

NTNU

PROJECT

Least Squares Finite Element Method

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Abstract

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Project

Least Squares Finite Element Method

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The Thesis Abstract is written here (and usually kept to just this page). The page is kept centered vertically so can expand into the blank space above the title too...

Notation

LAH List Abbreviations **Here**

Chapter 1

Theory

1.1 Informal introduction to LSFEM

The Least Squares Finite Element Method is a numerical method similar to mixed galerkin. But with a different approach regarding defining the bilinear functional. Let us look at a system of first order differential equations on the form

$$Au = f \text{ in } \Omega \quad (1.1)$$

$$Bu = g \text{ on } \partial\Omega. \quad (1.2)$$

Where A is a partial differential operator defined as

$$A = \sum_{i=1}^n A_i \frac{\partial}{\partial x_i} + A_0. \quad (1.3)$$

n beeing the number of dimensions of the domain Ω . Let us initially assume without loss of generality that $g = 0$. Further we require $f \in L_2(\Omega)$ and choose $V = \{v \in L_2(\Omega) | v = 0 \text{ on } \partial\Omega\}$. A residual is defined

$$R(v) = Av - f, \quad (1.4)$$

and a functional

$$J(v) = \frac{1}{2} \|R(v)\|_0^2. \quad (1.5)$$

The solution u and its gradient needs to be in L^2 for the functional to make sense, hence u is restricted to the space $H^1(\Omega)$. By minimizing J we obtain

$$\lim_{t \rightarrow 0} \frac{d}{dt} J(u + tv) = \int_{\Omega} (Av)^T (Au - f) d\Omega = 0, \quad \forall v \in V. \quad (1.6)$$

We can now write a variational formulation of the least-squares method: Find $u \in V$ such that

$$Q(u, v) = F(v), \quad \forall v \in V, \quad (1.7)$$

where

$$Q(u, v) = (Au, Av), \quad (1.8)$$

$$F(v) = (f, Av). \quad (1.9)$$

[?] Notice that the bilinear form Q is symmetric, this is an important advantage LSFEM has over regular galerkin methods. The bilinear form that surged from a first-order problem by the LSFEM leads us to a variational formulation similar to the one obtained from a second order problem by regular FEM. Generally the bilinear form from LSFEM will correspond to a bilinear form of a problem of twice the order obtained using FEM. In order to avoid problems of large complexity a higher order PDE should therefore be transformed to a system of first order PDE's (similar to a mixed galerkin approach) before defining the least squares functional.

[?]

1.2 Formal formulation of LSFEM

Let us look at a general boundary value problem where $f \in Y(\Omega)$, $g \in B(\partial\Omega)$, $\mathcal{B}: X(\partial\Omega) \rightarrow B(\partial\Omega)$ and $\mathcal{L}: X(\Omega) \rightarrow Y(\Omega)$. Find $u \in X(\Omega)$ such that

$$\mathcal{L}u = f \quad \text{in } \Omega \quad (1.10)$$

$$\mathcal{B}u = g \quad \text{on } \partial\Omega. \quad (1.11)$$

Whenever this BVP has a unique solution, a least-squares functional can be defined as

$$J(u; f, g) = \|\mathcal{L}u - f\|_Y^2 + \|\mathcal{B}u - g\|_B^2 \quad (1.12)$$

and the corresponding minimization problem is then given as

$$\min_{u \in X} J(u; f, g) \quad (1.13)$$

For any well-posed problem $\exists \alpha, \beta > 0$ such that

$$\alpha \|u\|_X^2 \leq J(u; 0, 0) = (\mathcal{L}u, \mathcal{L}u)_Y \leq \beta \|u\|_X^2. \quad (1.14)$$

The fact that our functional is norm-equivalent is of crucial importance to a successful LS-method. It is therefore important that the spaces X, Y and B is chosen such that the LS-functional defines a norm is equivalent to $\|\cdot\|_X$. **Need to show this ?** Minimizing this functional is equivalent to solving the Euler-lagrange equations formulated as

$$\text{find } u \in X \text{ such that } Q(u, v) = F(v) \quad \forall v \in X \quad (1.15)$$

Where $Q(u, v)$ is the continuous bilinear form given as $(\mathcal{L}u, \mathcal{L}v)_Y$ and $F(v)$ is the bounded linear functional given as $(\mathcal{L}v, f)_Y$. Notice that $Q(u, v)$ defines an inner product and $Q(u, u)^{1/2}$ defines the corresponding norm.

