

Research Article

Sequential Quadratic Programming Method for Nonlinear Least Squares Estimation and Its Application

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In this study, we propose a direction-controlled nonlinear least squares estimation model that combines the penalty function and sequential quadratic programming. The least squares model is transformed into a sequential quadratic programming model, allowing for the iteration direction to be controlled. An ill-conditioned matrix is processed by our model; the least squares estimate, the ridge estimate, and the results are compared based on a combination of qualitative and quantitative analyses. For comparison, we use two equality indicators: estimated residual fluctuation of different methods and the deviation between estimated and true values. The root-mean-squared error and standard deviation are used for quantitative analysis. The results demonstrate that our proposed model has a smaller error than other methods. Our proposed model is thereby found to be effective and has high precision. It can obtain more precise results compared with other classical unwrapping algorithms, as shown by unwrapping using both simulated and real data from the Jining area in China.

1. Introduction

Most problems faced during the course of processing measurement data are nonlinear in nature. In the classical least squares approach, the nonlinear approximation function is expanded at an approximate value. This transformation can linearize the data, allowing the problem to be solved through the linear least squares approach. The precondition of this model is that initial parameter values must reach the adjusted value; otherwise, the model error is significant.

Nonlinear least squares is an optimization method used to solve nonlinear problems [1–4]. Sequential quadratic programming is another important and effective optimization method [5–7]. A nonlinear least squares problem can be transformed into a sequential quadratic programming model and then solved [8–11].

A direction-controlled estimation model is proposed in this paper. The nonlinear least squares problem is transformed into a sequential quadratic programming model. The iterative point enters the feasible region via the penalty function, after which a problem's optimal solution can be obtained through sequential quadratic programming [12–14]. In the model, the penalty function is used to reduce the constraints of the initial values [15-19]. Sequential quadratic programming possesses quadratic convergence. The model is contrasted with least squares, and ridge estimate and simulated data are used to check model performance [20-23]. The root-mean-squared error (RMSE) and standard deviation between the estimated and true values, and the ε value representing the quality of unwrapping, are taken as indices of performance [24, 25]. We conducted experiments that showed the feasibility and effectiveness of the model. Unwrapping experiments using real data from the Jining area in China show that our proposed algorithm achieves more precise results than the least squares unwrapping algorithms. Quantitative indexes include differences in RMSE between rewrapped results and the original wrapped phase, computation time, and $\tilde{\epsilon}$ values [26–28]. The sequential quadratic programming method achieves better results with respect to three indices.

2. Sequential Quadratic Programming Model

Quadratic optimization with equality constraints is defined as follows:

min
$$f(x) = \frac{1}{2}x^{T}Gx + g^{T}x$$

s.t. $Ax = b$, (1)
 $x \ge 0$

where $G \in R_{n \times n}$ is a symmetric matrix, $g \in R_{n \times 1}, A \in R_{m \times n}, b \in R_{m \times 1}$.

The Lagrange function of (1) is

$$L(x, u) = \frac{1}{2}x^{T}Gx + g^{T}x + v^{T}(Ax - b)$$
(2)

where v is the Lagrange multiplier. At \hat{x} , the Kuhn–Tucker (K–T) condition of (2) is

$$G\hat{x} + g + A^{T}v = 0$$

$$A\hat{x} = b$$
(3)

If G is a positive definite matrix, then the unique solution of (3) is

$$\hat{x} = -G^{-1}g + G^{-1}A^T \left(AG^{-1}A^T\right) \left(AG^{-1}g + b\right)$$
(4)

In the sequential quadratic programming model, an optimal subproblem is set up to search for the next feasible point in the current iterative point. d_k is the solution of the optimal subproblem, that is

$$x^{(k+1)} = x^{(k)} + \alpha_k d_k$$
 (5)

where $\alpha_k \in (0, 1]$. The process for more iterations is repeated, and the optimization solution can then be obtained.

