

Homework Problem Sheet 3

Problem 3.1 Fourier Spectral Galerkin Scheme for Two-Point Boundary Value Problem

In [NPDE, Section 1.5.2.1] you learned about the discretization of 2-point boundary value problems based on global polynomials using integrated Legendre polynomials as basis. The implementation for a linear BVP was presented in [NPDE, § 1.5.48].

This problem is focused on another variant of spectral Galerkin discretization, which relies on non-polynomial trial and test spaces and, again, employs globally supported basis functions. This time the solution will be approximated by linear combinations of trigonometric functions. You will be asked to implement the Galerkin discretization in MATLAB and to study its convergence in a numerical experiment, cf [NCSE, Thm. 9.1.4].

We consider the linear variational problem: seek $u \in \mathcal{C}_{0,\text{pw}}^1([0, 1])$ such that

$$\int_0^1 \sigma(x) \frac{du}{dx}(x) \frac{dv}{dx}(x) dx = \int_0^1 f(x)v(x) dx, \quad \forall v \in \mathcal{C}_{0,\text{pw}}^1([0, 1]), \quad (3.1.1)$$

cf. [NPDE, Eq. (1.4.23)]. For the discretization of (3.1.1) we may use a so-called *Fourier-spectral Galerkin method*, which boils down to a Galerkin method using the trial and test space

$$V_{N,0} := \text{span}\{\sin(\pi x), \sin(2\pi x), \dots, \sin(N\pi x)\}, \quad (3.1.2)$$

and the basis function already given in the definition (3.1.2). It is related the spectral Galerkin scheme discussed in [NPDE, Section 1.5.2.1].

(3.1a) Show that the functions specified in (3.1.2) really provides a basis of $V_{N,0}$.

HINT: First establish orthogonality of the basis functions with respect to the inner product

$$(f, g)_{L^2(0,1)} := \int_0^1 f(x)g(x) dx.$$

Then recall that a basis of a finite dimensional vector space is a maximal *linearly independent* subset.

Solution: We only have to show that the sines are linearly independent. In fact, they are orthogonal. For $i \neq j$:

$$\int_0^1 \sin(i\pi x) \sin(j\pi x) dx = \frac{1}{2} \int_0^1 [\cos((i-j)\pi x) - \cos((i+j)\pi x)] dx = 0,$$

as both $i+j$ and $i-j$ are nonzero.

(3.1b) Which basis should be used for the Fourier spectral scheme, if we had to approximate functions in the space $\mathcal{C}_{0,\text{pw}}^1([a, b])$ for fixed $a < b$, instead of the space $\mathcal{C}_{0,\text{pw}}^1([0, 1])$.

HINT: Read [NPDE, § 1.5.41].

Solution: The transformed basis in this case would be

$$V_{N,0}([a, b]) = \text{span}\{\sin(\pi(x - a)/h), \sin(2\pi(x - a)/h), \dots, \sin(N\pi(x - a)/h)\},$$

with $h = b - a$.

(3.1c) Since (3.1.1) is a linear variational problem, any Galerkin discretization will lead to a linear system of equations. For the case $\sigma \equiv 1$ compute its matrix for the Galerkin scheme relying on $V_{N,0}$ and the trigonometric basis specified in (3.1.2). Discuss structural properties of the matrix like sparsity, symmetry, regularity.

Solution: We need to compute the matrix entries

$$\mathbf{A}_{i,j} = ij\pi^2 \int_0^1 \cos(i\pi x) \cos(j\pi x) dx.$$

Using the identity

$$\cos(i\pi x) \cos(j\pi x) = \frac{1}{2}[\cos((i - j)\pi x) + \cos((i + j)\pi x)],$$

we see as in (3.1a) that the off-diagonal terms are all zero. Then

$$\mathbf{A}_{i,i} = \frac{i^2\pi^2}{2} \int_0^1 (1 + \cos(2i\pi x)) dx = \frac{i^2\pi^2}{2}.$$

So

$$\mathbf{A} = \frac{\pi^2}{2} \begin{pmatrix} 1 & & & & \\ & 4 & & & \\ & & 9 & & \\ & & & \ddots & \\ & & & & N^2 \end{pmatrix}.$$

This matrix is diagonal, thus both sparse and symmetric.

(3.1d) Write a MATLAB function

```
function A = getGalMat(sigma,N)
```

that computes the Galerkin matrix for the Fourier spectral Galerkin scheme for (3.1.1). Here `sigma` is a handle to the coefficient function σ . The evaluation of the integrals should be done by means of a $3N$ -point Gaussian quadrature formula on $[0, 1]$.

HINT: The nodes and weights of the Gaussian quadrature rules on $[a, b]$ can be computed by the MATLAB function `[x,w] = gauleg(a,b,n,tol)`, which is available for download.

Solution: See Listing 3.1.

Listing 3.1: Computation of the Galerkin matrix

```

1 function C = getGalMat(sigma, N)
2
3     [x,w] = gauleg(0,1,3*N);
4
5     C = cos(pi*x*(1:N));
6     % C(k,i) = cos(pi i x_k)
7
8     D = bsxfun(@times, C, sigma(x));
9     D = bsxfun(@times, D, w);
10    % D(k,i) = cos(pi i x_k) sigma(x_k) w_k
11
12    C = C' * D;
13    % C(i,j) = sum_k cos(pi i x_k) cos(pi j x_k) sigma(x_k)
14                w_k
15
16    C = bsxfun(@times, C, 1:N);
17    C = bsxfun(@times, C, (1:N)');
18    C = pi^2 * C;
19 end

```

(3.1e) Write a function

```
function phi = getrhsvector(f,N)
```

that computes the right-hand side vector for the Fourier spectral Galerkin discretization with N basis functions. The routine should rely on $3N$ -point Gaussian quadrature for the evaluation of the integrals.

Solution: See [Listing 3.2](#).

Listing 3.2: Computation of the right-hand side vector

```

1 function phi = getrhsvector(f, N)
2
3     [x,w] = gauleg(0,1,3*N);
4
5     C = sin(pi*x*(1:N));
6     C = bsxfun(@times, C, f(x));
7     phi = C' * w;
8
9 end

```

(3.1f) For $\sigma(x) = \frac{1}{\cosh(\sin(\pi x))}$, $f(x) = \pi^2 \sin(\pi x)$, determine an approximate solution of (3.1.1) by means of the Fourier spectral scheme introduced above. Create a suitable plot of the L^2 -norm and L^∞ -norm of the discretization error versus the number N of unknowns for $N = 2, 3, \dots, 14$ and, thus, investigate the convergence of the method [[NCSE](#), Remark 9.1.4].

