

Contents

Preface		ix
0.1	Life is complex	ix
0.2	Readings	ix
0.3	Least Squares	ix
0.4	Linear Algebra and Matrix Analysis	x
0.5	Numerical Linear Algebra	x
0.6	Discussions on Least Squares	x
I	Rudiments	1
1	Least Squares Problems	3
1.1	Linear Systems	3
1.1.1	$\ \mathbf{A}x - b\ = 0$	3
1.1.2	$\ \mathbf{A}x - b\ > 0$	4
1.2	Least Squares Solutions	4
1.2.1	Zonal Approximation	5
1.2.2	Modal Approximation	7
1.2.3	Errors	8
1.3	Least Squares Problem	8
2	Least Squares Solutions	9
2.1	Fundamental Theorem of Linear Algebra	9
2.2	Singular Value Decomposition - I	12
2.2.1	SVD Theorem	13
2.2.2	SVD and Least Squares	14
2.3	Singular Value Decomposition - II	16
2.3.1	Σ Gymnastics	16
2.3.2	Fundamental Projectors	18
2.4	Least Squares Solution - Again	19
2.5	Reflection Invariance	19

II	Modal Example	23
3	Solution Via Calculus	27
3.1	Modal Approximation	27
3.2	Bevington Example	28
3.2.1	Problem Statement	28
3.2.2	Normal Equations via Calculus	29
3.3	Numerical Results	32
3.3.1	Exact Form	32
3.3.2	Computed Form	33
3.4	Visualization	34
3.4.1	Seeing the Solution	34
3.4.2	Seeing the Uncertainty	36
3.4.3	Digging Deeper	37
4	Solution via SVD	41
4.1	Computing the SVD	41
4.1.1	Singular values	42
4.1.2	Domain matrix	42
4.1.3	Codomain matrix	43
4.1.4	Error terms	44
4.2	Visualization	45
5	Solution Via Other Methods	51
5.1	Normal Equations from Vectors	51
5.1.1	Composing the normal equations	52
5.2	QR Decomposition	53
5.2.1	Computing the QR Decomposition	53
6	Finer Points	55
6.1	Invariances	55
6.1.1	Translation Invariance	55
6.1.2	Reflection Invariance	57
6.2	Fitting To Higher Orders	58
7	Tricks	61
7.1	Where Do Parallel Lines Cross?	61
7.1.1	Intersecting Lines	61
III	Zonal Example	63
8	Zonal Example	65
8.1	Problem	65
8.1.1	Zonal Subsection	65
9	Lines	67

9.1	Face-centered cubic lattice	67
9.2	Model	68
9.3	Solution	71
9.4	Problem Statement	72
9.5	Data	72
9.6	Results	72
9.6.1	Least Squares Results	72
9.6.2	Apex Angles	72
9.6.3	Qualitative Results	72
10	Crystals	79
11	Stitching Local Maps	81
11.1	What is stitching?	81
11.2	Stitch ϕ	81
11.2.1	Genesis	81
11.2.2	Data	81
11.2.3	Data and results	85
11.2.4	Linear System	85
11.2.5	Least Squares Arbitration	87
11.3	Stitch $\nabla\phi$	90
12	Gradient I	93
12.1	One Dimension	93
IV	Applications: Nonlinear Problems	95
13	Linearization	97
13.1	Powers Laws and Exponentials	97
13.2	Learning Curve	97
13.2.1	Problem Statement	98
13.2.2	Solution	99
13.2.3	Results	101
13.3	What Not To Do	101
13.3.1	Logarithmic Transform	101
13.3.2	Linear Transformation	104
13.3.3	Reflection Test Fails	106
13.4	Radioactive Decay	106
13.4.1	Theory	106
13.4.2	Problem Statement	106
13.4.3	Results	106
14	Population Growth	111
14.1	Model	111

14.2	Problem Statement	112
14.3	Data	112
14.4	Example	112
14.5	Polynomials	112
V	Appendices	121
A	Exercises	123
A.1	Linear systems	123
A.1.1	6.1.1	123
VI	Backmatter	125
	Bibliography	127