In order to solve this numerically we define the function spaces X^h, Y^h and B^h and the corresponding variational formulation is then written as

$$\text{find } u^h \in X^h \text{ such that } Q(u^h, v^h) = F(v^h) \quad \forall v^h \in X^h. \quad (1.16)$$

[?]

1.2.1 Error analysis

Let u be the analytical solution of a problem of the type 1.11, u^h is our numerical solution to 1.16 and u_\perp^h is the orthogonal projection of u in X_h .

$$Q(u - u^h, u - u^h) = Q(u - u^h, u - u_\perp^h) + Q(u - u^h, u_\perp^h - u^h) \quad (1.17)$$

$$= Q(u - u^h, u - u_\perp^h) \quad (1.18)$$

$$\leq \beta \|u - u^h\|_{X_h} \|u - u_\perp^h\|_{X_h}. \quad (1.19)$$

The first equality is due to adding and subtracting u_\perp^h , because both u^h and u_\perp^h solves the variational formulation we can cancel the last term, and by using the norm-equivalency from 1.14 and schwartz inequality we get the last expression. Now by applying the first inequality of 1.14 we end up with

$$\|u - u^h\|_{X_h} \leq \frac{\beta}{\alpha} \|u - u_\perp^h\|_{X_h} = \min_{w^h \in X_h} \frac{\beta}{\alpha} \|u - w^h\|_{X_h}. \quad (1.20)$$

Hence we can show that the Least squares method provides a convergence result of similar order as the Finite element method.

1.3 Example - Poisson problem

The poisson problem is defined as

$$-\Delta u = f \text{ in } \Omega \quad (1.21)$$

$$u = g \text{ on } \partial\Omega \quad (1.22)$$

Let us first consider the homogenous case. The straight forward LSFEM approach is to define $w = -\nabla u$ and solve the system of equations

$$w + \nabla u = 0 \text{ in } \Omega \quad (1.23)$$

$$\nabla w = f \text{ in } \Omega \quad (1.24)$$

$$u = 0 \text{ on } \partial\Omega. \quad (1.25)$$

which can be written in the same form as 1.11 with $u = w \oplus u$, $f = (0, 0, f)$, $g = 0$, $\mathcal{B} = (0, 0, 1)^T$ and \mathcal{L} given as

$$\mathcal{L} = \begin{bmatrix} 1 & 0 & \partial/\partial x \\ 0 & 1 & \partial/\partial y \\ \partial/\partial x & \partial/\partial y & 0 \end{bmatrix} \quad (1.26)$$

We define the search space $X = H^1(\Omega; \text{div}) \times H_0^1(\Omega)$ and the solution space $Y \times B = [L^2(\Omega)]^3 \times L^2(\Omega)$ and the functional can then be defined as in 1.12. The variational formulation of the problem can be stated. Find $u \in X$ s.t.

$$Q(u, \phi) = F(\phi) \quad \forall \phi \in X. \quad (1.27)$$

We require that $f \in Y$. Notice that the spaces X and Y chosen as described above fulfill the condition 1.14.

1.3.0.1 non-homogenous boundary conditions

If $g \neq 0$ then we simply define a lifting function $R_g \in X$ such that $R_g(\partial\Omega) = g(\partial\Omega)$. By defining $\tilde{u} = u - R_g$ we can replace u in the variation formulation and get

$$Q(\tilde{u} + R_g, \phi) = F(\phi) \quad (1.28)$$

$$Q(\tilde{u}, \phi) + Q(R_g, \phi) = F(\phi) \quad (1.29)$$

$$Q(\tilde{u}, \phi) = F(\phi) - Q(R_g, \phi) \quad (1.30)$$

$$Q(\tilde{u}, \phi) = \tilde{F}(\phi) \quad (1.31)$$

1.4 Example - Diffusion convection problem

The diffusion convection problem to be analyzed is given as

$$-\Delta u + b \cdot \nabla u = f \text{ in } \Omega \quad (1.32)$$

$$u = g \text{ on } \partial\Omega \quad (1.33)$$

The straight forward LSFEM approach is to define $w = -\nabla u$ and solve the system of equations

$$w + \nabla u = 0 \text{ in } \Omega \quad (1.34)$$

$$\nabla \cdot w - b \cdot w = f \text{ in } \Omega \quad (1.35)$$

$$u = 0 \text{ on } \partial\Omega. \quad (1.36)$$

which can be written in the same form as 1.11 with $u = w \oplus u$, $f = (0, 0, f)$, $g = 0$, $\mathcal{B} = (0, 0, 1)^T$ and L given as

$$\mathcal{L} = \begin{bmatrix} 1 & 0 & \partial/\partial x \\ 0 & 1 & \partial/\partial y \\ \partial/\partial x - b_1 & \partial/\partial y - b_2 & 0 \end{bmatrix} \quad (1.37)$$

