# Conformal Mapping Using the Schwarz-Christoffel Transform 2007-2008 

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#### Abstract

The Schwarz-Christoffel transform is a conformal mapping from the upper half of the complex plane to a polygonal domain. It allows many physical problems posed on two-dimensional, polygonal regions, such as heat flow, fluid flow, and electrostatics, to be solved numerically. This type of problem cannot generally be solved in closed form; the Schwarz-Christoffel transform provides an exceptionally accurate method of solution. This paper describes the implementation of a working software unit that efficiently and accurately calculates Schwarz-Christoffel transforms and inverses. The program incorporates graphical, easy-to-use interfaces and will contain resources to aid in solving physical problems. Future research into mathematical extensions to the Schwarz-Christoffel transform, such as the inclusion of simple curves, will be conducted.


Keywords: Schwarz-Christoffel transform, conformal mapping, numerical analysis, Laplace's equation, fluid flow, heat flow

## 1 Introduction

Many physical problems are expressed as differential or boundary value problems over a surface. Often, these surfaces are or can be approximated by two-dimensional polygons. When this occurs, one method of determining accurate solutions is by assuming the polygonal domain exists in the complex plane and determining a conformal map, which preserves the structure of Laplace's equation, that re-
states the problem in a simpler domain. Here, the upper half-plane is used. A solution to the problem, now easy to solve analytically or numerically, is then mapped back to the original domain. For such polygonal domains, a method of determining the specific transform needed is provided by the following formula, known as the Schwarz-Christoffel transform:

$$
\begin{equation*}
f(z)=A \int_{0}^{z} \prod_{j=1}^{n}\left(\zeta-x_{j}\right)^{-\theta_{j} / \pi} d \zeta+B \tag{1}
\end{equation*}
$$

In this formula, $\zeta$ is an independent complex variable in the upper half-plane, the $\theta_{j}$ are the exterior angles of the polygon, the $x_{j}$ are prevertices of the mapping (given along the real axis), $n$ is the number of verticies of the polygon, and $A$ and $B$ are complex constants that specify the location, size, and orientation of the image polygon in the complex plane. The $\theta_{j}$ must satisfy

$$
\begin{equation*}
\sum_{j=1}^{n} \theta_{j}=2 \pi, \tag{2}
\end{equation*}
$$

which ensures the completeness of the image polygon [2]. Unfortunately, the Schwarz-Christoffel formula is not easy to evaluate, and requires both effective integration algorithms and an efficient, convergent method to solve a specific nonlinear equation. Implementation of such numerical routines is not a trivial problem, and is the subject of this paper.

The initial project may be divided into four separate problems. First, a method to effectively evaluate integrals of the form found in the Schwarz-Christoffel formula is required. Second, a numerical algorithm to solve the so-called Schwarz-Christoffel parameter problem, a system of nonlinear equations for the prevertices, must be developed. Third, methods to eval-
uate the forward and backward transforms based on given prevertices must be coded. Fourth, a user interface is needed, which should be robust and accessible to allow nonspecialists to systematically solve various physical problems. The four components may be coded simultaneously or in series, as they are by nature almost entirely separable problems. In this project, each part was coded in series.

Subsequent research will be done into improvements and optimizations to the numerical algorithms for various subproblems and extensions. These include the problem of mapping polygons with large aspect ratios, which are generally highly ill-conditioned, and the extension of the Schwarz-Christoffel formula to simple curves. The goal of the project is thus twofold: to produce a piece of software that will be useful in the solution of real, physical problems, and to improve upon current algorithms for producing the Schwarz-Christoffel transform.

## 2 Background

The Schwarz-Christoffel transform was first discovered independently in the late 1860s by Elwin Christoffel and Hermann Schwarz. Schwarz used some of the ideas of the transform to provide a more rigorous proof of the Riemann Mapping Theorem, which he had previously shown to be incomplete, but the majority of this work was on a purely theoretical level [5]. The usefulness of the transform was mitigated by the formula's unwieldiness, as the mappings for all but the simplest domains could not be calculated in closed form. Numerical estimates, especially for nonsymmetric polygons with four or more verticies, could not be effectively calculated by hand. Application to physical problems, therefore, was limited at best until the advent of the computer. A computer algorithm to compute the Schwarz-Christoffel transform was first written in the 1960s, and others have been written and modified since then [3].

