Accelerating ordered subsets image reconstruction for X-ray CT using spatially non-uniform optimization transfer

Donghwan Kim, Student Member, IEEE, Debashish Pal, Member, IEEE, Jean-Baptiste Thibault, Member, IEEE, and Jeffrey A. Fessler, Fellow, IEEE

Abstract—Statistical image reconstruction algorithms in X-ray CT provide improved image quality for reduced dose levels but require substantial computation time. Iterative algorithms that converge in few iterations and that are amenable to massive parallelization are favorable in multiprocessor implementations. The separable quadratic surrogate (SQS) algorithm is desirable as it is simple and updates all voxels simultaneously. However, the standard SQS algorithm requires many iterations to converge. This paper proposes an extension of the SQS algorithm that leads to spatially non-uniform updates. The non-uniform (NU) SQS encourages larger step sizes for the voxels that are expected to change more between the current and the final image, accelerating convergence, while the derivation of NU-SQS guarantees monotonic descent. Ordered subsets (OS) algorithms can also accelerate SQS, provided suitable “subset balance” conditions hold. These conditions can fail in 3D helical cone-beam CT due to incomplete sampling outside the axial region-of-interest (ROI). This paper proposes a modified OS algorithm that is more stable outside the ROI in helical CT. We use CT scans to demonstrate that the proposed NU-OS-SQS algorithm handles the helical geometry better than the conventional OS methods and “converges” in less than half the time of ordinary OS-SQS.

Index Terms—Statistical image reconstruction, computed tomography, parallelizable iterative algorithms, ordered subsets, separable quadratic surrogates

I. INTRODUCTION

STATISTICAL image reconstruction methods can improve resolution and reduce noise and artifacts by minimizing either penalized likelihood (PL) [1]–[3] or penalized weighted least-squares (PWLS) [4]–[6] cost functions that model the physics and statistics in X-ray CT. The primary drawback of these methods is their computationally expensive iterative algorithms. This paper describes new accelerated minimization algorithms for X-ray CT statistical image reconstruction.

There are several iterative algorithms for X-ray CT. Coordinate descent (CD) algorithms [7] (also known as Gauss Siedel algorithms [8, p. 507]) and block/group coordinate descent (BCD/GCD) algorithms [9]–[11], update one or a group of voxels sequentially. These can converge in few iterations but can require long computation time per iteration [6], [12]. Considering modern computing architectures, algorithms that update all voxels simultaneously and that are amenable to parallelization are desirable, such as ordered subsets based on separable quadratic surrogate (OS-SQS) [13]–[15] and preconditioned conjugate gradient (PCG) algorithms [16]. However, those highly parallelizable algorithms require more iterations than CD algorithms [6], [12], and thus it is desirable to reduce the number of iterations needed to reach acceptable images. Splitting techniques [17] can accelerate convergence [18], but require substantial extra memory.

In this paper, we propose an enhanced version of a highly parallelizable SQS algorithm that accelerates convergence. SQS algorithms are optimization transfer methods that replace the original cost function by a simple surrogate function [19], [20]. Here we construct surrogates with spatially non-uniform curvatures that provide spatially non-uniform step sizes to accelerate convergence.

Spatially non-homogeneous (NH) approach [7] accelerated the CD algorithm by more frequently visiting the voxels that need updates. This approach is effective because the differences between the initial and final images are non-uniform. Inspired by such ideas, we propose a spatially non-uniform (NU) optimization transfer method that encourages larger updates for voxels that are predicted to be farther from the optimal value, using De Pierro’s idea in SQS [21]. We provide a theoretical justification for the acceleration of NU method by analyzing the convergence rate of SQS algorithm (in Section II-D). The NH approach also balanced homogeneous and non-homogeneous updates for fast overall convergence rate [7]. Section III-C discusses similar considerations for the proposed NU approach.

Ordered subsets (OS), also known as incremental gradient methods [22], [23] or block iterative methods [24], can accelerate gradient-based algorithms by grouping the projection data into (ordered) subsets and updating the image using each subset. OS algorithms are most effective when a properly scaled gradient of each subset data-fit term approximates the gradient of the full data-fidelity term, and then it can accelerate convergence by a factor of the number of subsets.
II. STATISTICAL IMAGE RECONSTRUCTION

A. Problem

We reconstruct a non-negative image $x = (x_1, \ldots, x_{N_N}) \in \mathbb{R}^{N_N}$ from noisy measured transmission data $Y \in \mathbb{R}^{N_d}$ by minimizing either penalized likelihood (PL) or penalized weighted least-squares (PWLS) cost functions:

$$\hat{x} = \arg\min_{x \geq 0} \Psi(x),$$ (1)

$$\Psi(x) \triangleq L(x) + R(x) = \sum_{i=1}^{N_N} h_i([Ax]_i) + \sum_{k=1}^{N_d} \psi_k([Cx]_k),$$ (2)

where $\hat{x}$ is a minimizer of $\Psi(x)$ subject to a non-negativity constraint. The function $L(x)$ is a negative log-likelihood term (data-fit term) and $R(x)$ is a regularizer. The matrix $A = \{a_{ij}\}$ is a projection operator ($a_{ij} \geq 0$ for all $i$, $j$) where $[Ax]_i \triangleq \sum_{j=1}^{N_p} a_{ij}x_j$, and $C = \{c_{k}\}$ is a finite differencing matrix considering 26 neighboring voxels in 3D image space. The function $\psi_k(t)$ is a (convex and typically non-quadratic) edge-preserving potential function. The function $h_i(t)$ is selected based on the chosen statistics and physics:

- Penalized likelihood (PL) for pre-log data $Y_i$ with Poisson model $[1]$–$[3]$ uses:

  $$h_i(t) = (b_i e^{-t} + r_i) - Y_i \log(b_i e^{-t} + r_i),$$ (3)

  where $b_i$ is the blank scan factor and $r_i$ is the mean number of background events. The function $h_i(\cdot)$ is non-convex if $r_i \neq 0$, or convex otherwise. A shifted Poisson model $[30]$ that partially accounts for electronic recorded noise can be used instead.

- Penalized weighted least squares (PWLS) for post-log data $y_i = \log(b_i/(Y_i - r_i))$ with Gaussian model $[4]$–$[6]$ uses a convex quadratic function:

  $$h_i(t) = \frac{1}{2} w_i (t - y_i)^2,$$ (4)

  where $w_i = (Y_i - r_i)^2/Y_i$ provides statistical weighting. We use the PWLS cost function for our experiments in Section IV and V.

The proposed NU-SQS algorithm, based on optimization transfer methods (in Section II-B), decreases the cost function $\Psi(x)$ monotonically for either (3) or (4).

B. Optimization transfer method

When a cost function $\Psi(x)$ is difficult to minimize, we replace $\Psi(x)$ with a surrogate function $\phi^{(n)}(x)$ at the $n$th iteration for computational efficiency. This method is called optimization transfer $[19]$, $[20]$, which is also known as a majorization principle $[31]$, and a comparison function $[32]$.

There are many optimization transfer algorithms such as expectation maximization (EM) algorithms $[33]$, $[34]$, separable

Each row of $C$ consists of a permutation of $(1, -1, 0, \ldots, 0) \in \mathbb{R}^{N_p}$ where the indices of the nonzero entries 1 and $-1$ correspond to adjacent voxel locations in 3D image space.
surrogate algorithms based on De Pierro’s lemma [35]–[37] and surrogate algorithms using Lipschitz constants [38], [39].

The basic iteration of an optimization transfer method is

$$x_{(n+1)} = \arg \min_{x \geq 0} \phi^{(n)}(x).$$

To monotonically decrease $\Psi(x)$, we design surrogate functions $\phi^{(n)}(x)$ that satisfy the following majorization conditions:

$$\Psi(x^{(n)}) = \phi^{(n)}(x^{(n)}),$$

$$\Psi(x) \leq \phi^{(n)}(x), \quad \forall x \in \mathbb{R}^{N_p}_+.$$  

Constructing surrogates with smaller curvatures while satisfying condition (6) is the key to faster convergence in optimization transfer methods [11].

Optimization transfer has been used widely in tomography problems. De Pierro developed a separable surrogate (SS) approach in emission tomography [35], [36]. Quadratic surrogate (QS) functions have been derived for non-quadratic problems, enabling monotonic descent [1]. QS algorithms combine SS and QS [14], and are the focus of this paper. Partitioned QS methods for multi-core processors have been proposed for separating the image domain by the number of processors and updating each of them separately while preserving the monotonicity [40]. In addition, replacing $\mathbb{R}^{N_p}_+$ in (6) by an interval that is known to include the minimizer $\hat{x}$ can reduce the surrogate curvature [7], [41].