The optimal subproblem of (1) can be defined as follows:

$$\min_{x} \quad \nabla f^{T} \left(x^{k} \right) \left(x - x^{k} \right) \\
+ \frac{1}{2} \left(x - x^{k} \right)^{T} G \left(x^{k} \right) \left(x - x^{k} \right) \\
s.t. \quad s \left(x^{k} \right) + \nabla s^{T} \left(x^{k} \right) \left(x - x^{k} \right) = 0$$
(6)

where s(x) = Ax - b, $\nabla f(x^k) = (\partial f/\partial x_1, \partial f/\partial x_2, \cdots, \partial f/\partial x_n)^T|_{x=x^k}$, $\nabla s(x^k) = (\partial s/\partial x_1, \partial s/\partial x_2, \cdots, \partial s/\partial x_n)^T|_{x=x^k}$

 $G(x^k)$ is the Hessian matrix of f(x):

$$G\left(x^{k}\right) = \nabla^{2} f\left(x^{k}\right) = \frac{\partial^{2} f}{\partial x^{2}}$$

$$= \begin{bmatrix} \frac{\partial^{2} f\left(x\right)}{\partial x_{1}^{2}} & \frac{\partial^{2} f\left(x\right)}{\partial x_{1} \partial x_{2}} & \cdots & \frac{\partial^{2} f\left(x\right)}{\partial x_{1} \partial x_{t}} \\ \frac{\partial^{2} f\left(x\right)}{\partial x_{1} \partial x_{2}} & \frac{\partial^{2} f\left(x\right)}{\partial x_{2}^{2}} & \cdots & \frac{\partial^{2} f\left(x\right)}{\partial x_{2} \partial x_{t}} \\ \cdots & \cdots & \cdots & \cdots \\ \frac{\partial^{2} f\left(x\right)}{\partial x_{1} \partial x_{t}} & \frac{\partial^{2} f\left(x\right)}{\partial x_{2} \partial x_{t}} & \cdots & \frac{\partial^{2} f\left(x\right)}{\partial x_{t}^{2}} \end{bmatrix} \Big|_{x=x^{k}}$$

$$(7)$$

The penalty function of model (1) is defined as follows:

$$\widetilde{P}(x,\pi) = \frac{1}{2}x^{T}Gx + g^{T}x + \pi ||Ax - b||$$
(8)

where $\pi > 0$ is the penalty coefficient.

If the objective and constraint function satisfy certain conditions, then the penalty method is convergent; i.e., the algorithm converges to the unique, optimal result [15, 29, 30].

The penalty function method is simple to compute. As the penalty coefficient increases, the initial point is iterated to the optimal solution. However, when the penalty coefficient is very large, computation time increases.

3. Sequential Quadratic Programming for Nonlinear Least Squares

Synthesizing both methods, the penalty function is simple to compute in earlier iterations, and the sequential quadratic programming method allows for quadratic convergence. A direction-controlled nonlinear least squares estimation algorithm using the penalty function and sequential quadratic programming is proposed herein. The least squares model is transformed into the quadratic optimization model. The iterative point enters the feasible region by the penalty function, and the optimal solution can then be obtained by sequential quadratic programming. Efficiency in problem solving can be improved by combining the two methods.

The nonlinear model is defined as follows:

$$L + \Delta = f(x) \tag{9}$$

where $L = (L_1, L_2, \dots, L_n)^T$ is an $n \times 1$ vector; f is a nonlinear function;

 $X = (X_1, X_2, \dots, X_t)^T$ is a $t \times 1$ vector to be estimated, and $\Delta = (\Delta_1, \Delta_2, \dots, \Delta_n)^T$ is an $n \times 1$ observational error vector. The error equation of (9) is

$$V(x) = f(x) - L \tag{10}$$

The least squares principle must be satisfied, that is,

$$\min F(x) = V^{T}(x) PV(x)$$

$$= [f(x) - L]^{T} P[f(x) - L]$$
(11)

Equation (11) can be converted to quadratic optimization with equality constraints as follows:

min
$$F(\Delta x^k) = V^T P V$$

s.t. $V = f(x^k) - L$
(12)

The subproblem of (12) can be obtained by linearizing the equality constraint at x^k . The expression is as follows:

min
$$F(\Delta x^{k}) = V^{T} P V$$

s.t. $V = B(x^{k}) \Delta x^{k} + C(x^{k}) - L$
(13)

	True value	Least squares method	Ridge method	Sequential quadratic programming method
	1	2.8552	0.8654	1.2205
	1	-1.3035	1.3872	1.2598
\hat{x}	1	1.2584	0.8283	1.1896
	1	2.3954	0.5180	0.9064
	1	-1.9306	1.1375	1.1034
RMSE		1.9672	0.2996	0.1852
The standard deviation		4.3989	0.6699	0.4141

TABLE 1: Estimated values by sequential quadratic programming, least squares, and the ridge estimate.

where $B(x^k) = \partial f(x)/\partial x|_{x=x^k}$, $C(x^k) = f(x^k)$, and the positive definite matrix *P* is the weight matrix. Equation (12) shows convex programming.