The first problem in calculating the SchwarzChristoffel mapping is the evaluation of the integral given by Eq. (1). The integrand contains singularities at each of the endpoints of the image polygon, which tend to render ordinary numerical integration
routines either useless or hopelessly slow. In addition, the presence of negative powers in $f$ means that domains of applicability for each of the subfunctions $\left(\zeta-x_{j}\right)^{-\theta_{j} / \pi}$ must be chosen so that the entire domain in and immediately around the image polygon is meromorphic. Although several quadrature routines have been used for this problem, the method of choice today is Gauss-Jacobi quadrature, which uses a specially-tailored weighting function to choose points of evaluation and weights for the points that maximize efficiency. In practice, the SchwarzChristoffel formula is altered so that the last prevertex, $x_{n}$, is chosen to be both $-\infty$ and $+\infty$ (the values are equivalent for a conformal map, which acts on the Riemann sphere). This can always be done due to the extra degrees of freedom contained in Eq. (1). The integrals that must be evaluated in practice in the course of the Schwarz-Christoffel transform are of the form

$$
\begin{equation*}
\int_{x_{i-1}}^{x_{i}} \prod_{j=1}^{n-1}\left(\zeta-x_{j}\right)^{-\theta_{j} / \pi} d \zeta \tag{3}
\end{equation*}
$$

These integrals can always be written as required for Gauss-Jacobi quadrature; that is, in the form

$$
\begin{equation*}
\int_{a}^{b}(z-a)^{\alpha}(z-b)^{\beta} \psi(z) d z \tag{4}
\end{equation*}
$$

where $\alpha$ and $\beta$ are real numbers greater than -1 .
The points and weights of a Gauss-Jacobi quadrature are calculated here using a routine from Numerical Recipes [4] which efficiently estimates and solves for the roots of the Jacobi polynomials, which form the sample points just as the roots of the Chebyshev polynomials form the sample points for standard Gaussian quadrature. These points, however, are uniformly calculated in the range $[-1,1]$, and the integrals must be adjusted slightly to conform to this range. During the calculation of the prevertices, discussed below, the $z$ in Eq. (4) are restricted to the real axis; however, in direct calculations once the prevertices have been found, the $z$ will generally be fully complex.

The second problem is the Schwarz-Christoffel parameter problem, where the $x_{j}$ in Eq. (1) are calcu-
lated. As described in [2], a series of nonlinear, constrained equations can be formed from the requirement that the image polygon and the desired polygon be similar (the constants $A$ and $B$ in Eq. (1) then ensure congruency). Written out, there are $n-3$ linear equations in $n-3$ unknowns, once the extra degrees of freedom have been taken care of by arbitrarily giving three of the $x_{j}$ precise values. Here, as in the literature, we take $x_{1}=-1$ and $x_{2}=0$ in addition to the already-defined $x_{n}= \pm \infty$. The equations to be solved, ensuring that the target polygon and the image of the Schwarz-Christoffel map are similar, are then

$$
\begin{equation*}
\frac{\left|\int_{x_{i-1}}^{x_{i}} \prod_{j=1}^{n}\left(\zeta-x_{j}\right)^{-\theta_{j} / \pi} d \zeta\right|}{\left|\int_{x_{1}}^{x_{2}} \prod_{j=1}^{n}\left(\zeta-x_{j}\right)^{-\theta_{j} / \pi} d \zeta\right|}-\frac{\left|w_{j}-w_{j-1}\right|}{\left|w_{2}-w_{1}\right|}=0 \tag{5}
\end{equation*}
$$

where the $w$ are the target vertices, $i=3,4, \ldots, n-1$, and all other variables are as defined above. However, there is an additional complication, as preserving the order of the prevertices on the real axis is important. The extra constraint can be expressed as

$$
\begin{equation*}
1<x_{3}<x_{4}<\ldots<x_{n-1}<\infty \tag{6}
\end{equation*}
$$

The unconstrained problem is relatively easy to solve; however, the constraint prevents a naive application of a Newton's Method variant to this problem. To get around this, Trefethen in [3] suggests a simple change of variables that ensures the inequalities of Eq. (6). Take a new series of variables, $\chi_{j}$, and let

$$
\begin{equation*}
\chi_{j}=\ln \left(x_{j}-x_{j-1}\right) \tag{7}
\end{equation*}
$$

The resulting $\chi_{j}$ will automatically obey Eq. (6), and the original $x_{j}$ are found by the simple inverse formula

$$
\begin{equation*}
x_{j}=x_{j-1}+e^{\chi_{j}} \tag{8}
\end{equation*}
$$

This new set of equations in the $\chi_{j}$ is readily solved by a variant of Newton's Method that does not require the calculation of the Jacobian matrix (which would be hopelessly complex), but rather uses progressive estimates.