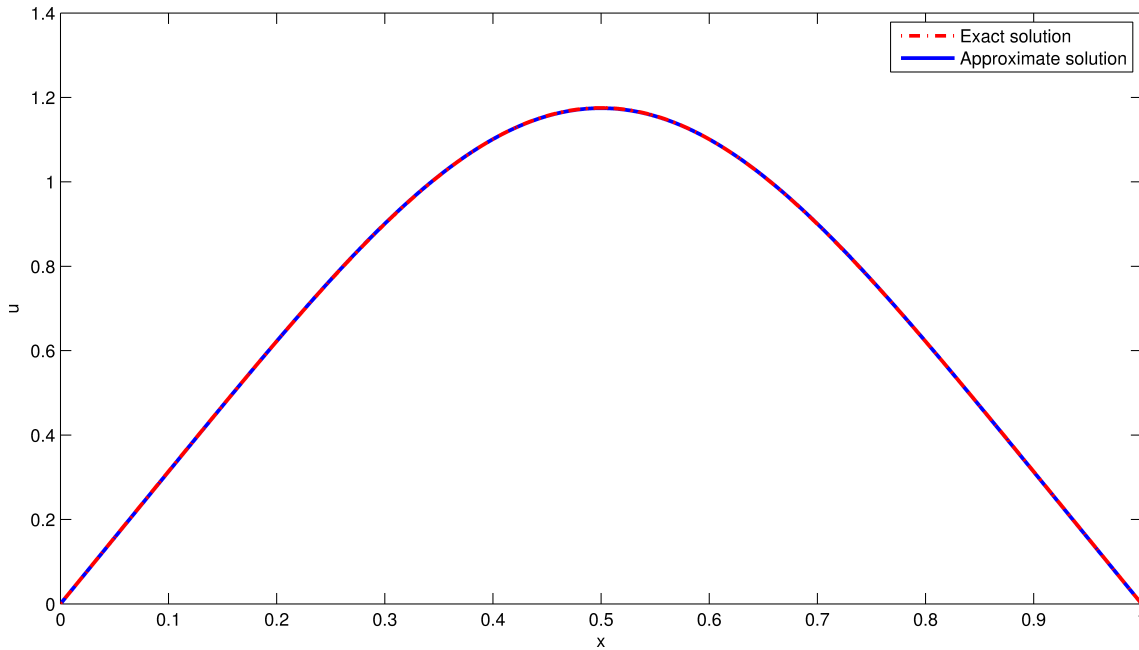


Figure 3.1: Exact and approximate solution for $N = 3$.

The computation of the norms should be done approximately by means of numerical quadrature (equidistant trapezoidal rule with 10^5 points) and sampling (in 10^5 equidistant points), respectively, see [NPDE, Rem. 1.6.19].

HINT: The exact solution of the 2-point boundary value problem is

$$u(x) = \sinh(\sin(\pi x)).$$

Solution:

See Figure 3.1 for the approximate and exact solution at $N = 3$, Figure 3.2 for the convergence plots and Listing 3.3 for the code used.

To compute the $L^2([a, b])$ -error (here $[a, b] = [0, 1]$) we use a composite trapezoidal rule with $N + 1$ equidistant points x_0, \dots, x_N

$$\int_b^a f(x) \, dx \approx \frac{b-a}{N} \left(\frac{f(x_0)}{2} + \sum_{i=1}^{N-1} f(x_i) + \frac{f(x_N)}{2} \right) = \frac{b-a}{N} \sum_{i=1}^N f(x_i),$$

where in the last step we used the periodicity of f , i.e., $f(x_0) = f(a) = f(b) = f(x_N)$.

To estimate the rate of the exponential convergence, we do the following consideration for the error e

$$e \approx e^{-\gamma n^\delta} \iff \log e \approx -\gamma n^\delta \iff \log \log e \approx \log(-\gamma) + \log(n^\delta) = \delta \log n + \log -\gamma$$

Then we do a polyfit of the last term and obtain the best matching polynomial in $x = \log n$ and obtain $\delta x + x_0$, and we recompute γ as $\gamma = -e^{x_0}$. We implement this in Listing 3.3 and obtain a similar convergence for both norms

L-inf convergence exponential with gamma: 1.4904 and delta: 1.2291

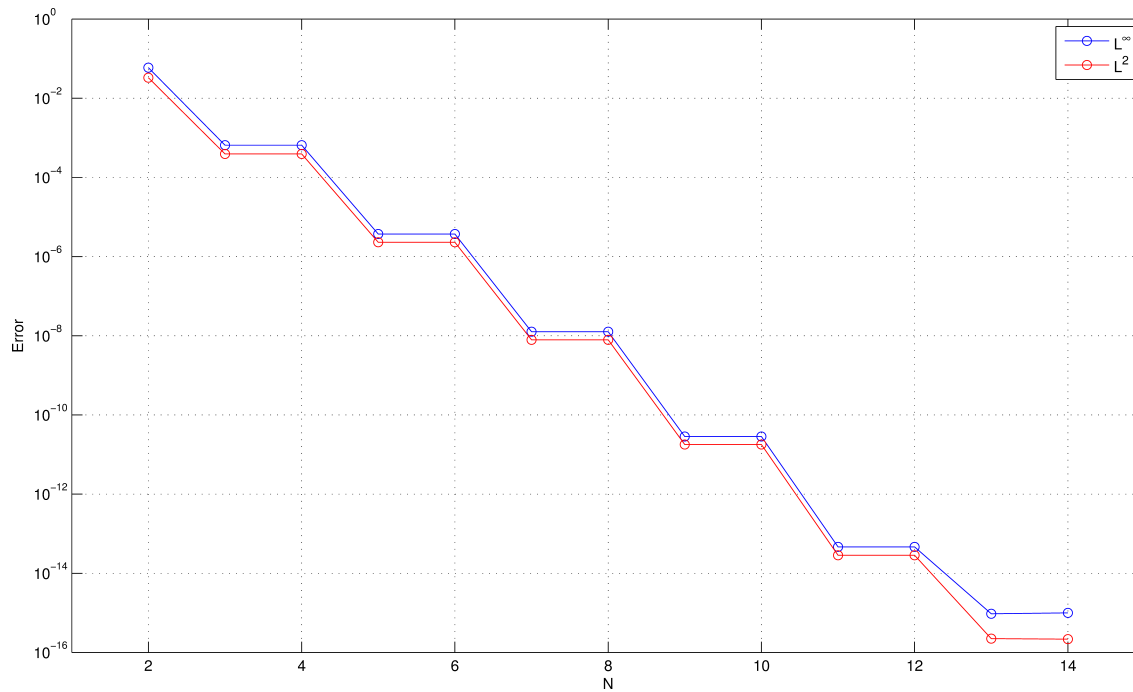


Figure 3.2: Convergence for subproblem (3.1f).

L^2 convergence exponential with gamma: 1.7698 and delta: 1.1627. Note that the parameter $0 < \delta < \infty$ (high values are good) is more important than γ (which is often omitted). The plots looks like a *staircase*, this comes from the fact, that this particular exact solution uses only every second function in the space $V_{N,0}$.

Listing 3.3: Convergence estimates for (3.1f)

```

1 nvals = 2:14;
2 M = 10^5;
3 sigma = @(x) ones(size(x))./cosh(sin(pi*x));
4 f = @(x) pi^2*sin(pi*x);
5 exact = @(x) sinh(sin(pi*x));
6
7 xpts = linspace(0,1,M+1)';
8 xpts = xpts(2:end-1);
9
10 l2err = [];
11 lierr = [];
12
13 for N = nvals
14     A = getGalMat(sigma, N);
15     L = getrhsvector(f, N);
16     mu = A \ L;
17
18     approx = evaltrigsum(mu, M);
19     err = abs(approx - exact(xpts));
20

```

```

21     l2err = [l2err; sqrt(sum(err.^2)/M)];
22     lierr = [lierr; max(err)];
23 end
24
25 semilogy(nvals, lierr, 'bo-');
26 hold on; grid on;
27 plot(nvals, l2err, 'ro-');
28 xlim([1, 15]);
29 xlabel('N');
30 ylabel('Error');
31 legend('L^\infty', 'L^2');
32
33 P = polyfit(log(nvals), log(log(lierr')), 1);
34 gamma_inf=-real(exp(P(2)));
35 delta_inf=P(1);
36
37 disp(['L-inf convergence exponential with gamma: '
       num2str(gamma_inf) ' and delta: ' num2str(delta_inf) ]);
38 Q = polyfit(log(nvals), log(log(l2err')), 1);
39 gamma_2=-real(exp(Q(2)));
40 delta_2=Q(1);
41 disp(['L-2 convergence exponential with gamma: '
       num2str(gamma_2) ' and delta: ' num2str(delta_2) ]);
42
43 plot(nvals, exp(-gamma_inf.*nvals.^delta_inf));
44 hold off;

```

(3.1g) Carry out the investigations requested in [subproblem \(3.1f\)](#) for

$$\sigma(x) = \begin{cases} 2 & \text{for } |x - \frac{1}{2}| < \frac{1}{4}, \\ 1 & \text{elsewhere,} \end{cases} \quad f \equiv 1,$$

this time using $N = 2, 4, 8, 16, 32, 64, 128, 256, 512, 1024$. What kind of convergence do you observe? Relate with the observation made in [subproblem \(3.1f\)](#) and try to explain.