List of Figures

1.1	The residual error $\ r\ _2$ given in (1.5).	5
1.2	Scalar function ϕ and approximations.	6
2.1	Decomposing the data vector.	16
2.2	Projections of the data vector.	19
2.3	Data vector resolved as $b = b_{\mathcal{R}} + b_{\mathcal{N}}$	20
2.4	Data vector resolved as $b + b_{\mathcal{N}} = b_{\mathcal{R}}$	20
2.5	The merit function for the learning curve showing the minimum and the value.	21
3.1	Measuring the temperature of a bar.	28
3.2	Solution plotted against data with residual errors shown in red. . .	35
3.3	Scatter plot of residual errors.	36
3.4	Scatter plot of residual errors with data points connected.	36
3.5	The merit function.	37
3.6	Another look at the merit function showing the primary error ellipse. .	38
3.7	Whisker plot showing 250 randomly sampled solutions.	38
3.8	Scatter plot showing sampling of solutions.	40
4.1	The solution vector is the mixture of u_1 and u_2 which eliminates error.	46
4.2	Measurement space for the Bevington example	48
4.3	Minimization occurs in the codomain.	49
4.4	Data vector resolved into range and null space components.	49
4.5	Decomposing $\ r = T_{\mathcal{N}}\ _2^2$ into residual error terms r_k^2 of table 4.3. .	49
6.1	Solution after translation along x axis: the slope is invariant. . . .	57
6.2	Scatter plot of residual errors.	58
6.3	The merit function after translation.	59
6.4	The merit function for the learning curve showing the minimum and the value.	59
7.1	Parallel lines and the least squares solution.	62
9.1	A slice of a face-centered cubic lattice showing a single crystal. . .	67

9.2	Simulation output showing atomic shades shaded by potential energy.	68
9.3	Full data set showing inset.	69
9.4	Sample data set showing fit parameters.	69
9.5	Solutions for three data sets.	73
9.6	Apex angles displayed in table ??	76
9.7	Merit functions for the three data sets.	77
11.1	Stitching local maps together to form a global map.	81
11.2	The ideal potential function showing five measurement zones and four overlap bands.	83
11.3	Waterfall diagram showing discretization within measurement zones with left and right zone overlaps.	83
11.4	Stitching unifies the data.	84
11.5	A set of piston adjustments which restores continuity across the domain.	84
11.6	Looking at the merit function on the $p_2 - p_3$ axis.	89
11.7	Pistons from the solution and pistons used to create the data . . .	90
11.8	A set of tilt adjustments which restores continuity of the gradient across the domain.	91
11.9	A function and its gradient.	91
13.1	Merit function for the learning curve in solution space.	99
13.2	Merit function constrained to one parameter b	101
13.3	Merit function for the learning curve in solution space showing the constrained a parameter as a dotted line.	102
13.4	Solution for equations (9.2) and (9.3) using data in table 9.2. . . .	102
13.5	The residual errors in figure 9.4.	103
13.6	Minimization of the merit function for the learning curve.	104
13.7	The merit function for the learning curve showing the minimum and the value	105
13.8	The merit function for the learning curve showing the minimum and the value	107
14.1	Residual error for (10.3) with a_1 , a_2 , and a_3 at optimal values. . .	113
14.2	Residual error for a_1 and a_2 fixed at optimal values.	114
14.3	Solution plotted against data.	114
14.4	Scatterplot of residual errors.	115
14.5	The merit function showing least squares solution.	115
14.6	The merit function showing least squares solution and the null cline.	116
14.7	Total error r^*r by order of fit.	119

List of Tables

0.1	Matrix manipulations for \mathbf{A}^* and \mathbf{A}^T	ix
2.1	The Fundamental Theorem of Linear Algebra	10
2.2	The Fundamental Theorem of Linear Algebra in pictures	11
2.3	Dimensions of the fundamental subspaces for $\mathbf{A} \in \mathbb{C}_\rho^{m \times n}$	12
2.4	Orthonormal spans for the invariant subspaces.	13
2.5	Fundamental Projectors using the pseudoinverse.	18
2.6	Fundamental Projectors using domain matrices.	18
2.7	Reflecting the data for $\mathbf{A}x - T = r$	21
3.1	Problem statement for linear regression.	30
3.2	Raw data and results.	31
3.3	Results for linear regression.	32
3.4	The solution parameters expressed as normal distributions.	39
3.5	Comparing samples to ideal normal distribution.	40
4.1	The column vectors of \mathbf{U}	48
4.2	Singular value decomposition for the system matrix \mathbf{A}	50
4.3	A summary of the residual errors and their contributions to $\ r\ _2$	50
9.1	Data sets and basic results	70
9.2	Problem statement for grain identification by rows (coupled linear regression).	73
9.3	Point membership in data sets shown in figure ??	74
9.4	Excerpted data set.	75
9.5	Least squares results for three axes.	75
9.6	Intermediate results: angles for the axes.	75
9.7	Final results: apex angle measurements	75
11.1	The input data in continuous and discrete form.	82
11.2	Sample showing two zones with overlap.	82
11.3	Measurements displaying the connection between overlap bands in figure ??	85
11.4	Computation of the zone shift values.	85
11.5	Computation of the zone shift values.	86