C. Separable quadratic surrogate (SQS) algorithm

We first construct a quadratic surrogate at the nth iteration for the non-quadratic cost function in (2):

$$\Psi(x) = L(x) + R(x) \leq Q_{L}^{(n)}(x) + Q_{R}^{(n)}(x),$$   

where $Q_{L}^{(n)}(x)$ and $Q_{R}^{(n)}(x)$ are quadratic surrogates for $L(x)$ and $R(x)$. Based on (2), the quadratic surrogate for $L(x)$ has the form:

$$Q_{L}^{(n)}(x) = \sum_{i=1}^{N_d} q_{i}^{(n)}(\{Ax^{(n)}\}_i),$$

$$q_{i}^{(n)}(t) = h_{i}(t_{i}^{(n)}) + \hat{h}_{i}(t_{i}^{(n)})(t - t_{i}^{(n)}) + \frac{\hat{c}_{i}^{(n)}}{2}(t - t_{i}^{(n)})^2,$$

where $t_{i}^{(n)} \triangleq [Ax^{(n)}]_i$, and $\hat{c}_{i}^{(n)} = \max \left\{ \hat{c}_{i}, \eta \right\}$ is the curvature of $q_{i}^{(n)}(t)$ for some small positive value $\eta$ that ensures the curvature $\hat{c}_{i}^{(n)}$ positive [1]. In PWLS problem, $h_{i}(.)$ is quadratic already, so $q_{i}^{(n)}(t) = h_{i}(t)$. The quadratic surrogate $Q_{R}^{(n)}(x)$ for $R(x)$ is defined similarly.

We choose curvatures $\{\hat{c}_{i}^{(n)}\}$ that satisfy the monotonicity conditions in (6). For PL, the smallest curvatures:

$$\hat{c}_{i}^{(n)} = \begin{cases} \frac{2h_{i}(0) - h_{i}(t_{i}^{(n)}) + t_{i}^{(n)}h_{i}(t_{i}^{(n)})}{\hat{h}_{i}(0)^2_{+}}, & t_{i}^{(n)} > 0, \\ \hat{h}_{i}(0)_{+}, & t_{i}^{(n)} = 0, \end{cases},$$

where $[t]_{+} = \max\{t, 0\}$, called “optimal curvatures,” lead to the fastest convergence rate but require an extra back-projection each iteration for non-quadratic problems [1]. Alternatively, we may use “maximum curvatures”:

$$\hat{c}_{i}^{(n)} = \max_{t \geq 0} \hat{h}_{i}(t)$$

that we can precompute before the first iteration [1].

Next, we generate a separable surrogate of the quadratic surrogate. For completeness, we repeat De Pierro’s argument in [14]. We first rewrite forward projection $[Ax]_i$ as follows:

$$[Ax]_i = \sum_{j=1}^{N_p} a_{ij}x_j = \sum_{j=1}^{N_p} \pi_{ij}(n)\left(\frac{a_{ij}}{\pi_{ij}(n)}(x_j - x_j^{(n)}) + [Ax^{(n)}]_i\right),$$

where a non-negative real number $\pi_{ij}(n)$ is zero only if $a_{ij}$ is zero for all $i$, $j$, and satisfies $\sum_{j=1}^{N_p} \pi_{ij}(n) = 1$ for all $i$. Using the convexity of $q_{i}^{(n)}(\cdot)$ and the convexity inequality yields

$$q_{i}^{(n)}([Ax]_i) \leq \sum_{j=1}^{N_p} \pi_{ij}(n) q_{i}^{(n)}\left(\frac{a_{ij}}{\pi_{ij}(n)}(x_j - x_j^{(n)}) + [Ax^{(n)}]_i\right).$$

Thus we have the following separable quadratic surrogate $\phi_{L}^{(n)}(x)$ (with a diagonal Hessian) for the data-fit term $L(x)$:

$$L(x) \leq Q_{L}^{(n)}(x) \leq \phi_{L}^{(n)}(x) \leq \sum_{j=1}^{N_d} \phi_{L,j}^{(n)}(x_j),$$

$$\phi_{L,j}^{(n)}(x_j) \triangleq \sum_{i=1}^{N_d} \pi_{ij}(n) q_{i}^{(n)}\left(\frac{a_{ij}}{\pi_{ij}(n)}(x_j - x_j^{(n)}) + [Ax^{(n)}]_i\right).$$

The second derivative (curvature) of the surrogate $\phi_{L,j}^{(n)}(x_j)$ is

$$d_{j}^{L,(n)} \triangleq \frac{\partial^2}{\partial x_j^2} \phi_{L,j}^{(n)}(x_j) = \sum_{i=1}^{N_d} \pi_{ij}(n) \frac{\hat{c}_{i}^{(n)}}{\pi_{ij}} C_{ij},$$

where $\pi_{kj}^{(n)}(n)$ have similar constraints as $\pi_{ij}^{(n)}$. $\psi_k(0)$ is maximum curvature [14], or $\psi_k(0)$ can be replaced by $\psi_k([CAx^{(n)}]^k) / [CAx^{(n)}]^k$ for Huber’s optimal curvature [32, Lemma 8.3, p.184].

Combining the surrogates for the data-fit term and regularizer and minimizing it in (5) leads to the following separable quadratic surrogate (SQS) method [14] that updates all voxels simultaneously with a “denominator” $d_{j}^{L,(n)} + d_{j}^{R,(n)}$:

$$x_{j}^{(n+1)} = \left[ x_j^{(n)} - \frac{1}{d_{j}^{L,(n)} + d_{j}^{R,(n)}} \partial_{x_j} \Psi(x^{(n)}) \right]_+, $$

where $\Psi(x)$ is an objective function.
where a clipping $[·]_+$ enforces the non-negativity constraint. This SQS decreases the cost function $\Psi(x)$ monotonically, and it converges based on the proof in [20]. If $\Psi(x)$ is convex, a sequence $\{x^{(n)}\}$ converges to $x^{(\infty)}$ that is a global minimizer $\hat{x}$. Otherwise, $\{x^{(n)}\}$ converges to a local minimizer $x^{(\infty)}$ which may or may not be a global minimizer $\hat{x}$ depending on the initial image $x^{(0)}$.

The implementation and convergence rate of SQS depend on the choice of $\pi_{ij}^{(n)}$. A general form for $\pi_{ij}^{(n)}$ is

$$\pi_{ij}^{(n)} = \frac{\lambda_{ij}^{(n)}}{\sum_{l=1}^{N_p} \lambda_{il}^{(n)}},$$

(18)

where a non-negative real number $\lambda_{ij}^{(n)}$ is zero only if $a_{ij}$ is zero. Then (15) can be re-written as

$$d_j^{L(n)} = \sum_{i=1}^{N_d} \frac{\lambda_{ij}^{(n)}}{\sum_{l=1}^{N_p} \lambda_{il}^{(n)}} a_{ij} \left(\sum_{l=1}^{N_p} a_{il}\right),$$

(19)

Summations involving the constraint $a_{ij} \neq 0$ require knowledge of the projection geometry, and thereby each summation can be viewed as a type of forward or back projection.

The standard choice [11], [14]:

$$\lambda_{ij} = a_{ij}, \quad \bar{\lambda}_{ij} = |c_{ij}|,$$

(20)

leads to

$$d_j^{L(n)} = \sum_{i=1}^{N_d} \bar{\lambda}_{ij} a_{ij} \left(\sum_{l=1}^{N_p} a_{il}\right),$$

(21)

and

$$d_j^{R(n)} = \sum_{k=1}^{N_r} \bar{\lambda}_{ik} c_{ij} \left(\sum_{l=1}^{N_p} c_{kl}\right).$$

(22)

This choice is simple to implement, since the (available) standard forward and back projections can be used directly in (21). (Computing $d_j^{R(n)}$ in (22) is negligible compared with (21).) The standard SQS generates a sequence $\{x^{(n)}\}$ in (17) by defining the denominator as

$$d_j^{(n)} \triangleq d_j^{L(n)} + d_j^{R(n)}.$$

(23)

However, we prefer choices for $\lambda_{ij}^{(n)}$ (and $\lambda_{ij}^{(n)}$) that provide fast convergence. Therefore, we first analyze the convergence rate of the SQS algorithm in terms of the choice of $\lambda_{ij}^{(n)}$ in the next section. Section III introduces acceleration by choosing better $\lambda_{ij}^{(n)}$ (and $\lambda_{ij}^{(n)}$) than the standard choice (20).