## 3 Development

The software has been written entirely in Java, although certain routines may be later written in C to increase speed if there is a bottleneck at any point in the process. The entire development of the program is designed to be achieved in stages by attacking the subproblems individually. The following is a list of classes, with short descriptions, written up to this point:

- class Complex - this class stores and performs arithmetic on complex numbers, which are not directly supported by Java. Several of the methods, including the multiplication and division algorithms, are designed to run as quickly as possible while avoiding intermediate overflow and floating-point error propagation. The multiplication method, for instance, requires only three real multiplications rather than four. (see Appendix for examples)
- class GaussJacobiWeights - this class calculates and stores the sample points and weights for a given Gauss-Jacobi quadrature over the interval $[-1,1]$. This routine uses Newton's Method to find the roots of the Jacobi polynomials, which are the sample points for the integral, and was taken and translated from [4].
- class SchwarzFunction - this class evaluates the integrand of a given real-valued SchwarzChristoffel integral, serving as a storage class for data of this kind.
- class GaussQuad - this class accepts as input $\psi$, $a, b, \alpha$, and $\beta$ from Eq. (4). For an arbitrary integral in that form, shifting and scaling the bounds produces the equivalent integral

$$
\begin{equation*}
c^{\alpha+\beta+1} \int_{-1}^{1}(\zeta-1)^{\alpha}(\zeta+1)^{\beta} \psi(c \zeta+m) d \zeta \tag{9}
\end{equation*}
$$

where $b=\frac{a+b}{2}$ and $c=b-m=m-a$. This integral is then evaluated using the sample points and weights given by the GaussJacobiWeights class and returned. For any GaussQuad object, varying numbers of sample points (and
thus varying accuracy) are accepted by its integrate() method. (see Appendix for example code)

- class RealNewtonRaphson - this class accepts an array of vertices and calculates the necessary prevertices as well as the constants $A$ and $B$ from Eq. (1). The method employs a standard Newton-Raphson method to solve the Eq. (5). At each step, an approximate Jacobian matrix for the function is calculated using a forwarddifference method in each dimension; the step vector is then solved for using an LU factorization on the equation

$$
\begin{equation*}
\mathbf{J} \delta \vec{x}=\vec{f} \tag{10}
\end{equation*}
$$

where $\mathbf{J}$ represents the Jacobian, $\delta \vec{x}$ the step vector, and $\vec{f}$ the current function vector. Note that by employing a forward-difference method to find the Jacobian, the number of function evaluations can be cut in half, as the current function vector can be reused in the Jacobian calculation. (see Appendix for example code)

- class ForwardGaussQuad - this class, using already-calculated values for the prevertices, evaluates the Schwarz-Christoffel integral at a given point. To minimize error caused by the presence of singularities near the path of the integral (the singularities at the endpoints are handled by the Gauss-Jacobi quadrature), the path of integration is divided recursively such that no segment is closer to a singularity than one-half its length, a technique employed in [3]. Such recursive subdivision is known as compound Gauss-Jacobi quadrature.
- class SchwarzChristoffel - this class runs the graphical user interface and calls RealNewtonRaphson and ForwardGaussQuad when necessary. The graph itself has the ability to show axes and manually adjust window parameters.