The penalty function of (12) is

$$\widetilde{P}(x,\pi) = V^T P V + \pi \left\| f(x) - L \right\|$$
(14)

The steps of the algorithm using the penalty function and sequential quadratic programming are as follows.

(1) The initial values are defined as $x^{(0)}$ and π_0 ; the control error is e > 0, $\varepsilon > 0$; and the number of iterations is k(k = 0).

(2) The initial value $x^{(0)}$ is substituted into (14) to calculate the iterative value $||f(x^{(k)}) - L||$. If $||f(x^{(k)}) - L|| < e$, the iteration terminates.

(3) If $||x^{(k+1)} - x^{(k)}|| \le \varepsilon$, the result is $x^* = x^{(k+1)}$; otherwise, set $\hat{x} = x^{(k)}$ and move to the next step.

(4) By taking $x^{(0)} = \hat{x}$ as the new initial value, (12) is solved using sequential quadratic programming. The correction $x^{(k)}$ value can then be obtained.

(5) If $||x^{(k+1)} - x^{(k)}|| \le \varepsilon$, the result is $x^* = x^{(k+1)}$; otherwise, return to step (4) until the condition is met.

In the course of running the algorithm, the penalty coefficient of (6) is as follows [31]:

$$\pi = \max\left(\left|\frac{\left[V^{T}\left(\Delta x^{k}\right)^{T}\right]\left(\begin{smallmatrix}P & 0\\ 0 & 0\end{smallmatrix}\right)\left(\begin{smallmatrix}V\\ \Delta x^{k}\end{smallmatrix}\right)}{\Delta x^{k}}\right|, 2\pi_{-} + \beta\right) \quad (15)$$

where $\beta \ge 0$ is the constant, and π_{-} is the penalty coefficient of the previous iteration.

4. Experimental Results and Analysis

4.1. Simulated Data. In this case, AX = L, where \widetilde{A}

ſ	1	

=

/ 2.0000	-5.0000	1.0000	1.0000	-9.5000	
-2.0000	4.0000	1.0000	-1.0500	8.5000	١
-2.0000	1.0000	1.0000	-1.0000	2.4000	
-1.0000	2.5000	4.0000	-0.5000	7.0000	
-1.0000	3.2000	4.0000	-0.5000	8.4000	(16)
1.0000	1.0000	-3.0000	0.4000	0.4900	·
3.0000	7.0000	-3.0000	1.5000	12.7000	
5.0000	-1.0000	-2.0000	2.5000	-3.0000	ļ
4.0000	2.0000	-2.0000	2.0100	3.0000	
4.0000	3.0000	-2.0000	2.0000	5.0000 /	



FIGURE 1: Ridge parameter from the L-curve.

The observations are $\tilde{L} = [-10.5 \ 10.45 \ 1.4 \ 12 \ 14.1 - 0.11 \ 21.2 \ 1.5 \ 9.01 \ 12]^T$.

 $X = [x_1 \ x_2 \ x_3 \ x_4 \ x_5]^T$ are unknown parameters. The true values are $\widehat{X} = [1 \ 1 \ 1 \ 1 \ 1]^T$.

The white noise is $\Delta \sim N(0, \sigma^2 I)$, where *I* denotes the identity matrix and $\sigma = 0.1$, $e = 10^{-2} \varepsilon = 10^{-2}$.

The condition number of A is 41397, so this is an illconditioned problem. Using sequential quadratic programming, least squares, and the ridge estimate, the estimated results can be compared, as shown in Table 1.

In Table 1, the ridge parameter of the ridge estimation is 1.4568. Figure 1 shows the ridge parameter from the L-curve [32].

In this ill-conditioned problem, all methods obtain estimated values. The distance between the initial and true values is large. Sequential quadratic programming obtains the closest solution (Table 1). The RMSE and standard deviation show that the sequential quadratic programming model is the best, and the ridge estimate method is better than the least squares method. In sequential quadratic programming, the iterative point enters the feasible region via the penalty function and the optimal solution can then be obtained via sequential quadratic programming. The constraints on the initial values of sequential quadratic programming are reduced; however, the computational cost is increased. Ridge



FIGURE 2: Unwrapping results obtained using the above-mentioned two methods.

estimation performs much better than least squares, but determining the ridge parameters is difficult.

4.2. Phase Unwrapping Experiments

4.2.1. Simulated Data. To verify the effectiveness of the algorithm, simulated data were used to conduct experiments. In a simulated interferogram with a size of 100 pixels × 100 pixels, a coherence coefficient of 0.85 and Gaussian noise with a standard deviation of 0.8160 radians were added to the image, $e = \varepsilon = 10^{-3}$. The least squares and sequential quadratic programming methods were used to unwrap images. Figure 2 shows the unwrapping result of each method.