D. Convergence rate of SQS algorithm

The convergence rate of the sequence $\{x^{(n)}\}$ generated by the SQS iteration (17) depends on the denominator $D^{(n)} \triangleq \text{diag}\{d_j^{(n)}\}$. This paper’s main goal is to choose $\lambda_{ij}^{(n)}$ so that the sequence $\{x^{(n)}\}$ converges faster.

The asymptotic convergence rate of a sequence $\{x^{(n)}\}$ that converges to $x^{(\infty)}$ is measured by the root-convergence factor defined as $R_1 \{x^{(n)}\} \triangleq \limsup_{n \to \infty} \left|\frac{x^{(n)} - x^{(\infty)}}{x^{(n-1)} - x^{(\infty)}}\right|^{1/n}$ in [31, p. 288]. The root-convergence factor at $x^{(\infty)}$ for SQS algorithm is given as $R_1 \{x^{(n)}\} = \rho (I - [D^{(n)-1}] H^{(\infty)})$ in [31, Linear Convergence Theorem, p. 301] and [42, Theorem 1], where the spectral radius $\rho(\cdot)$ of a square matrix is its largest absolute eigenvalue and $H^{(\infty)} \triangleq \nabla^2 \Psi(x^{(\infty)})$, assuming that $D^{(n)}$ converges to $D^{(\infty)}$. Faster convergence, we want $R_1 \{x^{(n)}\}$ and $\rho(\cdot)$ to be smaller. We can reduce the root-convergence factor based on$^2$ [42, Lemma 1], by using a smaller denominator $D^{(n)}$ subject to the majorization conditions in (6) and (13).

However, the asymptotic convergence rate does not help us design $D^{(n)}$ in the early iterations, so we consider another factor that relates to the convergence rate of SQS:

**Lemma I**: For a fixed denominator $D$ (using the maximum curvature (10)), a sequence $\{x^{(n)}\}$ generated by an SQS algorithm (17) satisfies

$$\Psi(x^{(n+1)}) - \Psi(x^{(\infty)}) \leq \frac{\left|x^{(0)} - x^{(\infty)}\right|^2_D}{2(n+1)},$$

(24)

for any $n \geq 0$, if $\Psi(x)$ is convex. Lemma I is a simple generalization of Theorem 3.1 in [39], which was shown for a surrogate with a scaled identity Hessian (using Lipschitz constant). The inequality (24) shows that minimizing $\left|x^{(0)} - x^{(\infty)}\right|^2_D$ with respect to $D$ will reduce the upper bound of $\Psi(x^{(n)}) - \Psi(x^{(\infty)})$, and thus accelerate convergence. (Since the upper bound is not tight, there should be a room for further acceleration by choosing better $D$, but we leave it as future work.)

We want to adaptively design $D^{(n)}$ to accelerate convergence at the nth iteration. We can easily extend Lemma 1 to Corollary 1 by treating the current estimate $x^{(n)}$ as an initial image for the next SQS iteration:

**Corollary I**: A sequence $\{x^{(n)}\}$ generated by an SQS algorithm (17) satisfies

$$\Psi(x^{(n+1)}) - \Psi(x^{(\infty)}) \leq \frac{\left|x^{(n)} - x^{(\infty)}\right|^2_{D^{(n)}}}{2}$$

(25)

for any $n \geq 0$, if $\Psi(x)$ is convex. The inequality (25) motivates us to use $\left|x^{(n)} - x^{(\infty)}\right|^2_{D^{(n)}}$ when selecting $d_j^{(n)}$ (and $\lambda_{ij}^{(n)}$) to accelerate convergence at nth iteration. We discuss this further in Section III-A. We fix $D^{(n)}$ after the $n_{\text{iter}}$ number of iterations to ensure convergence of SQS iteration (17), based on [20]. In this case, $D^{(n)}$ must be generated by the maximum curvature (10) to guarantee the majorization condition (6) for subsequent iterations.

From (17) and (19), the step size $\Delta_j^{(n)}$ of the SQS iteration (17) has this relationship:

$$\Delta_j^{(n)} \triangleq \frac{x^{(n+1)} - x^{(n)}}{d_j^{(n)}} \propto \frac{1}{d_j^{(n)}},$$

(26)

where smaller $d_j^{(n)}$ (and relatively larger $\lambda_{ij}^{(n)}$) values lead to larger steps. Therefore, we should encourage $d_j^{(n)}$ to be small ($\lambda_{ij}^{(n)}$ to be relatively large) to accelerate the SQS algorithm.

$^2$If $D^{-1} \geq D_1^{-1} \geq H^{(\infty)} \geq 0$, then $\rho(I - D_1^{-1}H^{(\infty)}) \leq \rho(I - D_1^{-1}H^{(\infty)}) < 1$. 


However, we cannot reduce \( d_j^{(n)} \) simultaneously for all voxels, due to the majorization conditions in (6) and (13). Lemma 1 (and Corollary 1) suggest intuitively that we should try to encourage larger steps \( \Delta_j^{(n)} \) (smaller \( d_j^{(n)} \)) for the voxels that are far from the optimum to accelerate convergence.

III. SPATIALLY NON-UNIFORM SEPARABLE QUADRATIC SURROGATE (NU-SQS)

We design surrogates that satisfy condition (6) and provide faster convergence based on Section II-D. We introduce the “update-needed factors” and propose a spatially non-uniform SQS (NU-SQS) algorithm.

A. Update-needed factors

Based on Corollary 1, knowing \( |x_j^{(n)} - x_j^{(\infty)}| \) would be helpful for accelerating convergence at the \( n \)th iteration, but \( x_j^{(\infty)} \) is unavailable in practice. NH-CD algorithm [7] used the difference between the current and previous iteration instead:

\[
\hat{u}_j^{(n)} \triangleq \max \left\{ |x_j^{(n)} - x_j^{(n-1)}|, \delta_j^{(n)} \right\},
\]

(27)

which we call the “update-needed factors” (originally named a voxel selection criterion (VSC) in [7]). Including the small positive values \( \{\delta_j^{(n)}\} \) ensures all voxels to have at least a small amount of attention for updates. This \( u_j^{(n)} \) accelerated the NH-CD algorithm by visiting voxels with large \( u_j^{(n)} \) more frequently.

B. Design

For SQS, we propose to choose \( \lambda_{ij}^{(n)} \) to be larger if the \( j \)th voxel is predicted to need more updates based on the “update-needed factors” (27) after the \( n \)th iteration. We select

\[
\hat{\lambda}_{ij}^{(n)} = a_{ij} u_j^{(n)},
\]

(28)

which is proportional to \( u_j^{(n)} \) and satisfies the condition for \( \lambda_{ij} \). This choice leads to the following NU-based denominator:

\[
d_j^{L,(n)} = \frac{1}{u_j^{(n)}} \sum_{i=1}^{N_d} c_i^{(n)} a_{ij} \left( \sum_{l=1}^{N_p} a_{il} u_l^{(n)} \right),
\]

(29)

which leads to spatially non-uniform updates \( \Delta_j^{(n)} \propto u_j^{(n)}. \)

If it happened that

\[
|x_j^{(n)} - x_j^{(\infty)}| \approx B|x_j^{(n)} - x_j^{(n-1)}| \text{ for all } j,
\]

(30)

where \( B \) is a constant, then the NU denominator \( d_j^{L,(n)} \) would minimize the upper bound of \( \Psi(x(n+1)) - \Psi(x(\infty)) \) in Corollary 1:

**Lemma 2:** The proposed choice \( d_j^{L,(n)} \) in (29) minimizes the following weighted sum of the denominators

\[
\sum_{j=1}^{N_p} (u_j^{(n)})^2 d_j^{L,(n)}
\]

(31)

over all possible choices of the \( d_j^{L,(n)} \) in (19).

**Proof:** In Appendix A.

The proposed \( d_j^{L,(n)} \) in (29) reduces to the standard choice \( d_j^{(n)} \) in (21) when \( \{u_j^{(n)}\} \) is uniform. Similar to the standard choice \( d_j^{(n)} \), the proposed choice \( d_j^{L,(n)} \) can be implemented easily using standard forward and back projection. However, since \( d_j^{L,(n)} \) depends on iteration \( n \), additional projections required for \( d_j^{L,(n)} \) at every iteration would increase computation. We discuss ways to reduce this burden in Section III-F.