11.6	Input data	86
11.7	Problem statement for linear regression.	87
11.8	Results for stitching with piston.	88
13.1	Problem statement for learning curve.	100
13.2	The simultaneous conditions defining $\nabla M(a, b) = 0$	100
13.3	Results for learning curve analysis.	103
13.4	Problem statement for radioactive decay.	106
13.5	Logarithmic scaling distorts errors.	108
13.6	Results for radioactive decay.	109
14.1	Problem statement for population model with linear and exponen- tial growth.	112
14.2	Data v. prediction.	113
14.3	Results for census analysis	116
14.4	Fitting the census data with low order polynomials.	117
14.5	Fitting the census data with higher order polynomials.	118
14.6	Projections, by order of fit, for population in 2010.	119

The Devil is in the details.

0.1 Life is complex

\mathbf{A}^* or \mathbf{A}^T

$$\mathbb{R} \in \mathbb{C}$$

Table 0.1. *Matrix manipulations for \mathbf{A}^* and \mathbf{A}^T .*

	\mathbf{A}^*	\mathbf{A}^T
Application	$\mathbf{A} \in \mathbb{C}^{m \times n}$ and $\mathbf{A} \in \mathbb{R}^{m \times n}$	only $\mathbf{A} \in \mathbb{R}^{m \times n}$
Fortran	<code>conjg(transpose(A))</code>	<code>transpose(A)</code>
Mathematica	\mathbf{A}^H <code>ConjugateTranspose[A]</code>	\mathbf{A}^T <code>Transpose[A]</code>
MATLAB	\mathbf{A}' <code>ctranspose(A)</code>	$\mathbf{A}.'$ <code>transpose(A)</code>
Python	<code>A.T.conj()</code>	<code>A.T</code>

0.2 Readings

There are many excellent books available examining many facets of the least squares problem. Fuller references are in the bibliography.

Carl Freidrich Gauss

Theory of the Combination of Observations Least Subject to Errors

0.3 Least Squares

The titles are ranked by brevity.

Ilse C. F. Ipsen

Numerical Matrix Analysis: Linear System and Least Squares (128 pp)

Charles L. Lawson, and Richard J. Hanson

Solving Least Squares Problems (337 pp)

Åke Björk

Numerical Methods for Least Squares Problems (408 pp)

0.4 Linear Algebra and Matrix Analysis

The titles are ranked by brevity.

Alan J. Laub

Matrix Analysis for Scientists and Engineers (157 pp)

Carl D. Meyer

Matrix Analysis and Applied Linear Algebra (718 pp)

0.5 Numerical Linear Algebra

The titles are ranked by brevity.

Alan J. Laub

Computational Matrix Analysis (154 pp)

Lloyd N. Trefethen, and David Bau, III

Numerical Linear Algebra (361 pp)

E. Anderson, Z. Bai, C. Bischof, S. Blackford, J. Demmel, J. Dongarra, J Du Croz, A. Greenbaum, S. Hammarling, A. McKenney, and D. Sorenson
LAPACK Users' Guide (407 pp)

G. W. Stewart

Matrix Algorithms:

Volume I: Basic Decompositions (460 pp)

Volume II: Eigensystems (474 pp)

0.6 Discussions on Least Squares

Books with chapters dedicated to the topic. Sorted by author.

Gene H. Golub, Charles F. Van Loan

Matrix Computations, ch. 5, 6

Nicholas J. Higham

Accuracy and Stability of Numerical Algorithms, ch. 20

Cleve B. Moler

Numerical Computing with MATLAB, ch. 5

David S. Watkins

Numerical Analysis: a mathematical introduction, ch. 5

David S. Watkins

Fundamentals of Matrix Computations, ch. 3

Part I

Rudiments

Chapter 1

Least Squares Problems

1.1 Linear Systems

This story begins with the archetypal matrix–vector equation

$$\mathbf{A}x = b. \quad (1.1)$$

The matrix \mathbf{A} has m rows, n columns, and has rank ρ ; the vector b encodes m measurements. The solution vector x represents the n free parameters in the model. In mathematical shorthand,

$$\mathbf{A} \in \mathbb{C}_\rho^{m \times n}, \quad b \in \mathbb{C}^m, \quad x \in \mathbb{C}^n \quad (1.2)$$

with \mathbb{C} representing the field of complex numbers. The system matrix \mathbf{A} and the data vector b are given, and the task is to find the solution vector x .