HINT: The exact solution is

$$u(x) = \begin{cases} \frac{3}{64} + \frac{1}{4}x - \frac{1}{4}x^2 & |x - \frac{1}{2}| < \frac{1}{4}, \\ \frac{1}{2}x - \frac{1}{2}x^2 & \text{elsewhere.} \end{cases}$$

Also recall the observations made in [\[NPDE, Exp. 1.6.31\]](#).

Solution:

See [Figure 3.3](#) for the convergence plots and [Listing 3.4](#) for the code used. The convergence is algebraic with rates 1 for the L^∞ -norm and 1.25 for the L^2 -norm.

Listing 3.4: Convergence estimates for [subproblem \(3.1f\)](#)

```

1 nvals = 2.^(1:10);

```

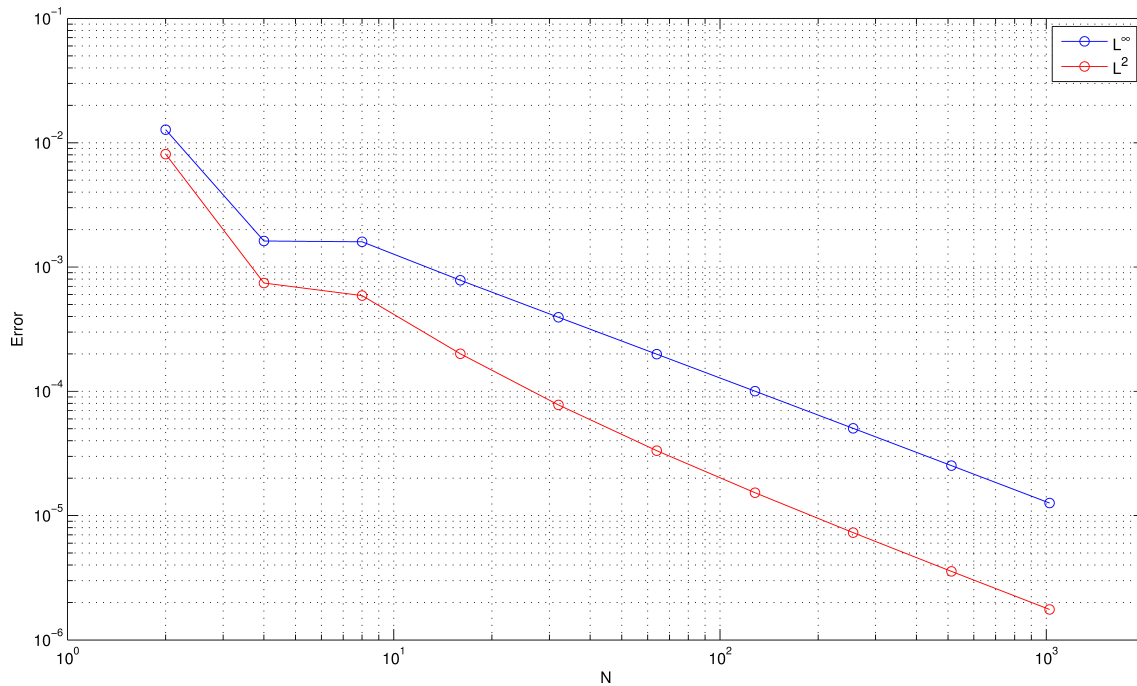


Figure 3.3: Convergence for subproblem (3.1f).

```

2 M = 10^5;
3 inside = @(x) abs(x-0.5) < 0.25;
4 outside = @(x) 1 - inside(x);
5 sigma = @(x) (1+inside(x)).*ones(size(x));
6 f = @(x) ones(size(x));
7 temp = @(x) 0.25*x - 0.25*x.^2;
8 exact = @(x) (1+outside(x)).*temp(x) +
    inside(x).*ones(size(x))*3/64;
9
10 xpts = linspace(0,1,M+1)';
11 xpts = xpts(2:end-1);
12
13 l2err = [];
14 lierr = [];
15
16 for N = nvals
17     A = getGalMat(sigma, N);
18     L = getrhsvector(f, N);
19     mu = A \ L;
20
21     approx = [evaltrigsum(mu, M)];
22     err = abs(approx - exact(xpts));
23
24     l2err = [l2err; sqrt(sum(err.^2)/M)];
25     lierr = [lierr; max(err)];
26 end

```

```

27
28 loglog(nvals, lierr, 'bo-');
29 hold on; grid on;
30 plot(nvals, l2err, 'ro-');
31 xlim([1, 2000]);
32 xlabel('N');
33 ylabel('Error');
34 legend('L^\infty', 'L^2');
35
36 P = polyfit(log(nvals), log(lierr'), 1);
37 disp(['L-inf convergence algebraic with rate: '
      num2str(-P(1))]);
38 P = polyfit(log(nvals), log(l2err'), 1);
39 disp(['L-2 convergence algebraic with rate: '
      num2str(-P(1))]);

```

Listing 3.5: Testcalls for [Problem 3.1](#)

```

1 M = 6;
2
3 sigma = @(x) (1./(cosh(sin(pi*x))));
4 N=5;
5 fprintf(' \n\n##getGalMat:')
6 getGalMat(sigma,N)
7
8 f = @(x) (pi^2 * sin(pi*x));
9 fprintf(' \n\n##getrhsvector:')
10 getrhsvector(f,N)

```

Listing 3.6: Output for Testcalls for [Problem 3.1](#)

```

1 test_call12
2
3 ##getGalMat:
4 ans =
5
6      4.4209      0.0000      1.4066      0.0000      0.2058
7      0.0000     16.1117      0.0000      3.4734     -0.0000
8      1.4066      0.0000     35.9390      0.0000      6.4682
9      0.0000      3.4734      0.0000     63.8443     -0.0000
10     0.2058     -0.0000      6.4682     -0.0000     99.7504
11
12 ##getrhsvector:
13 ans =
14
15      4.9348
16      0.0000
17      0.0000
18     -0.0000
19      0.0000

```


Problem 3.2 Linear Finite Elements for the Brachistochrone Problem

In this problem we focus on the Galerkin discretization of the variational formulation for the brachistochrone problem by means of linear finite elements as introduced in [NPDE, Section 1.5.2.2].