Similar to the data-fit term, we derive the denominator of NU-SQS for the regularizer term to be:

\[
d_j^{R,(n)} = \frac{1}{u_j^{(n)}} \sum_{k=1}^{N_p} \psi_k(0) |c_{kj}| \left( \sum_{l=1}^{N_p} |c_{kl}| u_l^{(n)} \right),
\]

(32)

from the choice \( \chi_k = |c_{kj}| u_j^{(n)} \) and the maximum curvature method in [14]. Alternatively, we may use Huber’s optimal curvature [32, Lemma 8.3, p.184] replacing \( \psi_k(0) \) in (32) by \( \psi_k([C x^{(\infty)}_k]/[C x^{(n)}_k]). \) The computation of (32) is much less than that of the data-fit term.

Defining the denominator in the SQS iteration (17) as

\[
d_j^{(n)} \triangleq d_j^{L,(n)} + d_j^{R,(n)}
\]

(33)

leads to the accelerated NU-SQS iteration, while the algorithm monotonically decreases \( \Psi(x) \) and is provably convergent [20]. We can further accelerate NU-SQS by ordered subsets (OS) methods [13, 14], while losing the guarantee of monotonicity. This algorithm, called ordered subsets algorithms based on a spatially non-uniform SQS (NU-OS-SQS), is explained in Section IV.

C. Dynamic range adjustment (DRA) of \( u_j^{(n)} \)

In reality, (30) will not hold, so (27) will be suboptimal. We could try to improve (27) by finding a function \( f^{(n)}(\cdot) : [\delta^{(n)}, \infty) \to [\epsilon, 1] \) based on the following:

\[
\arg \min_{f^{(n)}(\cdot)} \sum_{j=1}^{N_p} \left( f^{(n)}(u_j^{(n)}) - \frac{|x_j^{(n)} - x_j^{(\infty)}|}{\max_j |x_j^{(n)} - x_j^{(\infty)}|} \right)^2,
\]

(34)

where \( \epsilon \) is a small positive value. Then we could use \( f^{(n)}(u_j^{(n)}) \) as (better) update-needed factors. However, solving (34) is intractable, so we searched empirically for good candidates for a function \( f^{(n)}(\cdot) \).

Intuitively, if the dynamic range of the update-needed factors \( u_j^{(n)} \) in (27) is too wide, then there will be too much focus on the voxels with relatively large \( u_j^{(n)} \), slowing the overall convergence rate. On the other hand, a narrow dynamic range of \( u_j^{(n)} \) will provide no speed-up, since the algorithm will distribute its efforts uniformly. Therefore, adjusting the dynamic range of the update-needed factors is important to achieve fast convergence. This intuition corresponds to how the NH-CD approach balanced between homogeneous update orders and non-homogeneous update orders [7].

To adjust the dynamic range and distribution of \( u_j^{(n)} \), we first construct their empirical cumulative density function:

\[
F_{cdl}^{(n)}(u) \triangleq \frac{1}{N_p} \sum_{j=1}^{N_p} I\{u_j^{(n)} \leq u\}
\]

(35)
to somewhat normalize their distribution, where \( I_B = 1 \) if \( B \) is true or 0 otherwise. Then we map the values of \( F^{(n)}_{\text{eff}}(u) \) by a non-decreasing function \( g(\cdot) : [0, 1] \to [\epsilon, 1] \) as follows

\[
\tilde{u}^{(n)}_j = f^{(n)}(u^{(n)}_j) = g \left( F^{(n)}_{\text{eff}}(u^{(n)}_j) \right),
\]

which controls the dynamic range and distribution of \( \tilde{u}^{(n)}_j \), and we enforce positivity in \( g(\cdot) \) to ensure that the new adjusted parameter \( \tilde{\lambda}^{(n)}_{ij} = a_{ij} \tilde{u}^{(n)}_j \) is positive if \( a_{ij} \) is positive. (We set \( \delta^{(n)}_j \) in (27) to zero here, since a positive parameter \( \epsilon \) ensures the positivity of \( \tilde{\lambda}^{(n)}_{ij} \) if \( a_{ij} \) is positive.)

The transformation (36) from \( u^{(n)}_j \) to \( \tilde{u}^{(n)}_j \) is called dynamic range adjustment (DRA), and two examples of such \( \tilde{u}^{(n)}_j \) are presented in Fig. 2. Then we use \( \tilde{u}^{(n)}_j \) instead of \( u^{(n)}_j \) in (28).

Here, we focus on the following function for adjusting the dynamic range and distribution:

\[
g(v) \triangleq \max \left\{ v^t, \epsilon \right\}
\]

where \( t \) is a non-negative real number that controls the distribution of \( \tilde{u}^{(n)}_j \) and \( \epsilon \) is a small positive value that controls the maximum dynamic range of \( \tilde{u}^{(n)}_j \). The function reduces to the ordinary SQS choice in (20) when \( t = 0 \). The choice of \( g(\cdot) \), particularly the parameters \( t \) and \( \epsilon \) here, may influence the convergence rate of NU-SQS for different data sets, but we show that certain values for \( t \) and \( \epsilon \) consistently provide fast convergence for various data sets.

Fig. 2. A shoulder region scan: \( \tilde{u}^{(2)}_j \) and \( \tilde{u}^{(8)}_j \) after dynamic range adjustment (DRA) for NU-OS-SQS(8/2 subsets), with the choice \( g(v) = \max \{ v^t, 0.05 \} \). NU-OS-SQS updates the voxels with large \( u^{(n)}_j \), whereas ordinary OS-SQS updates all voxels equivalently.

**D. Related work**

In addition to the standard choice (20), the choice

\[
\lambda^{(n)}_{ij} = a_{ij} \max \left\{ x^{(n)}_j, \delta \right\},
\]

with a small non-negative \( \delta \), has been used in emission tomography problems [35], [36] and in transmission tomography problems [11], [37]. This choice is proportional to \( x^{(n)}_j \) and thereby provides a relationship \( \Delta^{(n)}_j \propto x^{(n)}_j \). This classical choice (38) can be also viewed as another NU-SQS algorithm based on “intensity”. However, intensity is not a good predictor of which voxels need more update, so (38) does not provide fast convergence based on the analysis in Section II-D.

**E. Initialization of \( u^{(0)}_j \)**

Unfortunately, \( u^{(0)}_j \) in (27) is available only for \( n \geq 1 \), i.e., after updating all voxels once. To define the initial update factors \( u^{(0)}_j \), we apply edge and intensity detectors to an initial filtered back-projection (FBP) image. This is reasonable since the initial FBP image is a good low-frequency estimate, so the difference between initial and final image will usually be larger near edges. We investigated one particular linear combination of edge and intensity information from an initial image. We used the 2D Sobel operator to approximate the gradient of the image within each transaxial plane. Then we normalized both the approximated gradient and the intensity of the initial image, and computed a linear combination of two arrays with a ratio : 1 for the initial update-needed factor \( u^{(0)}_j \), followed by DRA method. We have tried other linear combinations with different ratios, but the ratio : 1 provided the fastest convergence rate in our experiments.

**F. Implementation**

The dependence of \( \lambda^{(n)}_{ij} \) on iteration \( n \) increases computation, but we found two practical ways to reduce the burden. First, we found that it suffices to update \( \tilde{u}^{(n)}_j \) (and \( \tilde{a}^{(n)}_j \)) every \( n_{\text{loop}} \geq 1 \) iterations instead of every iteration. This is reasonable since the update-needed factors usually change slowly with iteration. In this case, we must generate a surrogate with the maximum curvature (10) to guarantee the majorization condition (6) for all iterations. Second, we compute the NU-based denominator (29) simultaneously with the majorization condition (6) for all iterations. Second, we compute the NU-based denominator (29) simultaneously with the data-fit gradient in (17). In 3D CT, we use forward and back-projectors that compute elements of the system matrix \( A \) on the fly, and those elements are used for the gradient \( \nabla L(x) \) in (17). For efficiency, we reuse those computed elements of \( A \) for the NU-based denominator (29). We implemented this using modified separable footprint projector subroutines [43] that take two inputs and project (or back-project) both. This approach required only 25\% more computation time than a single forward projection rather than doubling the time (see Table I). Combining this approach with \( n_{\text{loop}} = 3 \) yields a NU-SQS algorithm that required only 13\% more computation time per iteration than standard SQS, but converges faster.