1.1.1 $\|\mathbf{A}x - b\| = 0$

The letters in (1.1) will change, but the operation remains the same: a matrix operates on an n –vector and returns an m –vector. We can think of the matrix as a map from vectors of dimension n to vectors of dimension m :

$$\mathbf{A}: \mathbb{C}^n \mapsto \mathbb{C}^m.$$

If the vector b can be expressed a combination of the columns of the matrix \mathbf{A} then there is a direct solution:

$$\mathbf{A}x = b \implies x_1 a_1 + \cdots + x_n a_n = b$$

and the residual error vanishes:

$$r = \mathbf{A}x - b = \mathbf{0}$$

where the zero vector $\mathbf{0}$ is a list of m zeros. The total error, the norm of the r vector, is 0:

$$\|r\|_2^2 = \|\mathbf{A}x - b\|_2^2 = 0.$$

For the problem where the system matrix \mathbf{A} is the identity matrix \mathbf{I}_2 :

$$\begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} b_1 \\ b_2 \end{bmatrix},$$

the solution is

$$\begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} b_1 \\ b_2 \end{bmatrix};$$

there is no residual error

$$r = \mathbf{A}x - b = \begin{bmatrix} 0 \\ 0 \end{bmatrix}.$$

1.1.2 $\|\mathbf{A}x - b\| > 0$

But what happens when the vector b is *not* in the column space of the matrix \mathbf{A} ? The solution criteria must relax. Instead of seeking zero residual error, seek the least residual error

$$\|r\| = \|\mathbf{A}x - b\|.$$

Instead of a perfect solution, ask for the best solution. One such class of solutions are least squares solutions.

1.2 Least Squares Solutions

In computation of approximations, the goal is to minimize the residual error and this work explores the minimal solutions under the 2-norm, the familiar norm of Pythagoras:

$$\left\| \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} \right\|_2 = \sqrt{x_1^2 + x_2^2}.$$

Let's construct a sample problem with $\|\mathbf{A}x - b\| > 0$ by modifying the previous example:

$$\begin{bmatrix} 1 & 0 \\ 0 & 0 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} b_1 \\ b_2 \end{bmatrix}. \quad (1.3)$$

When $b_2 \neq 0$ there is no exact solution, a solution with $\|\mathbf{A}x - b\| = 0$. Consider the solutions given by

$$x_* = \begin{bmatrix} x_1 \\ 0 \end{bmatrix}; \quad (1.4)$$

the error is $\mathbf{A}x_* - b = -\begin{bmatrix} 0 \\ b_2 \end{bmatrix}$ which has a norm

$$\|r\|_2 = \|\mathbf{A}x_* - b\|_2 = |b_2|, \quad (1.5)$$

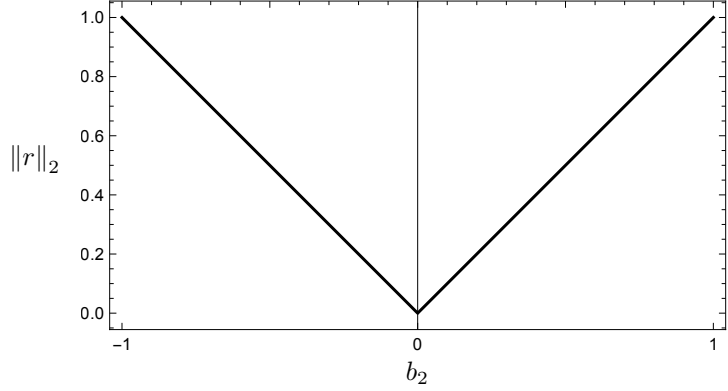


Figure 1.1. The residual error $\|r\|_2$ given in (1.5).

plotted in figure 1.1. This is the least possible error for the problem and (1.4) is the best solution. The transition from an exact solution to an inexact solution is continuous.

Think of the solutions to the linear system

$$\mathbf{A}x = b$$

as being described by the inequality

$$\|\mathbf{A}x - b\|_2 \geq 0.$$

The equality is attained when the data vector lies within the column space of \mathbf{A} .

Least squares solutions are classified by the interpretation of the output. In the first case, *zonal approximation*, the output represents data at a physical zone, a point or a region. In the second case, *modal approximation*, the output represents an amplitude, a contribution for a mode. Basic examples follow.

1.2.1 Zonal Approximation

Consider the vector field F described by the gradient of a scalar field ϕ .

$$F = \nabla\phi$$

The forward problem involves taking a given potential ϕ and computing the forces F . We are instead interested in the inverse problem: measure the forces F and compute the potential ϕ .