In future iterations of the project, a new set of routines will be implemented to calculate continuous Schwarz-Christoffel problems. Immediately following
from Eq. (3) above, we have

$$
\begin{equation*}
f^{\prime}(z)=A \prod_{j=1}^{n-1}\left(\zeta-x_{j}\right)^{-\theta_{j} / \pi} \tag{11}
\end{equation*}
$$

To change this into a continuous problem, we can rewrite this as

$$
\begin{equation*}
f^{\prime}(z)=A e^{\frac{1}{\pi} \sum_{j=1}^{n-1}-\theta_{j} \ln \left(z-x_{j}\right)} \tag{12}
\end{equation*}
$$

Then, defining the natural logarithm function as single-valued in the upper half-plane, except where $x_{i}=z, f^{\prime}$ becomes an analytic function in the required domain. To formulate the continuousboundary problem, we simply replce the sum in Eq. (12) with an integral, and integrate the entire function to find $f(z)$ :

$$
\begin{equation*}
f(z)=A \int_{0}^{z} e^{\frac{1}{\pi} \int_{-\infty}^{\infty}-\theta(x) \ln \left(\zeta-x_{j}\right) d x} d \zeta+B \tag{13}
\end{equation*}
$$

where $\theta(x)$ represents the amount of turning per unit length on the real axis, such that

$$
\begin{equation*}
\int_{-\infty}^{\infty} \theta(x) d x=2 \pi \tag{14}
\end{equation*}
$$

The continuous problem therefore has an extra subproblem to solve, namely, the solution of the integral equation, Eq. (13), to find $\theta(x)$ at every $x$.

The majority of testing of the program is specific to a single numerical routine; that is, each of the algorithmic components are tested individually. To calculate the GaussQuad routines, for instance, randomly generated sample problems are solved by MATLAB to provide an approximate check on the accuracy of solutions, then precision is achieved by manipulating the number of sample points used for the quadrature.

Shortly, a program specifically designed to track approximate error propagation and runtimes of each of the components will be developed. For the majority of the routines used in the program, strict error bounds can be calculated, and for the remaining algorithms error can be accurately estimated. For all subproblems, as well as the entire routine, plots of runtime versus precision will be generated to examine the efficacy of each routine. For the final program, random polygonal generation will be used to dynamically test the program for a range of inputs.

## 4 Results and Discussion

The purpose of this project was to calculate and display Schwarz-Christoffel transforms, which conformally map the upper half-plane to an arbitrary polygon, efficiently and accurately. In addition, additional research into the Schwarz-Christoffel transform itself, including its extension to curved target domains, was investigated. The evaluation of the Schwarz-Christoffel formula involves several parts, including the efficient calculation of a certain class of integrals as well as a solver of nonlinear systems of equations. Solving the continuous-parameter problem will require numerical solutions to a certain class of integral equations.

Results on accuracy versus time data are forthcoming; however, preliminary results have confirmed that answers are correct to within the modest relative error tolerance of about $10^{-4}$. Nevertheless, far greater accuracy is expected once dynamic testing begins.

The first problem, that of numerical integration, has been solved and refined, and a basic user interface has been designed. The second problem, that of a nonlinear equation solver to calculate the prevertices, has also been completed to satisfaction; current research focuses on correct implementation of the forward transform using given prevertices. Preliminary results indicate the general correctness but inexactitude of the forward transform in the absence of compound quadrature, especially near the boundaries of the given polygon. It is hoped that a full implementation of the compound quadrature will ameliorate these concerns.

The completed program will be useful on several levels: as a teaching aid, and as a tool for researchers solving certain equations on polygonal regions. Once the basic Schwarz-Christoffel problem is numerically solved, the program can form an easy basis for testing research in numerical analysis and mathematics that deals with improving or expanding the SchwarzChristoffel transform.