We remind that the variational problem reads as: Find $\mathbf{u} \in V$ so that

$$\int_0^1 \left(\frac{\mathbf{u}' \cdot \mathbf{v}'}{\sqrt{-u_2} \|\mathbf{u}'\|} - \frac{v_2 \|\mathbf{u}'\|}{2\sqrt{-u_2} u_2} \right) d\xi = 0 \quad \text{for all } \mathbf{v} \in V_0, \quad (3.2.1)$$

$V_0 := \{ \mathbf{v} \in (C_{\text{pw}}^1([0, 1]))^2 \mid \mathbf{v}(0) = \mathbf{v}(1) = \mathbf{0} \}$ and $V := \{ \mathbf{v} \in (C_{\text{pw}}^1([0, 1]))^2 \mid \mathbf{v}(0) = \mathbf{a}, \mathbf{v}(1) = \mathbf{b} \}$. Throughout we use equidistant meshes $\mathcal{M} = \{]x_{j-1}, x_j[: x_{j-1} := \frac{j-1}{M}, x_j := \frac{j}{M}, j = 1, \dots, M \}$ of $[0, 1]$ and the standard “tent function” basis \mathfrak{B} of the Galerkin trial space

$$V_{N,0} = (\mathcal{S}_{1,0}^0(\mathcal{M}))^2 = \left\{ \mathbf{v} \in (C^0([0, 1]))^2 : \mathbf{v}|_{[x_{i-1}, x_i]} \text{ linear}, \right. \\ \left. i = 1, \dots, M, \mathbf{v}(0) = \mathbf{v}(1) = \mathbf{0} \right\}. \quad (3.2.2)$$

This is explained in detail in [NPDE, § 1.5.82], see, in particular, [NPDE, Eq. (1.5.83)]. It is recommended to use the ordering of the basis functions implied by [NPDE, Eq. (1.5.83)], though you are free to use any other scheme.

(3.2a) Determine a “ \mathbf{u} -dependent coefficient function” $\sigma(\xi) = \sigma(\mathbf{u})(\xi)$ and a “ \mathbf{u} -dependent source function” $\mathbf{f}(\xi) = \mathbf{f}(\mathbf{u})(\xi)$ such that the variational formulation (3.2.1) can be written as

$$\mathbf{u} \in V : \quad \int_0^1 \sigma(\mathbf{u})(\xi) \mathbf{u}'(\xi) \cdot \mathbf{v}'(\xi) d\xi = \int_0^1 \mathbf{f}(\mathbf{u})(\xi) \cdot \mathbf{v}(\xi) d\xi \quad \forall \mathbf{v} \in V_0. \quad (3.2.3)$$

The point of recasting (3.2.1) in this form is the reduction to the structure of a linear variational problem. For the elastic string model this has proved highly useful as regards the implementation of Galerkin discretizations in [NPDE, § 1.5.55], Code [NPDE, Code 1.5.58], and [NPDE, § 1.5.82], Code [NPDE, Code 1.5.92]. Please study the latter example and code again, in case you do not remember the rationale behind (3.2.3).

Solution: We have:

$$\sigma(\mathbf{u}) = \frac{1}{\sqrt{-u_2} \|\mathbf{u}'\|}, \quad \mathbf{f}(\mathbf{u}) = \frac{\|\mathbf{u}'\|}{2\sqrt{-u_2} u_2} \begin{pmatrix} 0 \\ 1 \end{pmatrix}.$$

(3.2b) Implement a MATLAB function

```
s = sigma(mu, xi)
```

where:

- `mu` is a $2 \times (M + 1)$ matrix containing the components $\{\mu_1, \mu_2, \dots, \mu_{M+1}\}$ of \mathbf{u}_N with respect to the basis representation of the curve (where $\mu_1 = \mathbf{u}(0)$ and $\mu_{M+1} = \mathbf{u}(1)$ are the pinning points);
- `xi` is a vector of evaluation points in the interval $]0, 1[$.

The output is a vector \mathbf{s} with the evaluations of the scalar function $\sigma = \sigma(\mathbf{u})$ at the mesh points \mathbf{x}_i .

HINT: You may use the MATLAB function `linterp` to get a piecewise linear interpolation of \mathbf{u}_N at the evaluation points.

Remember that the mesh on which you have the coefficients μ is equispaced.

The derivative of \mathbf{u}_N is piecewise constant. You don't have to worry if an evaluation point is also a mesh point, where the derivative is discontinuous; in this case, you can take either the left or the right derivative.

A reference implementation `sigma_ref` is available in the file `sigma_ref.p`.

Solution: See listing 3.7 for the code.

Listing 3.7: Implementation for `sigma`

```

1 function s = sigma(mu, xi)
2
3     mesh = linspace(0, 1, length(mu));
4     h = mesh(2) - mesh(1);
5     d = (mu(:, 2:end) - mu(:, 1:end-1)) / h; % derivative of u
6     dnorm = sqrt(d(1, :).^2 + d(2, :).^2); % norm of the
           derivative
7     mu_xi = linterp(mesh, mu(2, :), xi); % interpolate the
           second component in the evaluation points
8
9     for j=1:length(xi)
10        index = find(mesh > xi(j), 1); % find the index of the
           first mesh point greater than xi(j)
11                                           % the derivative in xi
                                           corresponds to the
                                           entry d(index-1)
12                                           % at mesh points, we
                                           consider the right
                                           derivative
13        s(j) = 1 / (sqrt(-mu_xi(j)) * dnorm(index-1));
14    end

```

(3.2c) Write a MATLAB function

$$\mathbf{f} = \text{sourcefn}(\mu, \mathbf{x}_i)$$

where the input arguments are the same as in [subproblem \(3.2b\)](#) and the output is a $2 \times K$ matrix, with K the length of \mathbf{x}_i , containing the evaluations of the \mathbf{u} -dependent source function $\mathbf{f} = \mathbf{f}(\mathbf{u})$ at the mesh points \mathbf{x}_i .

If an evaluation point coincides with a mesh point, where the derivative of \mathbf{u}_N is not uniquely defined, consider either the left or the right derivative.

Solution: See listing 3.8 for the code. operations as far as possible.

Listing 3.8: Implementation for sourcefn

```

1 function f = sourcefn(mu,xi)
2
3     M = length (mu) -1;
4     mesh = linspace (0,1, length (mu));
5     h = mesh (2)-mesh (1);
6     d = (mu(:,2:end)-mu(:,1:end-1))/h; % derivative of u
7     dnorm = sqrt (d(1,:).^2 + d(2,:).^2); % norm of the
           derivative
8     mu_xi = linterp(mesh,mu(2,:),xi);
9
10    f(1,:) = zeros (1, length (xi));
11
12    for j=1:length (xi)
13        index = find (mesh>xi(j),1); % find the index of the
           first mesh point greater than xi(j)
14                                   % the derivative in xi
                                   corresponds to the entry
                                   d(index-1)
15                                   % at mesh points, we
                                   consider the right
                                   derivative
16        f(2,j) = dnorm(index-1) / (2*sqrt (-mu_xi(j))*mu_xi(j));
17    end

```

(3.2d) Implement a MATLAB function

```
t = travelttime (mu)
```

which accepts as input the vector `mu` as in (3.2b), and returns the approximate value of the functional

$$J(\mathbf{u}_N) = \int_0^1 \frac{\|\mathbf{u}'_N(\xi)\|}{\sqrt{-(\mathbf{u}_N)_2(\xi)}} d\xi \quad (3.2.4)$$

representing the time needed to go from $\mathbf{u}(0)$ to $\mathbf{u}(1)$ along the curve \mathbf{u}_N .

For the computation of the integral in (3.2.4), use the midpoint rule (see [NPDE, Eq. (1.5.77)]). Note that the trapezoidal rule would be inappropriate because of the singularity of the functional at the origin.

HINT: A reference implementation `travelttime_ref` is available in the file `travelttime_ref.p`.

Solution: See listing 3.9 for the code.