Zonal Problem

In practice one measures the vector field and solves the inverse problem by computing values for the potential at the zone boundaries. The input and outputs are represented in 1.2. The physical field is $\phi(x)$, $0 \leq x \leq 2$, the approximation is φ_{x_k} ,

$k = 0, 1, 2$. For a cleaner presentation let $\varphi_{x_k} \rightarrow \varphi_k$. The first measurement x_1 represents the potential change between $\phi(0)$ and $\phi(1)$; the second measurement x_2 the change between $\phi(1)$ and $\phi(2)$.

$$\varphi_1 - \varphi_0 \approx \delta_1$$

$$\varphi_2 - \varphi_1 \approx \delta_2$$

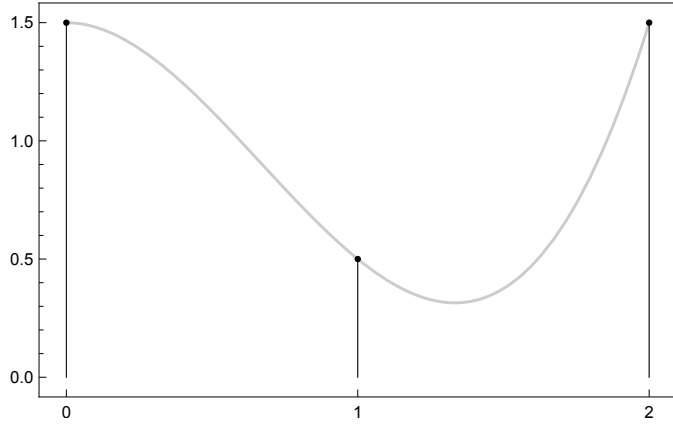


Figure 1.2. *Scalar function $\phi(x)$ (curve) and approximation φ_k (sticks).*

The system matrix

$$\mathbf{A} = \begin{bmatrix} -1 & 1 & 0 \\ 0 & -1 & 1 \end{bmatrix} \in \mathbb{R}_2^{2 \times 3}.$$

There are $m = 2$ measurements, $n = 3$ measurement locations, and the matrix rank is $\rho = 2$. Because the rank is less than the number of columns, $\rho < n$, this problem is *underdetermined*.

The linear system is

$$\begin{bmatrix} -1 & 1 & 0 \\ 0 & -1 & 1 \end{bmatrix} \begin{bmatrix} \varphi_0 \\ \varphi_1 \\ \varphi_2 \end{bmatrix} = \begin{bmatrix} \delta_1 \\ \delta_2 \end{bmatrix}. \quad (1.6)$$

Zonal Solution

The solutions for the linear system in (1.6) which minimize $\|\mathbf{A}x - b\|_2$ are

$$\begin{bmatrix} \varphi_0 \\ \varphi_1 \\ \varphi_2 \end{bmatrix} = \begin{bmatrix} -2 & -1 \\ 1 & -1 \\ 1 & 2 \end{bmatrix} \begin{bmatrix} \delta_1 \\ \delta_2 \end{bmatrix} + \gamma \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix}, \quad \gamma \in \mathbb{C}.$$

The color blue represents range space vectors, red null space vectors. In this way, the fundamental spaces spring to life.

There is a continuum of solutions due to the fact that

$$\mathbf{A} \left(x + \gamma \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \right) = \mathbf{A}x + \mathbf{A} \left(\gamma \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \right) = \mathbf{A}x + \gamma \mathbf{A} \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} = \mathbf{A}x.$$

An quick demonstration confirms the red vector is in the null space:

$$\mathbf{A} \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} = \begin{bmatrix} -1 & 1 & 0 \\ 0 & -1 & 1 \end{bmatrix} \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \end{bmatrix}.$$

1.2.2 Modal Approximation

Modal Problem

In the modal approximation, the user first selects a set of basis functions to describe measurements. Popular basis functions include orthogonal polynomials, trigonometric functions, or monomials. For example, a linear regression implies a basis set of two elements: a constant function, and a linear function. This leads to the familiar equation for a straight line:

$$y(x) = a_0 + a_1x$$

The $n = 2$ parameters represent the intercept (a_0), and the slope (a_1); each of the m measurements represents a straight line:

$$a_0 + a_1x_1 = y_1$$

$$\vdots$$

$$a_0 + a_1x_m = y_m.$$

The goal is to simultaneously solve the set of equations.