We define the search space $X = H^1(\Omega; \text{div}) \times H_0^1(\Omega)$ and the solution space $Y \times B = [L^2(\Omega)]^3 \times L^2(\Omega)$ and the functional can then be defined as in 1.12. The variational formulation of the problem can be stated. Find $u \in X$ s.t.

$$Q(u, \phi) = F(\phi) \quad \forall \phi \in X. \quad (1.38)$$

We require that $f \in Y$. Notice that the spaces X and Y chosen as described above fulfill the condition 1.14.

Chapter 2

New Theory

2.1 Using the Least squares method to gain stability

In regular galerkin FEM for the Diffusion-Transport equation you end up with the variational formulation

$$a(u, v) = (f, v)_0 \quad \forall v \in V \quad (2.1)$$

where the bilinear functional is given as

$$a(u, v) = \mu \int_{\Omega} \nabla u \cdot \nabla v d\Omega + \int_{\Omega} v(b \cdot \nabla u) d\Omega \quad (2.2)$$

We can define a new norm from this functional in the following manner

$$a(u, u) = \mu \int_{\Omega} \nabla u \cdot \nabla u d\Omega + \int_{\Omega} u(b \cdot \nabla u) d\Omega \quad (2.3)$$

$$= \mu \|\nabla u\|_0^2 + \frac{1}{2} \int_{\Omega} b \cdot \nabla u^2 d\Omega \quad (2.4)$$

$$= \mu \|\nabla u\|_0^2 - \frac{1}{2} \int_{\Omega} u^2 (\nabla \cdot b) d\Omega \quad (2.5)$$

$$(2.6)$$

Let $\gamma_0 \leq \nabla \cdot b \leq \gamma_1$, we can now make a lower and upper bound for the bilinear form

$$\mu \|\nabla u\|_0^2 - \frac{\gamma_1}{2} \|u\|_0^2 \leq a(u, u) \leq \mu \|\nabla u\|_0^2 - \frac{\gamma_0}{2} \|u\|_0^2 \quad (2.7)$$

should I change signs in the equation??

It is clear that for small μ and large γ_1 the bilinear form is no longer coercive and thus our convergence requirements are no longer valid. Now, let us explore what happens if we add the variational formulation obtained from the LSFEM to the standard FEM formulation

$$a(u, v) + Q(u, v) = (f, v) + F(v) \quad (2.8)$$

$$\mathring{a}(u, v) = \mathring{f}(v) \quad (2.9)$$

Let us derive an upper bound for the norm created by the combined bilinear form $a(\cdot, \cdot)$.

$$\mathring{a}(u, u) = \mathring{f}(u) \quad (2.10)$$

$$= (f, u) + (\mathcal{L}u, f) \quad (2.11)$$

$$\leq \|f\|_0^2 \|u\|_0^2 + \|\mathcal{L}u\|_0^2 \|f\|_0^2 \quad (2.12)$$

$$= (\|u\|_0^2 + \|\mathcal{L}u\|_0^2) \|f\|_0^2 \quad (2.13)$$

Chapter 3

Implementation

For the general problem [1.11](#) the functional Q will take the form

$$Q(u, v) = \int_{\Omega} (\mathcal{L}v)^T (\mathcal{L}u) d\Omega. \quad (3.1)$$

Implementing Q requires two sets of basis functions $\{N_i\}$ that describes the search and solution space. In this project assignment the search and solution space will be described by the the same set of basis functions which will depend on the method applied. u is discretized as

$$u_h = \sum_{I=0}^K a_I N_I. \quad (3.2)$$

Since equation [1.15](#) requires equality for all test functions in the search space we simply solve the equation for each basis function. We are therefore left with a system of K

equations. Equation 3.1 can then be written for each test function as

$$Q(u_h, N_I) = \int_{\Omega} (\mathcal{L}N_I)^T (\mathcal{L}u_h) d\Omega \quad (3.3)$$

$$= \int_{\Omega} (\mathcal{L}N_I)^T (\mathcal{L} \sum_{J=1}^K a_J N_J) d\Omega \quad (3.4)$$

$$= \sum_{J=1}^K \int_{\Omega} (\mathcal{L}N_I)^T (\mathcal{L}a_J N_J) d\Omega \quad (3.5)$$

$$= \sum_{J=1}^K \int_{\Omega} (\mathcal{L}N_I)^T (\mathcal{L}a_J N_J) d\Omega \quad (3.6)$$

$$= \sum_{J=1}^K \int_{\Omega} (\mathcal{L}N_I)^T (\mathcal{L}N_J) d\Omega \cdot a_J. \quad (3.7)$$