Listing 3.9: Implementation for travelttime

```

1 function t = travelttime (mu)
2

```

```

3  mesh = linspace (0,1,length (mu)) ;
4  h = mesh (2)-mesh (1) ;
5  d = (mu (:,2:end)-mu (:,1:end-1))/h;
6  dnorm = sqrt (d (1,:) .^2 + d (2,:) .^2) ;
7  % TRAPEZOIDAL RULE
8  %     mu2 = mu (2,:) ;
9  %     t =
10     sum ( (dnorm (1:end-1)+dnorm (2:end)) ./ (sqrt (-mu2 (2:(end-1)))) ) + ...
11     ...+dnorm (end)/sqrt (-mu2 (end)) ;
12 %     if abs (mu2 (1))>eps
13 %         t = t+dnorm (1)/sqrt (-mu2 (1)) ;
14 %     end
15 % MIDPOINT RULE
16 xi = h/2:h:(1-h/2) ;
17 mu_xi = linterp (mesh,mu (2,:),xi) ;
18 t = sum (dnorm ./ (sqrt (-mu_xi))) *h;

```

(3.2e) Proceeding as in [NPDE, § 1.5.82], the discretization of (3.2.3) leads to a *nonlinear* system of equations of the form

$$\begin{pmatrix} \mathbf{R}(\vec{\mu}) & 0 \\ 0 & \mathbf{R}(\vec{\mu}) \end{pmatrix} \vec{\mu} = \begin{pmatrix} \vec{\varphi}_1(\vec{\mu}) \\ \vec{\varphi}_2(\vec{\mu}) \end{pmatrix}, \quad (3.2.5)$$

with $\mathbf{R}(\vec{\mu}) \in \mathbb{R}^{M-1,M-1}$ and $\vec{\varphi}_i(\vec{\mu}) \in \mathbb{R}^{M-1}$. Write a MATLAB function

`R = Rmat (mu)`

such that, given the coefficients $\vec{\mu} = \text{mu}$ in input (as in [subproblem \(3.2a\)](#)), returns the matrix $\mathbf{R} = \mathbf{R}(\vec{\mu})$.

For the evaluation of the integrals, use the midpoint rule [NPDE, Eq. (1.5.77)].

HINT: Compute the matrix including also the rows and columns referring to the two basis functions for the offset function. In this way, your matrix \mathbf{R} will have dimensions $(M+1) \times (M+1)$. Then, in [subproblem \(3.2g\)](#), where you will have to solve the linear system, you have to consider just the entries relative to the inner nodes (i.e. you have to exclude the first and last columns and rows of \mathbf{R}). The reason for doing this is that with such \mathbf{R} it will be easier, in [subproblem \(3.2g\)](#), to modify the right hand side to take into account the boundary conditions.

A reference implementation `R_ref` is available in the file `R_ref.p`.

Solution: See listing 3.10 for the code.

Listing 3.10: Implementation for `Rmat`

```

1  function R = Rmat (mu)
2
3      mesh = linspace (0,1,length (mu)) ;
4      h = mesh (2)-mesh (1) ;
5      xi = h/2:h:(1-h/2) ; % midpoints, where we will evaluate
                           sigma

```

```

6   M = length(mu)-1;
7   % Computation of the r_j
8   s = sigma(mu,xi);
9   r = s./h;
10
11  % Assemble tridiagonal matrix R
12  R = gallery('tridiag', [-r(1), -r(2:M-1), -r(M)], [r(1),
    r(1:M-1)+r(2:M), r(M)], [-r(1), -r(2:M-1), -r(M)]);

```

(3.2f) Write a MATLAB function

```
phi = rhs(mu)
```

which, given the coefficients `mu` in input, returns as output the right hand side vector from (3.2.5)

$$\vec{\varphi}(\vec{\mu}) = \begin{pmatrix} \vec{\varphi}_1(\vec{\mu}) \\ \vec{\varphi}_2(\vec{\mu}) \end{pmatrix} \in \mathbb{R}^{2M-2}. \quad (3.2.6)$$

For integration, consider the composite trapezoidal quadrature rule [NPDE, Eq. (1.5.72)]. At the origin (where the source function is singular), consider the integrand to be zero.

HINT: A reference implementation `rhs_ref` is available in the file `rhs_ref.p`.

Solution: See listing 3.11 for the code.

Listing 3.11: Implementation for `rhs`

```

1  function phi = rhs(mu)
2
3      mesh = linspace(0,1,length(mu));
4      h = mesh(2)-mesh(1);
5      % TRAPEZOIDAL RULE
6      xxi = h:h:(1-h);
7      phi =
          (sourcefn(mu,xxi(1:end))+sourcefn(mu,xxi(1:end)))*h/2;
8  % MIDPOINT RULE: DOESN'T WORK!
9  % xi = h/2:h:(1-h/2);
10 % sourcefn(mu,xxi); % for debugging
11 % 2*sourcefn(mu,xxi) % for debugging
12 % source = sourcefn(mu,xi);
13 % source(:,1:end-1) + source(:,2:end) % for debugging
14 % phi = source(:,1:end-1)*h/2 + source(:,2:end)*h/2;

```

(3.2g) The nonlinear system (3.2.6) can be solved by *fixed point iteration*. In this way, an approximate solution is computed solving, at each iteration, a *linear* system of equations (see [NPDE, § 1.5.82], [NPDE, Code 1.5.92]).

Implement a MATLAB function

```
mufinal = solvebrachlin(mu0, tol)
```

to compute the approximate solution (i.e. the coefficients `mu`final with respect to the basis functions) using the fixed point iteration.

The input arguments are the initial guess `mu0` and the tolerance `tol` for the relative error in the fixed point iteration algorithm (see [NPDE, Code 1.5.92], line 41).

For each iteration, plot the shape of the curve (see [NPDE, Code 1.5.92], lines 19-22).

Plot the travel time as defined in [subproblem \(3.2d\)](#) with respect to the number of iteration steps.

Remark: Remember to take into account the boundary conditions.

HINT: The solution should look like the cycloid

$$\mathbf{u}(\xi) = \begin{pmatrix} \pi\xi - \sin(\pi\xi) \\ \cos(\pi\xi) - 1 \end{pmatrix}, \quad 0 \leq \xi \leq 1, \quad (3.2.7)$$

that was shown in the previous assignment to be a strong solution.

A reference implementation `solvebrachlin_ref` is available in the file `solvebrachlin_ref.p`.

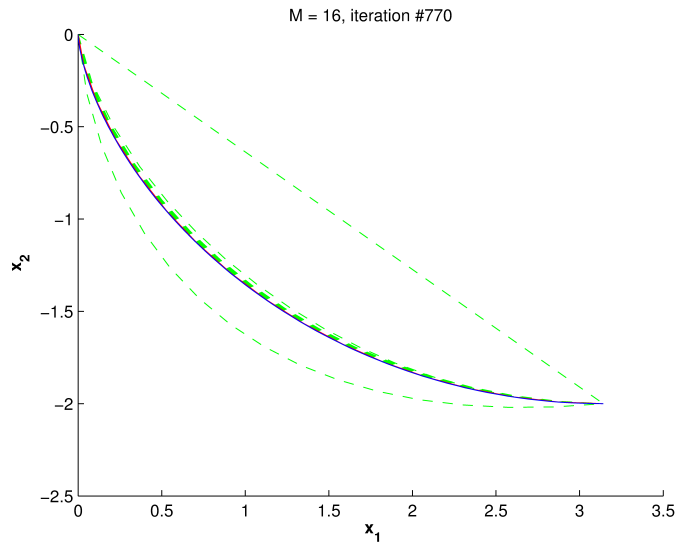
Solution: See listing 3.12 for the code.