Modal Solution

The first step is to compose the system

$$\begin{array}{ccc} \mathbf{A} & \alpha & = & y \\ \begin{bmatrix} 1 & x_1 \\ \vdots & \vdots \\ 1 & x_m \end{bmatrix} & \begin{bmatrix} \alpha_0 \\ \alpha_1 \end{bmatrix} & = & \begin{bmatrix} y_1 \\ \vdots \\ y_m \end{bmatrix}, \end{array} \quad (1.7)$$

which can be expressed using the column vectors

$$\mathbf{1} = \begin{bmatrix} 1 \\ \vdots \\ 1 \end{bmatrix}, \quad x = \begin{bmatrix} x_1 \\ \vdots \\ x_m \end{bmatrix}, \quad y = \begin{bmatrix} y_1 \\ \vdots \\ y_m \end{bmatrix}.$$

The columns of the system matrix $\mathbf{A} = \begin{bmatrix} \mathbf{1} & x \end{bmatrix}$. The solution parameters can be expressed in terms of the column vectors:

$$\begin{bmatrix} \alpha_0 \\ \alpha_1 \end{bmatrix} = \left((\mathbf{1}^T \mathbf{1}) (x^T x) - (\mathbf{1}^T x)^2 \right)^{-1} \begin{bmatrix} x^T x & -\mathbf{1}^T x \\ -\mathbf{1}^T x & \mathbf{1}^T \mathbf{1} \end{bmatrix} \begin{bmatrix} \mathbf{1}^T y \\ x^T y \end{bmatrix}. \quad (1.8)$$

1.2.3 Errors

When measurements are not exact, solutions are not exact. A great beauty of the method of least squares is that the quality of the solution can be quantified. An ability to discern answers like 3.0 ± 1.0 from 3.000 ± 0.0010 is invaluable. The machinery needed to compute uncertainties will be developed in following chapters.

1.3 Least Squares Problem

Emboldened by solutions to two basic problems, we turn attention towards formalities. Starting with a the linear system $\mathbf{A}x = b$ where the matrix $\mathbf{A} \in \mathbb{C}^{m \times n}$, the data vector $b \in \mathbb{C}^m$, the least squares solution x_{LS} is defined as the set

$$x_{LS} = \left\{ x \in \mathbb{C}^n : \|\mathbf{A}x - b\|_2^2 \text{ is minimized} \right\}. \quad (1.9)$$

The least squares solution may be a point or it may be a hyperplane. The general solution is a combination of a particular solution (in blue) and a homogenous solution (in red):

$$\begin{aligned} x_{LS} &= \mathbf{A}^\dagger b + \left(\mathbf{I}_n - \mathbf{A}^\dagger \mathbf{A} \right) y, & y \in \mathbb{C}^n \\ &= x_{\dagger} + x_{\mathcal{N}} \end{aligned} \quad (1.10)$$

where the matrix \mathbf{A}^\dagger is the pseudoinverse and y is an arbitrary vector.

Chapter 2

Least Squares Solutions

Bolstered from producing concrete results, attention now turns to an examination of solution methods through the lens of the Fundamental Theorem.

2.1 Fundamental Theorem of Linear Algebra

The Fundamental Theorem is a beautiful and powerful statement which helps to organize techniques and results in linear algebra. It is the stage upon which the stories of existence and uniqueness play out.

An exemplar matrix,

$$\mathbf{A} = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix} \in \mathbb{C}^{3 \times 4},$$

demonstrates salient points with an immediate subspace decomposition. The matrix \mathbf{A} maps 4-vectors from the domain \mathbb{C}^4 to 3-vectors in the codomain \mathbb{C}^3 . One way to express the Fundamental Theorem is to say that the matrix \mathbf{A} induces a 4 dimensional column space and a 3 dimensional row space.

For this case, the domain can be described as

$$\mathbb{C}^4 = \text{sp} \left\{ \begin{bmatrix} 1 \\ 0 \\ 0 \\ 0 \end{bmatrix}, \begin{bmatrix} 0 \\ 1 \\ 0 \\ 0 \end{bmatrix} \right\} \oplus \text{sp} \left\{ \begin{bmatrix} 0 \\ 0 \\ 1 \\ 0 \end{bmatrix}, \begin{bmatrix} 0 \\ 0 \\ 0 \\ 1 \end{bmatrix} \right\},$$

and the codomain as

$$\mathbb{C}^3 = \text{sp} \left\{ \begin{bmatrix} 1 \\ 0 \\ 0 \end{bmatrix}, \begin{bmatrix} 0 \\ 1 \\ 0 \end{bmatrix} \right\} \oplus \text{sp} \left\{ \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix} \right\}$$

The symbol \oplus specifies that the vectors spaces for the range and null space are orthogonal.

There are many statements of the Fundamental Theorem. One way describes the decomposition of the domain and codomain into the four fundamental subspaces: There are many statements of the Fundamental Theorem. One way describes the decomposition of the domain and codomain into the four fundamental subspaces: Each domain space, domain and codomain, are expressed as an orthogonal decom-

Table 2.1. *The Fundamental Theorem of Linear Algebra for $\mathbf{A} \in \mathbb{C}^{m \times n}$*

$$\begin{array}{llll} \text{domain: } \mathbb{C}^n & = & \mathcal{R}(\mathbf{A}^*) & \oplus & \mathcal{N}(\mathbf{A}) \\ \text{codomain: } \mathbb{C}^m & = & \mathcal{R}(\mathbf{A}) & \oplus & \mathcal{N}(\mathbf{A}^*) \end{array}$$

position of a range space and a null space.