**TABLE I**

<table>
<thead>
<tr>
<th>Run time of one iteration of NU-OS-SQS(8/2 subsets) for different choice of ( n_{\text{loop}} ) for GE performance phantom.</th>
</tr>
</thead>
<tbody>
<tr>
<td>( n_{\text{loop}} )</td>
</tr>
<tr>
<td>-------------------</td>
</tr>
<tr>
<td>1 Iter. [sec]</td>
</tr>
<tr>
<td>( n_{\text{loop}} )</td>
</tr>
</tbody>
</table>

Computing \( \tilde{u}^{(n)}_j \) and the corresponding NU-based denominator requires one iteration each. In the proposed algorithm, we computed \( \tilde{u}^{(n)}_j \) during one iteration, and then computed the NU-based denominator (29) during the next iteration combined with the gradient computation \( \nabla L(x) \). Then we used the denominator for \( n_{\text{loop}} \) iterations and then compute \( \tilde{u}^{(n)}_j \) again to loop the process (see outline in Appendix B).
IV. IMPROVED ORDERED SUBSETS (OS) ALGORITHM FOR HELICAL CT

Ordered subsets (OS) methods can accelerate algorithms by a factor of the number of subsets in early iterations, by using a subset of the measured data for each subset update. However, in practice, OS methods break the monotonicity of SQS and NU-SQS, and typically approach a limit-cycle looping around the optimum. This section describes a simple idea that reduces this problem, only slightly affecting the convergence rate unlike previous convergent OS algorithms. In helical CT geometries, we observed that conventional OS algorithms for PL and PWS problem are unstable for large subset numbers as they did not consider their non-uniform sampling. Thus, we describe an improved OS algorithm that is more stable for helical CT.

A. Ordinary OS algorithm

An OS algorithm (with \(M\) subsets) for accelerating the SQS or NU-SQS updates (17) has the following \(m\)th sub-iteration within the \(n\)th iteration using the denominator\(^3\) \(\hat{d}^{(n)}_j\) in (33):

\[
x^{(n+m+1)}_j = x^{(n+m)}_j - \frac{1}{\hat{d}^{(n)}_j} \left( \hat{\gamma}^{(n+m)}_j \frac{\partial}{\partial x_j} L_m \left( x^{(n+m)}_j \right) + \frac{\partial}{\partial x_j} R \left( x^{(n+m)}_j \right) \right),
\]

where \(\hat{\gamma}^{(n+m)}_j\) scales the gradient of a subset data-fit term \(L_m(x) = \sum_{i \in S_m} h_i \left( |Ax|_i \right)\), and \(S_m\) consists of projection views in \(m\)th subset for \(m = 0, 1, \ldots, M-1\). We count one iteration when all \(M\) subsets are used once, since the projection \(A\) used for computing data-fit gradients is the dominant operation in SQS iteration.

If we use many subsets to attempt a big acceleration in OS algorithm, some issues arise. The increased computation for the gradient of regularizer in (39) can become a bottleneck.

\(^3\)We consider the maximum curvature (10) here for computational efficiency in OS methods.

B. Proposed OS algorithm in helical CT

The constant scaling factor \(\gamma = M\) used in the ordinary regularized OS algorithm is reasonable when all the voxels are sampled uniformly by the projection views in all the subsets. But in geometries like helical CT, the voxels are non-uniformly sampled. In particular, voxels outside the ROI are sampled by fewer projection views than voxels within the ROI (see Fig. 1). So some subsets make no contribution to such voxels, i.e., very imbalanced subsets. We propose to use a voxel-based scaling factor \(\gamma_j\) that considers the non-uniform sampling, rather than a constant factor \(\gamma\).

Fig. 3. Effect of gradient scaling in regularized OS-SQS algorithm with GE performance phantom (GEPP) in helical CT: Each image is reconstructed after running 20 iterations of OS algorithm with 328 subsets, using ordinary and proposed scaling approaches. Standard deviation \(\sigma\) of a uniform region (in white box) is computed for comparison. We compute full-width half maximum (FWHM) of a tungsten wire (red arrow) to measure the resolution. (The result of a convergent algorithm is shown for reference. Images are cropped for better visualization.)
After investigating several candidates, we focused on the following scaling factor:

$$\gamma_j = \sum_{m=0}^{M-1} I \left\{ \sum_{i \in S_m} \tilde{c}_i^{(n_i)} a_{ij} \left( \sum_{i=1}^{N_p} a_{ij} x_i^{(0)} \right) > 0 \right\}, \quad (42)$$

where $I_B = 1$ if $B$ is true or 0 otherwise. As expected, $\gamma_j < M$ for voxels outside the ROI and $\gamma_j = M$ for voxels within the ROI. The scaling factor (42) has small computer overhead as it can be computed simultaneously with the precomputation of the initial data-fit denominator (29) by rewriting it as

$$\tilde{d}_j^{(0)} = \frac{1}{a_{ij}^{(0)}} \left( \sum_{m=0}^{M-1} \sum_{i \in S_m} c_i^{(n_i)} \left( \sum_{l=1}^{N_p} a_{il} \tilde{x}_l^{(0)} \right) \right). \quad (43)$$

We store (42) as a short integer for each voxel outside the ROI only, so it does not require very much memory.

We evaluated the OS algorithm with the proposed scaling factors (42) using the GE performance phantom. Fig. 3 shows that the OS algorithm using the proposed scaling factors (42) leads to more stable reconstruction than the ordinary OS approach which diverges outside the ROI. The instability seen with the ordinary OS approach may also degrade image quality within the ROI as seen by the noise standard deviations in Fig. 3. The results in Fig. 4 further show that the ordinary OS algorithm exhibits more variations within the ROI due to the instability outside ROI, whereas the proposed OS algorithm is robust.

This section discusses one practical method that can reduce noise and RMSD. As a practical alternative, we investigated an approach where the final image is formed by averaging all of the sub-iterations at the final iteration $n_{\text{end}}$ of the OS algorithm (after it approaches its limit cycle). A memory-efficient implementation of this approach uses a recursive in-place calculation:

$$\bar{x}^{(m+1)} = \frac{m}{m+1} \bar{x}^{(m)} + \frac{1}{m+1} \bar{x}^{(n_{\text{end}}-1 + m+1)}, \quad (44)$$

where $\bar{x}^{(0)}$ is an initial zero image, and $\bar{x}^{(1)}$ is the final averaged image. There was a preliminary simulation investigation of averaging the final iteration in [27], and we applied the averaging technique to CT scans here. In Table II, we investigated this averaging method using a scan of the GEPP phantom and quantified the noise and resolution properties (as described in Fig. 3), and evaluated root mean square difference (RMSD) between current and converged image within ROI. Table II shows that the averaging technique successfully reduces the noise and RMSD.

<table>
<thead>
<tr>
<th></th>
<th>Smoothed</th>
<th>OS-SQS(328)</th>
<th>Conv.</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>FBP</td>
<td>w/o averaging</td>
<td>w/ averaging</td>
</tr>
<tr>
<td>Mean [HU]</td>
<td>1127.7</td>
<td>1123.3</td>
<td>1123.8</td>
</tr>
<tr>
<td>Std. Dev. [HU]</td>
<td>2.3</td>
<td>8.0</td>
<td>7.2</td>
</tr>
<tr>
<td>FWHM [mm]</td>
<td>1.4</td>
<td>0.7</td>
<td>0.7</td>
</tr>
<tr>
<td>RMSD [HU]</td>
<td>9.4</td>
<td>3.4</td>
<td>0.8</td>
</tr>
</tbody>
</table>

| TABLE II | GE PERFORMANCE PHANTOM: NOISE, RESOLUTION AND RMSD BEHAVIOR OF OS-SQS(328 SUBSETS) AFTER 20 ITERATIONS FOLLOWED BY AVERAGING. |

Overall, we have enhanced the standard OS-SQS algorithm into the NU-OS-SQS method for 3D helical CT. First, we accelerated the standard OS-SQS algorithm by non-uniform (NU) approach, encouraging larger step sizes for the voxels that need more updates. We modified the algorithm to handle the helical CT geometry by introducing the scaling factor $\gamma_j$. We also averaged all sub-iterations at the final iteration to reduce noise. The outline of the proposed algorithm is presented in Appendix B. We investigate the performance of the NU-OS-SQS algorithm for various CT scans in the next section.

V. EXPERIMENTAL RESULTS

We investigated the proposed NU-OS-SQS algorithm for PWLS image reconstruction with a non-negativity constraint. The PWLS cost function is strictly convex and has a unique global minimizer [46]. We implemented the NU-OS-SQS algorithm in C and executed it on a Mac with two 2.26GHz quad-core Intel Xeon processors and a 16GB RAM. We used 16 threads, and projection views were grouped and assigned to each thread.