Listing 3.12: Implementation for `solvebrachlin`

```

1 function [mu_new,figsol] = solvebrachlin(init_guess,tol)
2 % M intervals, M-1 interior nodes
3 M = length(init_guess)-1;
4 h= 1/M; %meshwidth
5
6 figsol = figure; hold on;
7 maxiter=10000;
8 mu_new = init_guess;
9 t = [];
10 for k=1:maxiter
11     mu = mu_new;
12     % Plot shape of string
13     plot(mu(1,:), mu(2,:), '--g'); drawnow;
14     title(sprintf('M = %d, iteration #%d', M,k));
15     xlabel('{\bf x_1}'); ylabel('{\bf x_2}');
16     t = [t; traveltime(mu)];
17
18     R = Rmat(mu);
19     % Computation of right hand side
20     phi = rhs(mu);
21     phi1 = phi(1,:);
22     phi2 = phi(2,:);
23
24     % Modify right hand side to take in account to the offset
25     function
26     phi1(1)= phi1(1) - R(2,1)*mu(1,1);
27     phi1(M-1) = phi1(M-1) -R(M,M+1)*mu(1,end);
28     phi2(1) = phi2(1) - R(2,1)*mu(2,1);
29     phi2(M-1) = phi2(M-1) -R(M,M+1)*mu(2,end);

```



```

30 % Solve linear system and compute new iterate
31 mu_new = [mu(:,1), [(R((2:end-1), (2:end-1))\phi1')';
32               (R((2:end-1), (2:end-1))\phi2')'], mu(:,end)];
33 % Check simple termination criterion for fixed point
    iteration.
34 if (norm(mu_new - mu, 'fro') < tol*norm(mu_new, 'fro')/M)
35     plot(mu_new(1,:), mu_new(2,:), 'r');
36     xi = linspace(0,1);
37     plot(pi*xi-sin(pi*xi), cos(pi*xi)-1, 'b');
38     break;
39 end
40
41 end
42 figure
43 plot(t);

```

The solution for $M=16$, with straight line as initial guess, should look like the following:

(3.2h) The fixed point iteration algorithm converges quite slowly to the approximate solution. To improve this, we use *nested iterations*:

- start from an initial guess μ_0 and an initial *coarse* mesh mesh_0 and compute the solution μ_1 ;
- consider the mesh μ_1 obtained from μ_0 inserting the midpoints of the interval as mesh points (thus doubling the number of mesh points) and compute the solution μ_2 considering μ_1 as initial guess;
- repeat point b) iteratively until the desired final meshwidth level L is reached.

In point b), to get the initial guess, one has to extend a piecewise linear function defined on a coarser grid to a finer nested grid. To achieve this, piecewise linear interpolation can be used.

Remark: The advantage of considering the relative error tolerance for the fixed point iteration adapted to the meshsize (see [NPDE, Code 1.5.92], line 41) is that in all iterations from point b) the same tolerance `tol` can be used.

Write a MATLAB function

```
mufinal = nestitbrachlin(uend,L,tol)
```

to solve the brachistochrone problem using nested iterations.

Here, `uend` is the right pinning point, while the left pinning point is consider to be the origin.

`L` is the number of refinement levels. Start from a mesh of $M=2$ intervals, corresponding to the level $L=0$.

HINT: For the linear interpolation for the initial guess in point b), use the MATLAB function `linterp`.

Solution: See listing 3.13 for the code.

Listing 3.13: Implementation for `nestitbrachlin`

```
1 function mufinal = nestitbrachlin(uend,L,tol)
2
3 u0 = [0;0];
4 mu = [u0, (u0+uend)/2, uend];
5 level=0;
6
7 while (level<L+1)
8     mu_new = solvebrachlin(mu,tol);
9     meshsize = 1/(2^(level+1));
10    level = level+1;
11    clear mu;
12    mu(1,:) =
13        linterp(0:meshsize:1,mu_new(1,:),0:meshsize/2:1);
14    mu(2,:) =
15        linterp(0:meshsize:1,mu_new(2,:),0:meshsize/2:1);
16 end
17 mufinal(1,:) = mu(1,:);
18 mufinal(2,:) = mu(2,:);
```

Listing 3.14: Testcalls for Problem 3.2

```
1 u0 = [0;0];
2 u1 = [pi;-2];
3 M = 4;
4 mesh = linspace(0,1,M+1);
5 h = mesh(2)-mesh(1);
6 mu = u0*(1-(0:1/M:1))+u1*(0:1/M:1);
7
8 fprintf('\n\n##sigma:')
9 xi = h/2:h:(1-h/2);
10 sigma(mu, xi)
```



```

11
12 fprintf('\n\n##sourcefn:')
13 xxi = h:h:(1-h);
14 sourcefn(mu,xxi)
15
16 fprintf('\n\n#traveltime:')
17 traveltime(mu)
18
19 fprintf('\n\n##Rmat:')
20 Rmat(mu)
21
22 fprintf('\n\n##rhs:')
23 rhs(mu)
24
25 fprintf('\n\n##solvebrachlin:')
26 solvebrachlin(mu,10^(-5))
27
28 fprintf('\n\n##nestitbrachlin:')
29 nestitbrachlin([pi;-2],3,10^(-5))

```

Listing 3.15: Output for Testcalls for [Problem 3.2](#)

```

1 >> test_call_fem
2
3 ##sigma:
4 ans =
5
6     0.5370     0.3101     0.2402     0.2030
7
8 ##sourcefn:
9 ans =
10
11         0         0         0
12    -5.2668    -1.8621    -1.0136
13
14 #traveltime:
15 ans =
16
17     4.4737
18
19 ##Rmat:
20 ans =
21
22    (1,1)     2.1481
23    (2,1)    -2.1481
24    (1,2)    -2.1481
25    (2,2)     3.3883
26    (3,2)    -1.2402
27    (2,3)    -1.2402
28    (3,3)     2.2009

```

```

29      (4,3)      -0.9607
30      (3,4)      -0.9607
31      (4,4)       1.7726
32      (5,4)      -0.8119
33      (4,5)      -0.8119
34      (5,5)       0.8119
35
36  ##rhs:
37  ans =
38
39      0          0          0
40     -1.3167    -0.4655    -0.2534
41
42  ##solvebrachlin:
43  ans =
44
45      0      0.3114      1.0169      1.9977      3.1416
46      0     -0.6794     -1.3570     -1.8269     -2.0000
47
48  ##nestitbrachlin:
49  ans =
50
51  Columns 1 through 9
52
53      0      0.0510      0.1019      0.2179      0.3339      0.4994
54      0.6650      0.8683      1.0717
55      0     -0.1679     -0.3358     -0.5287     -0.7215     -0.9031
56      -1.0847     -1.2425     -1.4002
57
58  Columns 10 through 17
59
60      1.3040      1.5364      1.7906      2.0447      2.3145      2.5843
61      2.8629      3.1416
62     -1.5280     -1.6558     -1.7500     -1.8442     -1.9022     -1.9602
63     -1.9801     -2.0000

```

Problem 3.3 Spline Collocation Method

[NPDE, Section 1.5.3.2] introduces the spline collocation method for 2-point boundary value problems based on natural cubic splines, see [NPDE, Def. 1.5.116]. In this problem we consider the simple BVP

$$-\frac{d^2u}{dx^2} = f(x) \quad \text{in } [0, 1], \quad u(0) = u(1) = 0, \quad (3.3.1)$$

$f \in C^0([0, 1])$, and its cubic spline collocation discretization based on the knot set

$$\mathcal{T} := \{jh\}_{j=0}^M, \quad h := 1/M, \quad M \in \mathbb{N}, \quad (3.3.2)$$

and on the set of collocation nodes $\{jh\}_{j=1}^{M-1}$.

(3.3a) Refresh your knowledge of the basic idea of discretization by collocation as explained in the beginning of [NPDE, Section 1.5.3].