Both methods obtain good unwrapping results, without lines or islands. The image obtained by sequential quadratic programming is smoother than that obtained by least squares and is more similar to the noise-free image.

Figure 3 shows a cross-sectional view of the unwrapping phase.

As shown in Figure 3, the abscissa represents the horizontal position (x-coordinate) of the pixel in the 50th line of the image. The ordinate denotes the phase. The maximum value of the cross-sectional view of sequential quadratic programming is close to the maximum value of the simulated surface. The maximum value of the cross-sectional view of the least squares method is far from the maximum value of



FIGURE 3: Fiftieth line cross-sectional view of the unwrapping phase using least squares and sequential quadratic programming.

TABLE 2.	Unwran	ning	time	RMSE	and	~ val	110
IABLE \angle .	Unwrap	ping	unic,	RIVIOE,	anu	e vai	uc.

Method	Computation time (s)	RMSE (rad)	$\widetilde{\epsilon}$ (rad)
Least squares	12.142	0.7594	0.6214
Sequential quadratic programming	11.866	0.5393	0.4228

TABLE 3: Selected parameters of ASAR image data.

Number	Sensor	Date	Туре	Latitude and longitude of center	Polarization	Ascending	Wavelength
022403	ENVISAT	20050306	SLC	116.314 6, 35.305 1	VV	Ascending	С
022404	ENVISAT	20050130	SLC	116.314 4, 35.305 1	VV	Ascending	С

the simulated surface. The curve obtained for of the least squares method is also much rougher than that of sequential quadratic programming.

Table 2 shows the running time, RMSE, and $\tilde{\epsilon}$ value of sequential quadratic programming and least squares obtained by phase unwrapping. Both methods result in relatively reliable unwrapping. The running time is similar, indicating that the efficiency of the two algorithms is almost the same. The RMSE of sequential quadratic programming is smaller than that of least squares, indicating that sequential quadratic programming is more precise. $\tilde{\epsilon}$ indicates the quality of the wrap; the smaller the $\tilde{\epsilon}$, the better the unwrapping quality. Thus, the sequential quadratic programming method provides better unwrapping than the least squares method.

The experiment with simulated data shows that sequential quadratic programming is a feasible method: the model is effective and has high precision. *4.2.2. Observational Data.* Satellite data collected by ENVI-SAT ASAR, covering the Jining area of Shandong province, China, were used to test the model's performance.

Table 3 shows the basic parameters of the data.

The images are 1024×1024 pixels, with a resolution of 30 m. Figure 4 shows the unwrapping result of each method, $e = \varepsilon = 10^{-3}$.

Both unwrapping methods were run successfully. In the interferogram, the stripe noise is large and the fringes are unclear. The phase ripple of sequential quadratic programming is smaller than that of least squares, demonstrating that sequential quadratic programming experiences less interference.

Table 4 shows the computation time, RMSE, and $\tilde{\varepsilon}$ value.

The RMSE and $\tilde{\epsilon}$ values of sequential quadratic programming are smaller than those of the least squares method. This



FIGURE 4: Unwrapping of ASAR images, using least squares and sequential quadratic programming.

TABLE 4: Unwrapping time, RMSE, and ε values.

Method	computation time (s)	RMSE (rad)	$\tilde{\epsilon}$ (rad)
Least squares	725.940	1.7594	0.6214
Sequential quadratic programming	1336.948	0.5393	0.4228

illustrates that the difference between the rewrapped results and the original wrapped phase is small. The unwrapping phases of the sequential quadratic programming method are smoother and more reliable. However, the computation time of sequential quadratic programming is greater than that of the least squares method. This is due to high levels of noise and the penalty function increasing the computation time of sequential quadratic programming.

5. Conclusions

A direction-controlled nonlinear least squares estimation algorithm using the sequential quadratic programming method is proposed herein. The least squares model is transformed into the nonlinear programming model, and the iteration direction can then be controlled by sequential quadratic programming. The ill-conditioned matrix and phase unwrapping are used to verify the method. The proposed algorithm achieves better results than alternative approaches in terms of RMSE and $\tilde{\epsilon}$ values, and its effectiveness is demonstrated using observational satellite data. In future research, the role change if the inequality constrains are taken into account will be considered.

Data Availability

The 20050130.slc and 20050306.slc data used to support the findings of this study are available from the corresponding author upon request.

Conflicts of Interest

The authors declare that they have no conflicts of interest.

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