Three 3D helical CT data sets are used in this section to compare the proposed NU-OS-SQS algorithm to the ordinary OS-SQS algorithm, and we used the GE performance phantom (GEPP) to measure the resolution. We used two other clinical data sets to investigate the performance of NU approach. We

$$\text{RMSD} \triangleq \frac{\| x^{(n)}_{\text{ROI}} - x^{(\infty)}_{\text{ROI}} \|_2}{\sqrt{N_{\text{ROI}}}} \text{ [HU]},$$

where $N_{\text{ROI}}$ is the number of voxels in the ROI.
investigated tuning the DRA function $g(\cdot)$ in (37) to provide fast convergence rate for various data sets. We also provide results from a simulation data set in a supplementary material for reproducibility.\textsuperscript{5}

We chose the parameters of the cost function $\Psi(x)$ in (2) to provide a good image. We defined an edge-preserving potential function as $\psi_k((Cx)_k) \triangleq \tilde{\omega}_k \psi((Cx)_k)$, where the function:\textsuperscript{6}

$$\psi(t) = \frac{\delta^2}{b^2} \left( \frac{a b^2}{2} \left| \frac{t}{\delta} \right|^2 + b(b-a) \left( \frac{t}{\delta} \right) + (a-b) \log \left( 1 + b \left| \frac{t}{\delta} \right| \right) \right)$$

is a generalized version of a Fair potential function in [47], and the spatial weighting $\tilde{\omega}_k$ [48] provides resolution properties that emulate the GE product “Veo”. We used $M = 82$ subsets for the OS algorithms, assigning 12 out of 984 projection views per rotation to each subset. We used the maximum curvature (10) for generating the denominator of surrogate function of the cost function $\Psi(x)$, and focused on $n_{\text{loop}} = 3$ which balances the convergence rate and run time, based on Table I.

In Section II-D, we recommended fixing the denominator $\tilde{d}_j^{(n)}$ (generated by the maximum curvature (10)) after $n_{\text{fix}}$ iterations in NU-SQS algorithm to guarantee convergence. This condition is less important theoretically when we accelerate the NU-SQS algorithm with OS methods that break the convergence property. However, we still recommend fixing $\tilde{d}_j^{(n)}$ after $n_{\text{fix}}$ iterations (before approaching the limit-cycle) in the NU-OS-SQS algorithm, because we observed some instability from updating $\tilde{d}_j^{(n)}$ and $\tilde{u}_j^{(n)}$ every $n_{\text{loop}}$ iterations near the limit-cycle in our experiments. We selected $n_{\text{fix}} = 7$ for GEPP, but we did not use $n_{\text{fix}}$ for other two cases because the algorithm did not reach a limit-cycle within $n_{\text{end}} = 20$ iterations, and we leave optimizing $n_{\text{fix}}$ as a future work.

In Section IV-B, we stabilized the OS-SQS algorithm outside ROI in helical geometry by using the factor $\gamma_j$ in (42).

\textsuperscript{5}The supplementary material is available at http://ieeexplore.ieee.org.

\textsuperscript{6}The gradient $\dot{\psi}(t) = \frac{a b^2}{2} \frac{t}{\delta} + b(b-a) \frac{1}{\delta} + (a-b) \frac{1}{\delta} \log \left( 1 + b \left| \frac{t}{\delta} \right| \right)$ avoids expensive power operations, saving computation for OS-type methods. The function reduces to a Fair potential function in [47] when $a = 0$ and $b = 1$. We used $a = 0.0558$, $b = 1.6395$, and $\delta = 10$ in our experiments.

However, we experienced some instability outside ROI in NU-OS-SQS methods even with (42), because a small NU denominator $\tilde{d}_j^{(n)}$ outside ROI is more likely to lead to instability than for voxels within the ROI due to the incomplete sampling outside ROI. Therefore, we prevent the denominator $\tilde{d}_j^{(n)}$ outside ROI from being very small. We empirically modified the DRA function in Section III-C, and used it for our experiments, improving stability outside ROI. We first modified the function (35) as follows:

$$F^{(n)}_{\text{cdf}}(u) \triangleq \frac{1}{N_P} \sum_{j=1}^{N_P} I_{\left\{ \gamma_j \tilde{u}_j^{(n)} \leq u \right\}} , \quad n = 1, 2, \ldots, 5$$

since the value of $\tilde{u}_j^{(n)}$ in (27) outside ROI was found to be relatively large due to the incomplete sampling. We further modified (36) and (37) to prevent $\tilde{d}_j^{(n)}$ from becoming very small outside ROI as follows with $g(v; \alpha) \triangleq \max\{\alpha v, c\}$:

$$\tilde{d}_j^{(n)} \triangleq \begin{cases} 
  g \left( F_{\text{cdf}}(\gamma_j \tilde{u}_j^{(n)}); 1 \right), & \text{if } j \text{th voxel within ROI}, \\
  g \left( F_{\text{cdf}}(\gamma_j \tilde{u}_j^{(n)}); 0.5 \right), & \text{otherwise}.
\end{cases}$$

A. GE performance phantom (GEPP)

We reconstructed $512 \times 512 \times 47$ images of the GEPP from a $888 \times 64 \times 3071$ sinogram (the number of detector columns $\times$ detector rows $\times$ projection views) with pitch 0.5. We evaluated the full width at half maximum (FWHM) of a tungsten wire (see Fig. 3). Fig. 5(a) shows the resolution versus run time and confirms that non-uniform (NU) approach accelerates the SQS algorithm. This dramatic speed-up in FWHM is promising since SQS-type algorithms are known to have slow convergence rate of high frequency components [6]. We also evaluated the convergence rate by computing RMSD between current and converged\textsuperscript{7} image versus run time, within ROI.

\textsuperscript{7}We ran 100 iterations of OS-SQS algorithm with 41 subsets, followed by each 100 iterations of OS-SQS algorithm with 4 subsets, and 2000 iterations of (convergent) SQS. We subsequently performed 100 iterations of (convergent) NH-ABCD-SQS [21] to generate (almost) converged images $x^{(\infty)}$.

Fig. 5. GE performance phantom: plots of (a) FWHM and (b) RMSD as a function of run time for different choice of DRA parameters $t$ for $\epsilon = 0.05$. The plot markers show each iteration. There are no changes during first iterations, since we consider precomputing the denominator using one forward and back projections as one iteration.
Fig. 6. A shoulder region scan: plot of RMSD versus run time for different choice of parameters (a) $t$ and (b) $\epsilon$ in $g(v) = \max\{v_t, \epsilon\}$.

Fig. 5(a) and 5(b) illustrate that increasing $t$ in $g(\cdot)$ in (37) accelerates the convergence of “update-needed” region, particularly the wire and edges in GEPP. However, highly focusing the updates on few voxels will not help speed up the overall convergence for all objects. Therefore, we further investigate the choice of $g(\cdot)$ using various patient CT scans.

The RMSD plots\(^8\) of NU-OS-SQS in Fig. 5(b) reached a limit-cycle after 1500 sec that did not approach zero. Averaging the sub-iterations at the final iteration improved the final image with small computation cost, yielding the drop in RMSD at the last 20th iteration in Fig. 5(b). The reduced noise was measurable in the reconstructed image as seen in Table II.

\(\text{B. A shoulder region scan}\)

In this experiment, we reconstructed a $512 \times 512 \times 109$ image from a shoulder region scan $888 \times 32 \times 7146$ sinogram with pitch 0.5. Fig. 6(a) and 6(b) show that the non-uniform approach accelerates convergence, depending on the choice of parameters in $g(\cdot)$. We investigated the relationship between the convergence rate and the DRA function $g(\cdot)$ by tuning both the parameters $t$ and $\epsilon$ in (37). Fig. 6(a) shows that increasing $t$ to 10 accelerated convergence, but larger $t$ values did not help as the choice of $t = 40$ was slower than $t = 10$. In Fig. 6(b), decreasing $\epsilon$ to 0.01 accelerated the algorithm in this shoulder

\(^8\)We also provide the plots of the cost function for GEPP and shoulder region scan in the supplementary material.

Fig. 7. A shoulder region scan: (a) A center slice of initial FBP, converged image and reconstructed image by OS-SQS(82) and NU-OS-SQS(82)-
$g(v) = \max\{v^{10}, 0.05\}$ after about 95 min. (b) Difference between the reconstructed and converged images are additionally shown to illustrate the acceleration of NU approach. (Images are cropped for better visualization.)
region scan, but not for the data set in Section V-C, so \( \epsilon = 0.05 \) appears to be a reasonable choice overall.

We averaged the sub-iterations at the last iteration, but Fig. 6(a) and 6(b) did not show a drop at the final iteration (which appeared in Fig. 5(b)), because the algorithm had not yet reached a limit-cycle. Even though averaging did not noticeably decrease the RMSD, the reconstructed image had measurable noise reduction in regions that already reached a limit-cycle like uniform regions. (Results not shown.)