The total system of equation for all test functions can then be written as a matrix equation

$$Au = F. \quad (3.8)$$

Where $A_{I,J} = \int_{\Omega} (\mathcal{L}N_I)^T (\mathcal{L}N_J) d\Omega$. For the poisson equation $A_{I,J}$ will be a 3-by-3 matrix on the form

$$A_{I,J} = \int_{\Omega} \begin{bmatrix} N_I N_J + N_{I,x} N_{J,x} & N_{I,x} N_{J,y} & N_I N_{J,x} \\ N_{I,y} N_{J,x} & N_I N_J + N_{I,y} N_{J,y} & N_I N_{J,y} \\ N_{I,x} N_J & N_{I,y} N_J & N_{I,x} N_{J,x} + N_{I,y} N_{J,y} \end{bmatrix} d\Omega$$

Notice that by doing the splitting of variables we obtain a system of equations three times as big as if we were to solve the equation directly.

3.1 LSFEM for poisson

finite element space X_h^1 ? quadrature...

3.2 LS spectral method for poisson

The spectral implementation is done using Gauss Lobatto nodes and quadrature and the lagrange functions based on the GL nodes as basis functions. Notice that the discrete

solution u_h consist of the discretizations of both u and $w = \nabla u$. u_h can be structured blockwise such that or nodewise. By choosing a blockwise representation the final system of equations can be written as

$$\begin{bmatrix} A_{1,1} & A_{1,2} & A_{1,3} \\ A_{2,1} & A_{2,2} & A_{2,3} \\ A_{3,1} & A_{3,2} & A_{3,3} \end{bmatrix} \begin{bmatrix} u^h \\ w_1^h \\ w_2^h \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ F^h \end{bmatrix}.$$

Where each block $A_{i,j}$ corresponds to calculating element i, j in the matrix 3 for all the indices I, J . In order to implement this matrix it is convenient to write it in a compact form using the kronecker tensor product. The components needed for this formulation is the $n \times n$ diagonal matrix W with the GLL-weights along the diagonal and the $n \times n$ matrix $(L)_{i,j} = l'_j(x_i)$ where l_j is the j th lagrange polynomial and x_i is the i th node in either x or y direction. Note that the formulation is based on a grid of GLL-nodes in both x and y direction.

$$A = \begin{bmatrix} W \otimes (L^T W L + W) & W L \otimes L^T W & W \otimes W L \\ L^T W \otimes W L & (L^T W L + W) \otimes W & W L \otimes W \\ W \otimes L^T W & L^T W \otimes W & L^T W L \otimes W + W \otimes L^T W L \end{bmatrix}$$

do I need to show how this is derivated?

Similarly without the reformulation of the PDE and with regular galerkin formulation the stiffness matrix will simply be $A_{3,3} = W \otimes L^T W L + L^T W L \otimes W$

Chapter 4

Results

4.1 The main differences

Using least squares will always give you a SPD system of equations which can be advantageous. However for second order equations this system is three times as big as if we were to solve it using more standard methods. Comparing the correctness of the solution as done in figure ?? shows that the convergence rate is the same as for standard methods, but the the value of the residual is slightly higher for the least squares method. This can be explained by the functional that is minimized. Notice that in the least squares methods you minimize the square of the residual. since the correctness of both methods are restricted by the smoothness of the solution and the number of discrete points LS-methods will minimize the residual squared down to a given precision and hence the residual itself to a slightly higher value. The condition number is also worth comparing ...

