000.
- [16] Marc Snir, Steve Otto, Steven Huss-Lederman, David Walker, and Jack Dongarra. MPI-The Complete Reference, Volume 1 : The MPI Core. MIT Press, Cambridge, MA, USA, 2nd. (revised) edition, 1998.
- [17] J. M. Tang, R. Nabben, C. Vuik, and Y. A. Erlangga. Comparison of two-level preconditioners derived from deflation, domain decomposition and multigrid methods. *Journal of Scientific Computing*, 39(3):340–370, 2009.
- [18] J.R. Wallis, R.P. Kendall, and T.E. Little. Constrained residual acceleration of conjugate residual methods. SPE 13563, 1985.



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