In Fig. 7(a), we illustrate that statistical image reconstruction can reduce noise and preserve image features compared to analytical FBP reconstruction. The reconstructed images of (NU-)OS-SQS show that NU approach helps OS-SQS to approach the converged image faster than the ordinary method. After the same computation (95 min.), the reconstructed image of OS-SQS still contains streaks from the initial FBP image, while NU-OS-SQS has reduced the streaks. This is apparent in the difference images between the reconstructed and converged images in Fig. 7(b).

By analyzing NU-OS-SQS in two CT scans, we observed that the parameters \( \ell = 10 \) and \( \epsilon = 0.05 \) consistently accelerated the algorithm by about a factor of more than two.\(^9\) (The choice \( \epsilon = 0.01 \) was too aggressive in our other experiments.) We also have observed more than two-fold accelerations in other experiments. (Results not shown.) Fig. 6(b) shows the

\(^9\)We used the run time and RMSD of standard OS-SQS after 20 iterations (without averaging) as a reference to compare with the NU-OS-SQS for each data set. Then we compared the run time of NU-OS-SQS that is required for achieving the reference RMSD with the reference run time, and confirmed that NU provided more than two-fold accelerations in two CT scans.

RMSD plot using the (practically unavailable) oracle update-needed factor \( \tilde{u}_j^{(0)} \) instead of our heuristic choice \( \tilde{u}_j^{(n)} \). This result suggests that additional optimization of the DRA method and initialization of \( \tilde{u}_j^{(0)} \) could further speed up the NU algorithm in future work.

C. A truncated abdomen scan

We also reconstructed a 390 × 390 × 239 image from a 888 × 64 × 3516 sinogram with pitch 1.0. This scan contains transaxial truncation and the initial FBP image has truncation artifacts [49] that can be reduced by iterative reconstruction. The choice of \( \tilde{u}_j^{(0)} \) described in Section III-E did not consider truncation effects, and we found that NU-OS-SQS did not reduce such artifacts faster than standard OS-SQS. (The large patient size may also have reduced the possible speed-up by the NU method, compared to the previous two scans.) Therefore, we investigated an alternative NU method that can reduce truncation artifacts faster than standard algorithm.

We designed a modified NU method using a few \( m_{\text{sub}} \) sub-iterations of standard OS-SQS to generate the initial update-needed factor \( \tilde{u}_j^{(0)} \), which may also be a reasonable approach for other scans. We perform initial sub-iterations \( x_j^{(m/M)} \) in (39) efficiently using two-input projectors (in Section III-F) and replacing the all-view denominator \( d_{j,\text{sub}}^{(m)} \) in (29) by a standard subset-based denominator [25]:

\[
\tilde{d}_{j,\text{sub}}^{(m)} = \gamma \sum_{i \in S_m} a_{ij} \left( \sum_{l=1}^{N_p} a_{il} \right),
\]  

(48)
where $S_m$ consists of projection views in $m$th subset. The scaling factor $\gamma_j$ in (42) is unavailable at this point, so we use $\gamma = M$ instead. After $m_{\text{sub}}$ sub-iterations, we compute the following initial update-needed factors:

$$u_j^{(0)} \triangleq f_{m_{\text{sub}}}^{(m_{\text{sub}})} \left( \left[ x_{j_{m_{\text{sub}}}}^{(m_{\text{sub}})} - x_j^{(0)} \right] \right),$$

(49)

where $f_{m_{\text{sub}}}^{(m_{\text{sub}})}(\cdot)$ is a DRA function in (36), and we use these to compute the NU denominators $d_j^{L,(0)}$ and $d_j^{R,(0)}$ that we use for first $n_{\text{loop}}$ outer iterations.

Fig. 8(a) shows that statistical image reconstruction provides better image quality than FBP reconstruction. Fig. 8(b) illustrates that this NUsub-OS-SQS approach reduces the truncation artifacts faster than the standard OS-SQS and NUOS-SQS. Although standard OS-SQS reduces noise faster than other two algorithms in Fig. 8(b), both NU-OS-SQS and NUsub-OS-SQS show better convergence near the spine, the boundary of patient, and other internal structures than OS-SQS at the same computation time (90 min.).

VI. CONCLUSION

This paper has presented a spatially non-uniform SQS algorithm that can efficiently minimize both PL and PWLS problems monotonically. The experimental results show that the proposed NU-SQS approach converged more than twice as fast as SQS. The OS algorithm, further applied to SQS method for acceleration, was enhanced to handle non-uniformly sampled geometries such as helical CT. The improvements showed promising results on large 3D helical CT data sets.

The key of the NU-SQS approach is designing “update-needed” factors $u_j^{(n)}$ in (27) that encourage larger step sizes for voxels that are predicted to need larger changes to reach the final image. Further optimization of these factors, e.g., by improving the initialization of $\tilde{u}_j^{(0)}$ and the DRA function in (36), should lead to further acceleration and stability of the proposed NU-SQS and NU-OS-SQS methods.

ACKNOWLEDGMENT

The first author would like to thank Sathish Ramani for suggesting Lemma 1. The authors would like to thank anonymous reviewers for their valuable comments. J. A. Fessler acknowledges helpful discussions with Johan Nuysts about the denominators for OS methods.

APPENDIX A

PROOF OF LEMMA 2

The proposed choice $\lambda_{ij}^{(n)} = a_{ij}u_j^{(n)}$ in (28) and its corresponding $d_j^{L,(n)}$ in (29) are a choice that minimizes

$$\sum_{j=1}^{N_p} \left( u_j^{(n)} \right)^2 d_j^{L,(n)}$$

among all possible $d_j^{L,(n)}$ in (19), i.e.,

$$\left\{ \lambda_{ij}^{(n)} \right\} = \arg \min_{\lambda_{ij}^{(n)}} \sum_{j=1}^{N_p} \left( u_j^{(n)} \right)^2 \left( \sum_{i=1}^{N_d} \right)_{a_{ij} \neq 0} \frac{c_{ij}^{(n)}}{\lambda_{ij}^{(n)}} \sum_{l=1}^{N_p} \lambda_l^{(n)},$$

subject to the positivity constraint on $\lambda_{ij}^{(n)}$ if $a_{ij} \neq 0$.

Proof: By the Schwarz inequality $\langle s, t \rangle^2 \leq ||s||^2 ||t||^2$, we have

$$\sum_{j=1}^{N_p} \left( a_{ij}u_j^{(n)} \right)^2 \leq \sum_{j=1}^{N_p} \left( u_j^{(n)} \right)^2 \frac{a_{ij}^2}{\lambda_{ij}^{(n)}} \sum_{l=1}^{N_p} \lambda_l^{(n)},$$

where $s_j = \sqrt{\sum_{l=1}^{N_p} \lambda_l^{(n)}}$ and $t_j = a_{ij}u_j^{(n)}$. Then,

$$\sum_{j=1}^{N_p} \left( u_j^{(n)} \right)^2 d_j^{L,(n)}$$

$$= \sum_{i=1}^{N_d} \sum_{j=1}^{N_p} \left( u_j^{(n)} \right)^2 a_{ij}^2 \frac{\sum_{l=1}^{N_p} \lambda_l^{(n)}}{\lambda_{ij}^{(n)}}$$

$$\geq \sum_{i=1}^{N_d} \sum_{j=1}^{N_p} a_{ij}u_j^{(n)}$$

$$= \sum_{j=1}^{N_p} \left( u_j^{(n)} \right)^2 d_j^{L,(n)}.$$

APPENDIX B

OUTLINE OF THE PROPOSED NU-OS-SQS ALGORITHM

Set $M$, $n_{\text{end}}$, $n_{\text{loop}}$ and initialize $x$ by an FBP image. Generate $u_j$ from an FBP image by edge and intensity detectors. Compute the maximum curvature $\tilde{c}_i = \max \left\{ h_i(0), \eta_j \right\}$.

$$d_j^L = 0, \gamma_j = 0, x_{j,\text{ref}} = x_j, \text{and the final image } \hat{x}_j = 0.$$

$$d_j^{R} = \frac{1}{u_j} \sum_{k=1}^{N} \psi_k(0)|c_{ij}^{(n)}| \sum_{l=1}^{N_p} c_{kl}|u_l|$$

(50)

$$n = 0$$

for $m = 0, 1, \ldots, M - 1$

$$d_j^{L,\text{sub}} = \frac{1}{u_j} \sum_{i \in S_m} \tilde{c}_ia_{ij} \sum_{l=1}^{N_p} a_{il}|u_l|$$