## Appendix

## Code for class Complex

```
public class Complex
{
    public Complex multiply(Complex z)
    {
        double temp1=x*z.real();
        double temp2=y*z.imag();
        return new Complex(temp1-temp2, (x+y)*(z.real()+z.imag())-temp1-temp2);
    }
    public Complex divide(Complex z)
    {
        double temp1=z.real()/z.imag();
        double temp2=z.imag()/z.real();
        if(Math.abs(z.real())>=Math.abs(z.imag()))
        {
            double denominator=z.real()+z.imag()*temp2;
            return new Complex((x+y*temp2)/denominator,(y-x*temp2)/denominator);
        }
        else
        double denominator=z.real()*temp1+z.imag();
            return new Complex((x*temp1+y)/denominator,(y*temp1-x)/denominator);
        }
    }
    public double modulus()
    {
        if ( }\textrm{x}==0&&&y==0
        return 0.0;
        else if (y==0)
            return Math.abs(x);
        else if ( }x==0\mathrm{ )
            return Math.abs(y);
        if(Math.abs(y)>=Math.abs(x))
            return Math.abs(x)*Math.sqrt(1.0+(y*y)/(x*x));
        else
            return Math.abs(y)*Math.sqrt(1.0+(x*x)/(y*y));
    }
    public Complex sqrt()
    {\mp@code{L}
        double w=0;
        if ( }\textrm{x}==\textrm{y}&&y==0\mathrm{ )
            return new Complex(0.0,0.0);
        else if(Math.abs(x)>=Math.abs(y))
            w=Math.sqrt(Math.abs(x))*Math.sqrt ((1.0+Math.sqrt (1.0+(y*y)/(x*x)))/2.0);
        else
            w=Math.sqrt(Math.abs(y))*Math.sqrt((Math.abs(x/y)+Math.sqrt(1.0+(x*x)/(y*y)
        if ( }x>=0\mathrm{ )
            return new Complex(w,y/(2*w));
        else if (y>=0)
            return new Complex(Math.abs(y)/(2*w),w);
        else
            return new Complex(Math.abs(y)/(2*w),-w);
    }
    public Complex power(double a)
    {
        double theta=this.argument();
        double r=this.modulus();
        Complex temp = new Complex(Math.cos(theta*a),Math.sin(theta*a))
                            .multiply(Math.pow(r,a));
        return temp;
    }
    public Complex ln()
    pub
```

```
    double theta=this.argument();
    double r=this.modulus();
    return new Complex(Math.log(r),theta);
}
public Complex exp()
{
    double etothex=Math.exp(x);
    return new Complex(etothex*Math.cos(y),etothex*Math.sin(y));
}
}
    double[] points = gjw.getPoints();
    double[] weights = gjw.getWeights();
    double m= (b+a)/2;
    double c= b-m;
    double sum = 0;
    {
    for(int i=0;i<weights.length;i++)
    sum+=f.value(c*points[i]+m)*weights[i];
    sum=sum*Math.pow(c,alpha+beta+1);
    return sum;
}
}
```


## Code for class ForwardGaussQuad

## Code for class RealNewtonRaphson

```
public class ForwardGaussQuad
{
```

public class RealNewtonRaphson
\{
public Complex recurse(int $n$, Complex a Complex b, double al,
\{
ddeffiveetise)double[] solvematrix(double[][] A, double[] b)
\{
if (a.imag ()-b.imag ()==0\&\&a.real()-b.real ()==0)
return new Complex (0,0);
int $\mathrm{l}=\mathrm{n}-3$;
double[][] L = new double[1][1];
Complex half = a.add(b).divide(2.0);
double[] $\mathrm{q}=$ new double[1];


return integrateSubinterval(N, a, b, al, be);
\}
$-[p][p]=1$;
for (int $p=0 ; p<1 ; p^{++}$)
for (int $r=p+1 ; r<1 ; r++$ )
public Complex integrateSubinterval(int N, Complex a, Complex b, double alpha, double beta)
\{
GaussJacobiWeights gjw = new GaussJacobiWeights(0.0,beta,N); \{
for (int $c=p+1 ; c<1 ; c++$ )
double[] points = gjw.getPoints();
$\mathrm{A}[\mathrm{r}][\mathrm{p}]=0$;
double [] weights $=$ gjw.getWeights ()$;$
double $m=$ (b.real() + a.real()) $/ 2$;
$\mathrm{A}[\mathrm{r}][\mathrm{c}]=\mathrm{A}[\mathrm{r}][\mathrm{c}]-\mathrm{L}[\mathrm{r}][\mathrm{p}] * \mathrm{~A}[\mathrm{p}][\mathrm{c}]$;
$\}^{A}$
double $m=$ (b.real() + a.real()) $/ 2$;
double $c=a . r e a l()-m ;$
\}
double bimag $=\mathrm{b}$. imag () ;
for (int $r=0 ; r<1 ; r++$ )
double bimag $=$ b.imag ()$;$
Complex sum $=$ new Complex ( $0.0,0.0)$;
\{
Complex sum $=$ new Complex ( $0.0,0.0$ ) ;
for (int $i=0 ; i<$ weights.length;i++)
Complex sum $=$ new Complex (0.0,0.0);
for (int $i=0 ; i<w e i g h t s . l e n g t h ; i++) ~$$\quad$ double sum=0;