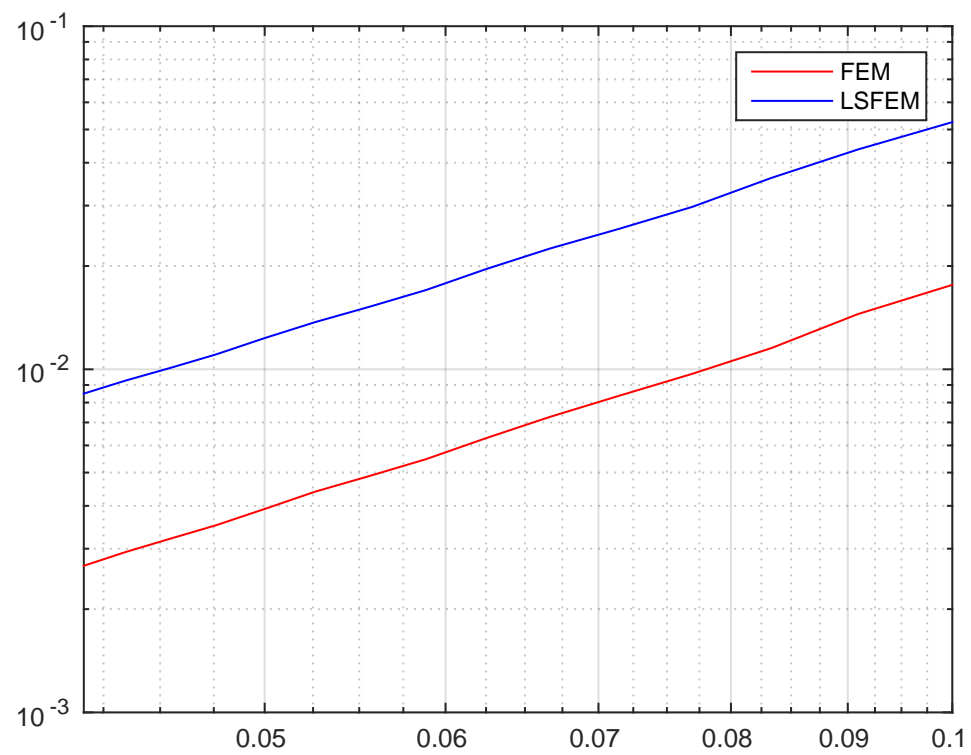


FIGURE 4.1: convergence of LSFEM and FEM

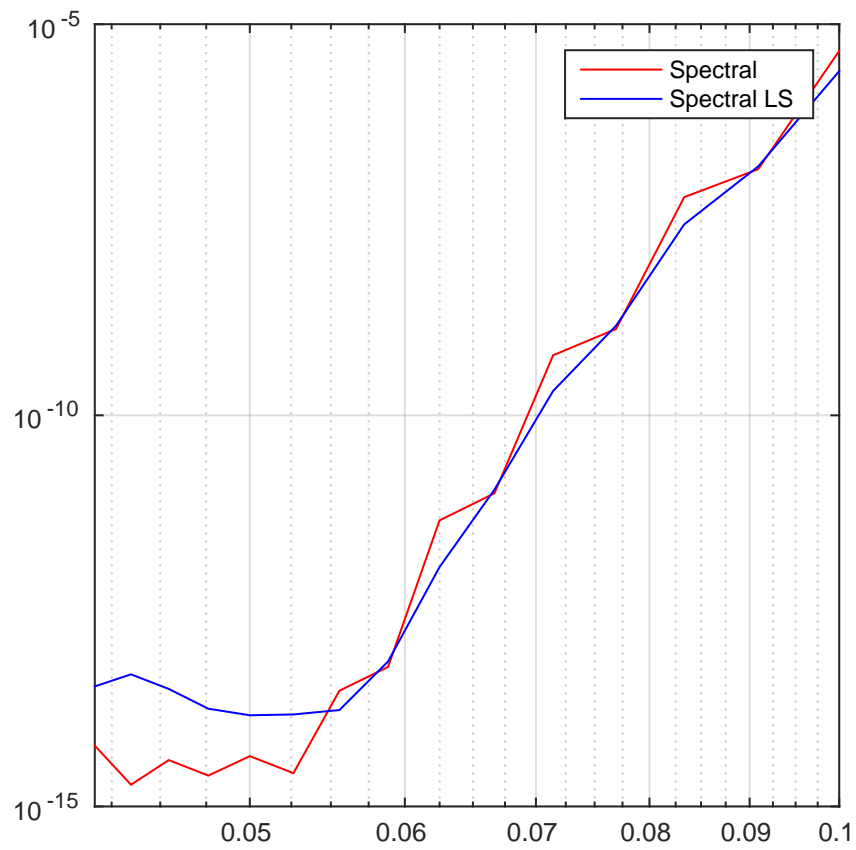


FIGURE 4.2: convergence of Spectral and LS-Spectral method

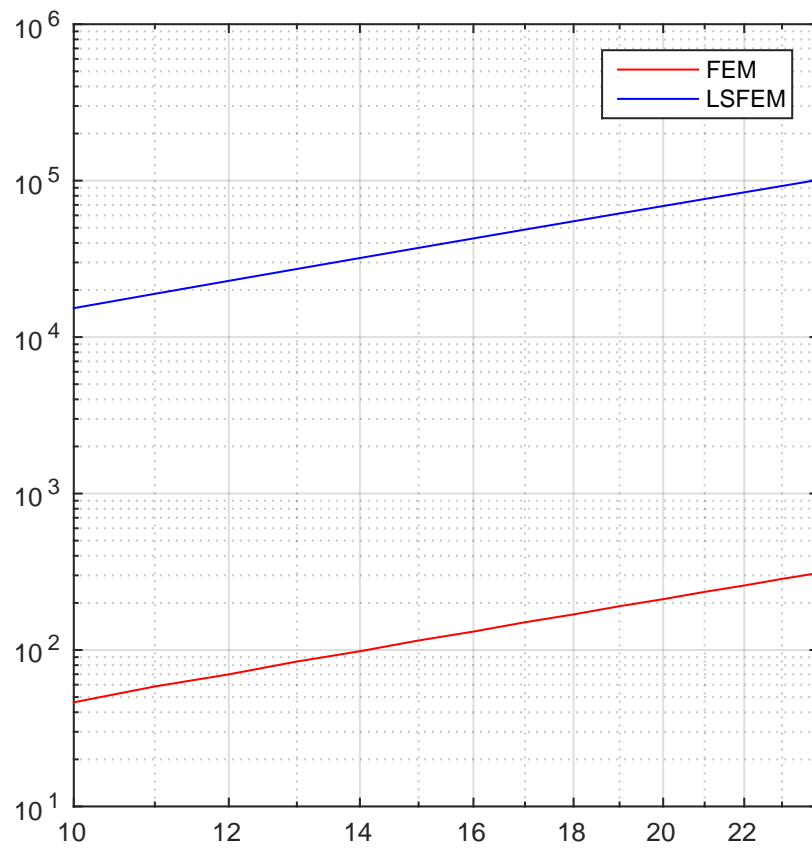