(51)

$$d_j^{L,+} = d_j^{L,\text{sub}} \text{ and } \gamma_j = 1 \{ d_j^{L,\text{sub}} > 0 \}
$$

end
for \( n = 1, 2, \ldots, n_{\text{end}} - 1 \)

if \( n \mod n_{\text{loop}} = 1 \) and \( n \leq n_{\text{fix}} \)

\[
d_{j}^{n} = d_{j}^{n_{\text{fix}}}, \quad \text{compute } d_{j}^{n} \text{ by (50)}
\]

else if \( n \mod n_{\text{loop}} = n_{\text{loop}} - 1 \) and \( n \leq n_{\text{fix}} - 2 \)

\[
x_{j, \text{ref}} = x_{j}
\]

else if \( n \mod n_{\text{loop}} = 0 \) and \( n \leq n_{\text{fix}} - 1 \)

\[
d_{j}^{n} = 0, \quad \text{and } u_{j} = g(F_{\text{data}}(x_{j} - x_{j, \text{ref}}))
\]

end

for \( m = 0, 1, \ldots, M - 1 \)

\[
x_{j, \text{new}} = x_{j}
\]

if \( n \mod n_{\text{loop}} \neq 0 \) or \( n > n_{\text{fix}} \)

\[
g_{\text{sub}}^{j} = \frac{\partial}{\partial x_{j}} L_{m}(x_{\text{prev}})
\]

else

compute both \( d_{j, \text{sub}}^{n} \) by (51) and \( g_{j, \text{sub}}^{n} \) by (52)

simultaneously using two-input projection function, and

\[
d_{j}^{n} = d_{j, \text{sub}}^{n}
\]

end

\[
x_{j} = \left[ x_{j, \text{prev}} - \frac{\gamma_{j} g_{j, \text{sub}}^{n} + \partial R(x_{\text{prev}})}{d_{j}^{n} + \partial R_{j}} \right] +
\]

if \( n = n_{\text{end}} - 1 \)

\[
x_{j} = \frac{m}{m + 1} x_{j} + \frac{1}{m + 1} x_{j}
\]

end

end

REFERENCES


Accelerating ordered subsets image reconstruction for X-ray CT using spatially non-uniform optimization transfer: Supplementary material

Donghwan Kim, Student Member, IEEE, Debashish Pal, Member, IEEE, Jean-Baptiste Thibault, Member, IEEE, and Jeffrey A. Fessler, Fellow, IEEE

This material extends the result section of [1] by providing cost function plots and a simulation study of a helical scan of the XCAT phantom [2].

References to equations, tables, figures, bibliography are within this material unless they are specified.

I. COST FUNCTION

In [1], we computed root mean square difference (RMSD) within region-of-interest (ROI) to evaluate the convergence rate of the proposed algorithm. Another way to assess the convergence rate is computing the cost function \( \Psi(x) \) in [1, Eqn. (2)] at each iteration. We used the following metric:

\[
\xi^{(n)} = 20 \log_{10} \left( \frac{\Psi(x^{(n)}) - \Psi(x^{(\infty)})}{\Psi(x^{(\infty)})} \right) \text{ [dB]} \quad (1)
\]

to better visualize how the cost function decreases each iteration. We used double precision and triple for loops when accumulating \( \Psi(x^{(n)}) \) to ensure high accuracy.

Fig. 2 shows plots of \( \xi^{(n)} \) for the choices of parameters used in [1, Fig. 5 and 6] for two real 3D scans: GE performance phantom (GEPP) and shoulder region scan. Fig. 2(a) shows that for the GEPP case, the NU-OS methods decreased the cost function at about the same rate than the ordinary OS method, or even perhaps slightly slower. In contrast, when we plotted RMSD distance to the converged image within the ROI [1, Fig. 5], NU-OS converged significantly faster. The reason for this different behavior is that the cost function plot considers all voxels, even those outside the ROI which are not of interest clinically. It is known that OS methods are not guaranteed to converge and apparently the non-ROI voxels are either not converging or perhaps approaching a larger limit-cycle, presumably due to the poor sampling in the padded slices outside the ROI, even with the stabilizing methods outside ROI described in [1, Section V]. Therefore, cost function plots may not provide practical measures of convergence rate for OS methods, particularly with acceleration. Future research on trying to further stabilize the NU-OS-SQS algorithm outside the ROI also may be helpful.

The final drops at the right in Fig. 2(a) show that averaging sub-iterations at the last iteration, as described in [1, Section IV.C], can compensate for the limit-cycle, particularly outside the ROI.

Unlike Fig. 2(a), the plots in Fig. 2(b) and 2(c) of shoulder region scan look similar to the plots of RMSD within ROI in [1, Fig. 6]. The scan geometry of each data set might explain these behavior of cost function in Fig. 2, where the shoulder region scan is a helical scan with pitch 1.0 and 7 helical turns and thus the corresponding image space has relatively few voxels outside the ROI, compared with GEPP data that is acquired by a helical scan with pitch 0.5 and 3 helical turns. Therefore, we can expect the cost function of shoulder region scan to be less affected by instability outside the ROI. Slower convergence of NU-OS-SQS algorithm at early iterations in Fig. 2(c) means that some choices of initial update-needed factor \( \tilde{u}_{(0)} \) were not good enough for voxels outside the ROI. The effect of averaging at the last iterations is apparent in Fig. 2(b) and 2(c), because the instability outside the ROI is suppressed by the averaging.

II. SIMULATION DATA

A. Simulation data

We simulated a helical CT scan data by using XCAT phantom [2]. We first acquired a \( 1024 \times 1024 \times 154 \) XCAT phantom for 500 [mm] transaxial field-of-view (FOV) at 70 [keV], where \( \Delta_x = \Delta_y = 0.4883 \) [mm] and \( \Delta_z = 0.6250 \) [mm]. (See Fig. 1.)

Fig. 1. A simulated XCAT phantom: a center slice of \( 1024 \times 1024 \times 154 \) XCAT phantom. (Images are cropped for better visualization.)
We simulated a helical scan using the blank scan factor $b_1 = 10^6$ and the mean number of background events $r_i = 0$ with Poisson noise. The sinogram data is in $888 \times 64 \times 2934$ (the number of detector columns $\times$ detector rows $\times$ projection views) space with pitch 1.0. Then, we reconstructed a $512 \times 512 \times 154$ image where $\Delta_x = \Delta_y = 0.9766$[mm] and $\Delta_z = 0.6250$[mm] using the proposed NU-OS-SQS algorithm.

B. Results

We use a cost function that is similar to the cost function used in [1, Section V]. We solve a PWLS function with a potential function $\psi_k(t) \triangleq \bar{\omega}_k \psi(t)$ in [1, Eqn. (45)] using a spatial weighting parameter:

$$\bar{\omega}_k \triangleq 50 \cdot \prod_{j=1}^{N_p} \max \{ \kappa_j, 0.01 \kappa_{\max} \}$$

that provides uniform resolution properties [3], where

$$\kappa_j \triangleq \sqrt{\sum_{i=1}^{N_z} \alpha_{ij} W_i} \sum_{i=1}^{N_z} \alpha_{ij}$$

and the value of $\kappa_{\max} \triangleq \max_j \kappa_j$ is used in (2) to avoid under-regularizing some voxels with very small $\kappa_j$. Fig. 3 illustrates both RMSD within ROI and $\zeta^{(n)}$ versus computation time, which we run the algorithm on the machine described in [1].

In Fig. 3(a), we evaluated the convergence rate using RMSD within ROI between current and converged image, where the converged image was generated by many iterations of a (convergent) SQS. We used parameters of DRA function that are used in [1, Fig. 5 and 6], and we observed similar trends. We also illustrate the plot of $\zeta^{(n)}$ versus run time in Fig. 3(b), which looks very similar to Fig. 3(a). This is because we regularized relatively more than two other experiments in this simulation experiment, and thus instability outside the ROI that can be caused by NU-OS-SQS methods is not apparent here.

In Fig. 4(a), the reconstructed images of (NU-)OS-SQS show that NU method accelerates OS-SQS and reaches closer to the converged image after the same computation time (88 min.). This is apparent when comparing the difference images between the reconstructed and converged images in Fig. 4(b), particularly around the spine.

REFERENCES

Fig. 3. A simulated XCAT phantom: plots of (a) RMSD and (b) $\xi^{(n)}$ versus run time for different choice of parameters $t$ for $\epsilon = 0.05$ in $g(v) = \max \{ v^t, \epsilon \}$.

Fig. 4. A simulated XCAT phantom: (a) A center slice of reconstructed image by OS-SQS(82) and NU-OS-SQS(82)-$g(v) = \max \{ v^{10}, 0.05 \}$ after about 88 min. (b) Difference between the reconstructed and converged images are additionally shown to illustrate the acceleration of NU approach. (Images are cropped for better visualization.)