(3.3b) Any cubic spline $s \in \mathcal{S}_{3,\mathcal{T}}$ has the local monomial representation

$$s(x) = a_i(x - x_{i-1})^3 + b_i(x - x_{i-1})^2 + c_i(x - x_{i-1}) + d_i, \quad x_{i-1} < x < x_i, \quad i = 1, \dots, M, \quad (3.3.3)$$

with coefficients $a_i, b_i, c_i, d_i \in \mathbb{R}$. This local monomial representation will be used in the sequel.

Denote $f_i := f(x_i)$, $i = 0, \dots, M$.

Show that the collocation conditions [NPDE, Eq. (1.5.100)] imply the following formulas

$$a_i = \frac{f_{i-1} - f_i}{6h}, \quad i = 1, \dots, M, \quad (3.3.4a)$$

$$b_1 = 0 \quad (3.3.4b)$$

$$b_i = -\frac{1}{2}f_{i-1}, \quad i = 2, \dots, M, \quad (3.3.4c)$$

for the coefficients of the local monomial representation of the spline solving the collocation equations (in the first equation we have set $f_M := s''(1) = 0$).

Solution: A natural cubic spline $s : [0, 1] \rightarrow \mathbb{R}$ satisfies

$$(i) \quad s \in \mathcal{C}^2([0, 1]),$$

$$(ii) \quad s|_{[x_{j-1}, x_j]} \in \mathcal{P}_3(\mathbb{R}),$$

$$(iii) \quad s''(0) = s''(1) = 0.$$

With (ii) we know that for $x_{i-1} \leq x \leq x_i$, $i = 1, \dots, M$ we have

$$\begin{aligned} s(x) &= a_i(x - x_{i-1})^3 + b_i(x - x_{i-1})^2 + c_i(x - x_{i-1}) + d_i, \\ s'(x) &= 3a_i(x - x_{i-1})^2 + 2b_i(x - x_{i-1}) + c_i, \\ s''(x) &= 6a_i(x - x_{i-1}) + 2b_i. \end{aligned}$$

Since $-s''(x_{i-1}) = f(x_{i-1})$ we have .

$$2b_i = -f(x_{i-1}) = -f_{i-1} \quad \Rightarrow \quad b_i = \frac{1}{2}f_{i-1}, \quad i = 2, \dots, M,$$

and $b_1 = 0$ since $s''(0) = 0$. We also have for $-s''(x_i) = f(x_i)$ that

$$-6a_i h - f_{i-1} = f_i \quad \Rightarrow \quad a_i = \frac{f_{i-1} - f_i}{6h} \quad i = 1, \dots, M.$$

(3.3c) The equations (3.3.4) are too few to determine all unknown coefficients. To obtain further equations, use the continuity conditions for the “zeroth” and first derivative of a cubic spline, the collocation and boundary conditions at $x = 0$ and $x = 1$, and the relationships (3.3.4) to establish

$$d_1 = 0 \quad (3.3.5a)$$

$$d_{i+1} - 2d_i + d_{i-1} = -\frac{h^2}{6}(f_i + 4f_{i-1} + f_{i-2}), \quad 2 \leq i \leq M-1, \quad (3.3.5b)$$

$$c_2 - c_1 = -\frac{h}{2}f_1, \quad (3.3.5c)$$

$$c_{i+1} - c_i = -\frac{h}{2}(f_i + f_{i-1}), \quad 2 \leq i \leq M-1, \quad (3.3.5d)$$

$$c_M h + d_M = \frac{h^2}{3}f_{M-1}, \quad (3.3.5e)$$

$$c_{M-1} h + d_{M-1} - d_M = \frac{h^2}{6}(2f_{M-2} + f_{M-1}). \quad (3.3.5f)$$

HINT: Recall the formulas for interpolating natural cubic splines from [NCSE, Sect. 3.8.1].

Solution: Since $s'(x)$ is continuous we have

$$\begin{aligned} c_{i+1} = s'(x_i) &= 3a_i h^2 + 2b_i h + c_i \\ &= 3 \frac{f_{i-1} - f_i}{6h} h^2 - 2 \frac{f_{i-1}}{2} h + c_i \\ &= -\frac{f_i + f_{i-1}}{2} h + c_i. \end{aligned}$$

Therefore $c_{i+1} - c_i = -\frac{f_i + f_{i-1}}{2} h$ for $i = 2, \dots, M-1$ and $c_{i+1} - c_i = -\frac{f_i}{2} h$ for $i = 1$.

Furthermore, since $s(x)$ is continuous, we obtain

$$\begin{aligned} d_i = s(x_i) &= a_{i-1} h^3 + b_{i-1} h^2 + c_{i-1} h + d_{i-1}, \\ \Rightarrow d_{i+1} - d_i &= a_i h^3 + b_i h^2 + c_i h; \end{aligned}$$

this implies

$$\begin{aligned} \Rightarrow d_{i+1} - 2d_i + d_{i-1} &= h^3 \left(\frac{f_{i-1} - f_i - f_{i-2} + f_{i-1}}{6h} \right) + h^2 \left(\frac{-f_{i-1} + f_{i-2}}{2} \right) + \frac{h^2}{2} (-f_{i-1} - f_{i-2}) = \\ &= -\frac{h^2}{6} (f_i + 4f_{i-1} + f_{i-2}), \end{aligned}$$

for $i = 2, \dots, M-1$, and

$$c_{M-1} h + d_{M-1} - d_M = -a_{M-1} h^3 - b_{M-1} h^2 = \frac{h^2}{6} (2f_{M-2} + f_{M-1}).$$

Additionally, the boundary conditions $s(0) = s(1)$ imply, respectively, $d_1 = 0$ and

$$c_M h + d_M = \frac{h^2}{3} f_{M-1}.$$

(3.3d) Write a MATLAB function

```
function [a,b,c,d] = natcubsplinecoll(M,f)
```

that computes the local monomial expansion coefficients of the natural cubic spline collocation solution of (3.3.1) with node set (3.3.2). These coefficients are to be returned in the row vectors a, b, c, d of length M. The argument f is a handle to the (continuous) right hand side function f.

Solution: See Listing 3.16.

Listing 3.16: Code for subproblem (3.3d).

```
1 function [a,b,c,d] = natcubsplinecoll(M,f)
2
3 % Computes the coefficients in the monomial expansion of the spline
  for
4 % equispace knots
5 % INPUT:
6 % M = number of intervals
7 % f = FHandle to right-hand side
8 % OUTPUT:
```

```

9  % column vectors of coefficients in the monomial expansion
10
11 h = 1/M;
12 x = linspace(0,1,M+1);
13 fval = f(x);
14 fval = fval(:);
15 a = (fval(1:end-1)-fval(2:end))/(6*h);
16 b = [0; -fval(2:end-1)/2];
17 A11 = diag([-ones(M-1,1);h])+diag(ones(M-1,1),1);
18 A12 = zeros(M,M);
19 A12(end,end) = 1;
20 A21 = zeros(M,M);
21 A21(end,end-1) = h;
22 A22 =
    diag([1;-2*ones(M-2,1);-1])+diag([0;ones(M-2,1)],1)+diag([ones(M-2,1);1],-1);
23 A = [A11 A12; A21 A22];
24 RHS = zeros(2*M,1);
25 RHS(1:M-1)=(fval(2:end-1)+fval(1:end-2))*(-h/2);
26 RHS(M) = fval(end-1)*h^2/3;
27 RHS(M+1) = 0;
28 RHS(M+2:end-1) = -h^2/6*(fval(3:M)+4*fval(2:M-1)+fval(1:M-2));
29 RHS(end) = h^2/6*(2*fval(M-1)+fval(M));
30 res = A\RHS;
31 c = res(1:M);
32 d = res(M+1:end);

```

(3.3e) Plot the approximate solution of (3.3.1) with $f(x) = \sin(2\pi x)$ obtained by natural cubic spline collocation on an equidistant node set (3.3.1) with $M = 5, 10, 20$.