FIGURE 4.3: condition number of LSFEM and FEM

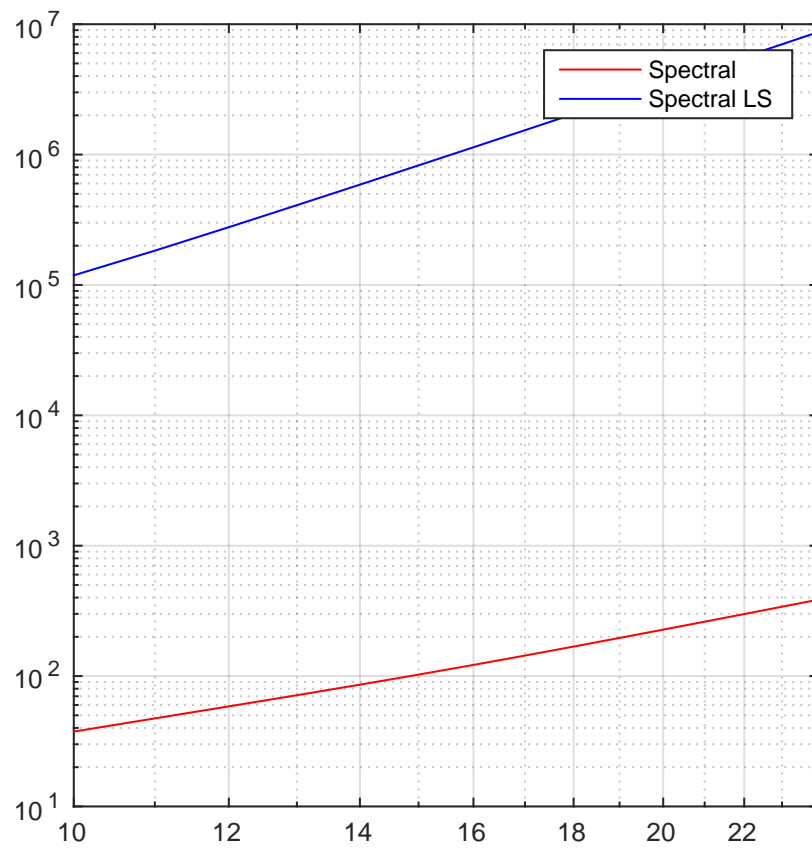


FIGURE 4.4: condition number of Spectral and LS-Spectral method

Appendix A

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