sum=sum.multiply ((new Complex (c,0)). power(beta+1.0));
sum $+=\mathrm{L}[\mathrm{r}][\mathrm{c}] * \mathrm{q}[\mathrm{c}]$;
sum=sum.multiply((new Complex(c,0)). power(
gjw $=$ new GaussJacobiWeights(alpha, $0.0, \mathrm{~N}$ );
$\mathrm{q}[\mathrm{r}]=(\mathrm{b}[\mathrm{r}]-$ sum $) ;$
gjw $=$ new GaussJacobiWeigh
points $=$ gjw.getPoints();
\}
points $=$ gjw.getPoints(); \}
weights $=$ gjw.getWeights () ; $\quad$ for (int $r=0 ; r<1 ; r++$ )
weights $=$ gjw. getWeights
$\mathrm{m}=(\mathrm{b} \cdot \operatorname{imag}()+\mathrm{a} \cdot \operatorname{imag}(\mathrm{)}) / 2 ;$
\{
$\mathrm{m}=(\mathrm{b} \cdot \operatorname{imag}(\mathrm{)}+\mathrm{a}$
$\mathrm{c}=\mathrm{a} \cdot \mathrm{imag}(\mathrm{m}) \mathrm{m} ;$
double sum=0;
$\mathrm{c}=\mathrm{a} . \mathrm{imag}^{()-m}$;
double areal $=$ a.real () ;
double sum=0;
for (int $c=1-r ; c<1 ; c++$ )
Complex sum2 = new Complex (0.0,0.0) ;
sum $+=A[1-r-1][c] * x[c]$;
$x[1-r-1]=(q[1-r-1]-$ sum $) / A[1-r-1][1-r-1] ;$
for (int $i=0 ; i<w e i g h t s . l e n g t h ; i++$ )
$\quad$ if (alpha! $=0.0)$
\}
(int $i=0 ; i<w e i g h t s . l e n g t h ; i++)$
if (alpha! $=0.0$ )

sum2=sum2.add(value(new Complex(areal, c*points[i]+m), $-1,-1$ ) pr.m
sum2=sum2.add(value(new Complex(areal, c*points[i]+m), -1, -1) p.
sum2=sum2.multiply(new Complex $(0, c) \cdot \operatorname{power}($ alpha+1.0));
if (a.imag ()==0)
sum2.imag (-sum2.imag());
\{
double[][] jac=new double[n][n] ;
for (int $j=0 ; j<(n-3) ; j++$ )
return sum.add(sum2);
\}
\{
double temp $=x[j+2]$;
double $\mathrm{h}=$ TOL*Math.abs (temp);
${ }^{\cdots}$
if ( $h==0$ )
if ( $h==0$ )
$h=T O L ;$
$x[j+2]=$ temp $+h$; //Reduces floating-point error

## Code for class GaussQuad

    \(\mathrm{h}=\mathrm{x}[\mathrm{j}+2]\)-temp;
    \(\mathrm{h}=\mathrm{x}[\mathrm{j}+2]\)-temp;
    double [] funcvalues2=function( x$)$;
public class GaussQuad
$x[j+2]=$ temp;
\{
for (int $i=0 ; i<(n-3) ; i++)$
$j a c[i][j]=(f u n c v a l u e s 2[i]-f u n c v a l u e s[i]) / h ;$
... $\quad$ public double integrate(int $N$ )
\}
public double integrate(int N) \}
\{ GaussJacobiWeights gjw = new GaussJacobiWeights(alpha,beta,N); \}

```
    private double[] function(double[] x)
    pr
        double[] f = new double[n-3];
        SchwarzFunction denom = new SchwarzFunction(x,angle,0);
        GaussQuad gq = new GaussQuad(denom, denom.alpha(), denom.beta(), denom.a(), denom.b());
        double den = Math.abs(gq.integrate(GQN));
        double den1 = vertex[1].subtract(vertex[0]).modulus();
        for(int i=0;i<(n-3);i++)
        {
            SchwarzFunction numer = new SchwarzFunction(x,angle,i+1);
            gq = new GaussQuad(numer, numer.alpha(), numer.beta(), numer.a(), numer.b());
            double num = Math.abs(gq.integrate(GQN));
            double num1 = vertex[i+2].subtract(vertex[i+1]).modulus();
            f[i]=num/den-num1/den1;
        }
        return f;
    }
;
```


## References

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