Solution: See Listing 3.17.

Listing 3.17: Code for subproblem (3.3e).

```

1  f = @(x) sin(2*pi.*x(:));
2
3  M = 20;
4  [a, b, c, d]=natcubsplinecoll(M,f);
5  x = linspace(0,1,M+1);
6  xx = linspace(0,1,4*M+1);
7  yy = [];
8  for i=1:M
9      yval = a(i)*(xx((i-1)*4+1:i*4+1)-x(i)).^3 +
            b(i)*(xx((i-1)*4+1:i*4+1)-x(i)).^2 +
            c(i)*(xx((i-1)*4+1:i*4+1)-x(i)) + d(i);
10     yy = [yy(1:end-1) yval];
11 end
12
13 plot(xx,yy,'o')

```

For $M = 20$ we obtain the plot shown in Figure 3.4.

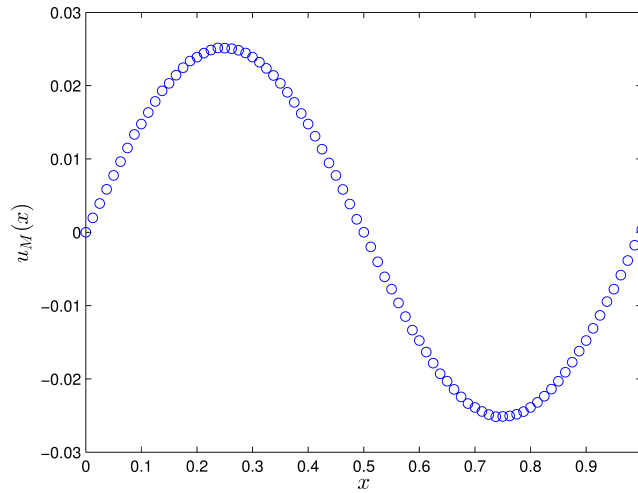


Figure 3.4: Spline collocation method for $M = 20$ mesh points.

(3.3f) As outlined in [NCSE, Rem. 9.1.4], investigate the convergence of the method on the same problem as in subproblem (3.3e) with $M = 5, 10, 20, 40, 80$ in the L^2 -norm. In order to approximate the integral for computation of the norm use the composite 2-point Gauss quadrature rule.

HINT: The nodes and weights for 2-point Gaussian quadrature on $[-1, 1]$ are $\zeta_1 = -\frac{1}{3}\sqrt{3}$, $\zeta_2 = \frac{1}{3}\sqrt{3}$, $\omega_1 = 1$, $\omega_2 = 1$. This quadrature rule has to be transformed to all mesh cells (x_{j-1}, x_j) , see [NCSE, Rem. 10.1.3].

Solution:

Listing 3.18: Implementation for subproblem (3.3f)

```

1  %%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
2  % SPLINECONV computes the convergence rate of a
3  %   solution of a model BVP problem
4  %   using the spline collocation method
5  %
6
7  % input data
8  M = 2.^(5:10)'; % number of nodes
9  f = @(x) sin(2*pi.*x(:)); % right-hand side
10 u_exact = @(x) 1/(4*pi^2)*sin(2*pi.*x(:)); % exact solution
11
12 % loop over M
13 l2error = zeros(length(M),1);
14 for i = 1:length(M)
15
16     % mesh
17     h = 1/M(i);
18     x = linspace(0,1,M(i)+1)';
19
20     % coefficients
21     [a, b, c, d]=natcubsplinecoll(M(i),f);
22
23     % compute quadrature points
24     xpts = repmat(x(1:length(x)-1)+h/2,2,1)'; xpts = xpts(:);

```

```

25     gqpts = repmat([-1;1]/sqrt(3),1,M(i))'*h/2; gqpts = gqpts(:) +
        xpts;
26     gqpts = sort(gqpts);
27
28     % compute solution in quadrature points
29     u = [];
30     for j=1:M(i)
31         uval = a(j)*(gqpts((j-1)*2+1:j*2)-x(j)).^3 +
                b(j)*(gqpts((j-1)*2+1:j*2)-x(j)).^2 +
                c(j)*(gqpts((j-1)*2+1:j*2)-x(j)) + d(j);
32         u = [u; uval];
33     end
34
35     % compute error
36     l2error(i) = sqrt(sum(h/2*(u-u_exact(gqpts)).^2));
37 end
38
39 % compute convergence rate
40 p = polyfit(log(M),log(l2error),1);
41 s = p(1);
42 fprintf('Convergence rate s = %2.4f\n', abs(s));
43
44 % plot convergence rate
45 figure(1)
46 h = axes;
47 loglog(M,l2error,'bo-')
48 hold on;
49 add_Slope(gca, 'NorthEast', p(1));
50 grid on
51 set(h, 'FontSize', 14);
52 xlabel('M')
53 ylabel('L^2-error')
54 %print -depsc splineconv.eps

```

We obtain the optimal convergence rate $s = 2$ as shown in [Figure 3.5](#).

Listing 3.19: Testcalls for [Problem 3.3](#)

```

1
2 fprintf(' \n\n##natcubsplinecoll:')
3
4 f = @(x) sin(2*pi.*x(:));
5 M = 5;
6 [a, b, c, d]=natcubsplinecoll(M,f)

```

Listing 3.20: Output for Testcalls for [Problem 3.3](#)

```

1 >> test_call
2
3 ##natcubsplinecoll:
4 a =

```

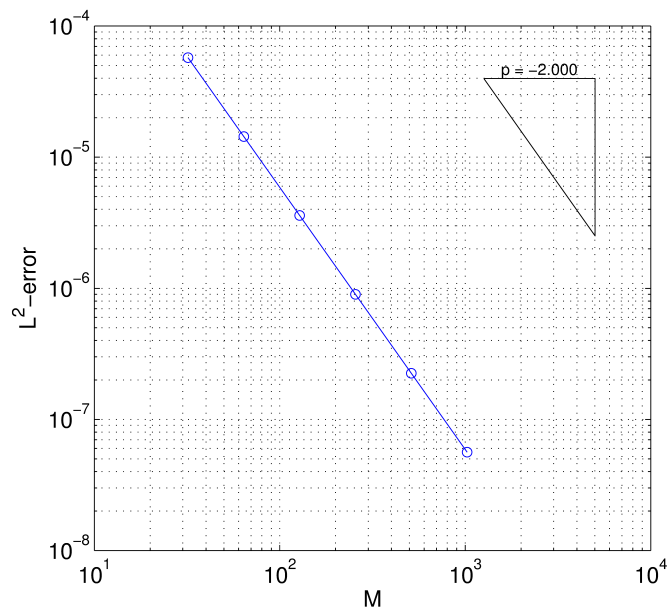


Figure 3.5: Convergence rate for the spline collocation method

```

5
6   -0.7925
7    0.3027
8    0.9796
9    0.3027
10   -0.7925
11
12  b =
13
14      0
15   -0.4755
16   -0.2939
17    0.2939
18    0.4755
19
20  c =
21
22    0.1376
23    0.0425
24   -0.1114
25   -0.1114
26    0.0425
27
28  d =
29
30      0
31    0.0212
32    0.0131
33   -0.0131

```


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References

[NPDE] [Lecture Slides](#) for the course “Numerical Methods for Partial Differential Equations”.SVN revision # 74075.

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