Table 2.2. The Fundamental Theorem of Linear Algebra and Least Squares for $A \in \mathbb{C}^{m \times n}$

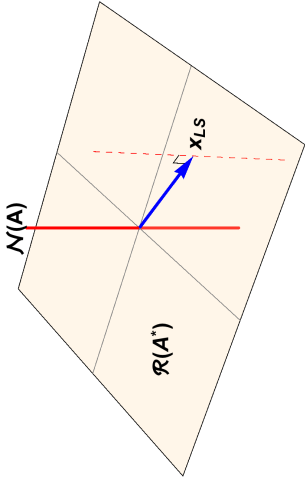
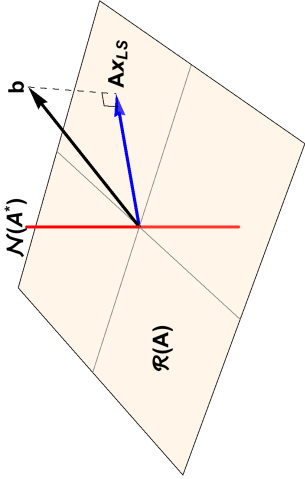
Domain	Mapping			Codomain
	$A: \mathbb{C}^n$	\mapsto	\mathbb{C}^m	
	\mathbb{C}^n	\leftarrow	$\mathbb{C}^m: A^*$	
$\mathbb{C}^n = \mathcal{R}(A^*) \oplus \mathcal{N}(A)$				
$\mathbb{C}^m = \mathcal{R}(A) \oplus \mathcal{N}(A^*)$				

Table 2.3. *Dimensions of the fundamental subspaces for $\mathbf{A} \in \mathbb{C}_\rho^{m \times n}$.*

$$\begin{array}{llll} \dim(\mathcal{R}(\mathbf{A})) & = & \rho & \dim(\mathcal{N}(\mathbf{A}^*)) & = & m - \rho \\ \dim(\mathcal{R}(\mathbf{A}^*)) & = & \rho & \dim(\mathcal{N}(\mathbf{A})) & = & n - \rho \end{array}$$

Example from (1.6)

$$\begin{bmatrix} -1 & 1 & 0 \\ 0 & -1 & 1 \end{bmatrix}$$

Begin with the column space, also known as the solution space because the solution parameters will inhabit this space:

$$\text{Column vectors} = \begin{bmatrix} -1 \\ 0 \end{bmatrix}, \begin{bmatrix} 1 \\ -1 \end{bmatrix}, \begin{bmatrix} 0 \\ 1 \end{bmatrix}.$$

The column vectors are a complete span of \mathbb{C}^2 , and the null space is trivial:

$$\mathbb{C}^2 = \text{sp} \left\{ \begin{bmatrix} -1 \\ 0 \end{bmatrix}, \begin{bmatrix} 1 \\ -1 \end{bmatrix}, \begin{bmatrix} 0 \\ 1 \end{bmatrix} \right\} \oplus \begin{bmatrix} 0 \\ 0 \end{bmatrix}$$

A minimal span for \mathbb{C}^2 is

$$\mathbb{C}^2 = \text{sp} \left\{ \begin{bmatrix} 1 \\ 0 \end{bmatrix}, \begin{bmatrix} 0 \\ 1 \end{bmatrix} \right\}.$$

Next up is the row space, or the measurement space

$$\text{Row vectors} = \begin{bmatrix} -1 \\ 1 \\ 0 \end{bmatrix}, \begin{bmatrix} 0 \\ -1 \\ 1 \end{bmatrix}$$

$$\mathbb{C}^3 = \text{sp} \left\{ \begin{bmatrix} -1 \\ 1 \\ 0 \end{bmatrix}, \begin{bmatrix} 0 \\ -1 \\ 1 \end{bmatrix} \right\} \oplus \text{sp} \left\{ \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \right\}$$

2.2 Singular Value Decomposition - I

The Fundamental Theorem describes the world as a orthogonal decomposition of the domain and codomain. Why not ask for an orthonormal decomposition? This is precisely what we get from the singular value decomposition. As noted by Gene Golub, the singular value decomposition is a singularly valuable decomposition. It not only finds an orthonormal basis for the fundamental subspaces, it also aligns the domain and codomain, and furthermore, resolves the scale factors needed to move between spaces. This first presentation offers a statement of the SVD theorem.

2.2.1 SVD Theorem

Given a matrix $\mathbf{A} \in \mathbb{C}_\rho^{m \times n}$, a matrix with complex entries with m rows, n columns, and matrix rank $0 < \rho \leq \min(m, n)$, then there exists a decomposition of the form

$$\mathbf{A} = \mathbf{U} \Sigma \mathbf{V}^*$$

where

1. column vectors of unitary matrix $\mathbf{V} \in \mathbb{C}^{n \times n}$ represent an orthonormal span of the domain,
2. column vectors of unitary matrix $\mathbf{U} \in \mathbb{C}^{m \times m}$ represent an orthonormal span of the codomain,
3. Diagonal entries of $\Sigma \in \mathbb{R}^{m \times n}$ contain the singular values; the ordered, nonzero eigenvalues of the product matrix $\mathbf{A}^* \mathbf{A}$.

In block form

$$\mathbf{A} = \mathbf{U} \Sigma \mathbf{V}^* = \left[\begin{array}{c|c} \mathbf{U}_{\mathcal{R}} & \mathbf{U}_{\mathcal{N}} \end{array} \right] \left[\begin{array}{c|c} \mathbf{S} & \mathbf{0} \\ \hline \mathbf{0} & \mathbf{0} \end{array} \right] \left[\begin{array}{c} \mathbf{V}_{\mathcal{R}}^* \\ \hline \mathbf{V}_{\mathcal{N}}^* \end{array} \right]$$

Column vectors span the subspaces:

$$\begin{aligned} \mathbf{V} &= \left[\begin{array}{ccc|ccc} v_1 & \dots & v_\rho & v_{\rho+1} & \dots & v_n \end{array} \right], \\ \mathbf{U} &= \left[\begin{array}{ccc|ccc} u_1 & \dots & u_\rho & u_{\rho+1} & \dots & u_m \end{array} \right]. \end{aligned}$$

$$\mathbf{U} \in \mathbb{C}^{m \times m},$$

$$\mathbf{V} \in \mathbb{C}^{n \times n},$$

$$\Sigma \in \mathbb{R}^{m \times n}.$$

Table 2.4. Orthonormal spans for the invariant subspaces.

domain			codomain		
$\mathcal{R}(\mathbf{A}^*)$	=	$\text{span}\{v_1, \dots, v_\rho\}$	$\mathcal{R}(\mathbf{A})$	=	$\text{span}\{u_1, \dots, u_\rho\}$
$\mathcal{N}(\mathbf{A})$	=	$\text{span}\{v_{\rho+1}, \dots, v_n\}$	$\mathcal{N}(\mathbf{A}^*)$	=	$\text{span}\{u_{\rho+1}, \dots, u_m\}$

$$u_j \cdot u_k = \delta_{jk},$$

$$v_j \cdot v_k = \delta_{jk}.$$

Decomposition for (1.6):

$$\mathbf{A} = \mathbf{U} \Sigma \mathbf{V}^*$$

$$\begin{bmatrix} -1 & 1 & 0 \\ 0 & -1 & 1 \end{bmatrix} = \frac{1}{\sqrt{2}} \begin{bmatrix} -1 & 1 \\ 1 & 1 \end{bmatrix} \begin{bmatrix} \sqrt{3} & 0 & 0 \\ 0 & 1 & 0 \end{bmatrix} \begin{bmatrix} \frac{1}{\sqrt{6}} & -\frac{2}{\sqrt{6}} & \frac{1}{\sqrt{6}} \\ -\frac{1}{\sqrt{2}} & 0 & \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{3}} & \frac{1}{\sqrt{3}} & \frac{1}{\sqrt{3}} \end{bmatrix}$$

$$\mathbf{S} = \begin{bmatrix} \sqrt{3} & 0 \\ 0 & 1 \end{bmatrix}$$

2.2.2 SVD and Least Squares

A direct implication of the singular value decomposition is the homogeneous solution.

Unitary transformation

The definition of the least squares problem in (1.9) shows that the target of minimization is the quantity

$$r^* r = r^2 = \|\mathbf{A}x - b\|_2^2.$$

One minimization strategy invokes a unitary transformation to create a simpler problem:

$$\|\mathbf{A}x - b\|_2^2 = \|\mathbf{U}^* (\mathbf{A}x - b)\|_2^2. \quad (2.1)$$

This remarkable insight opens a door to solution. Rearranging the singular value decomposition

$$\mathbf{U}^* \mathbf{A} = \Sigma \mathbf{V}^*,$$

and using the block form in (2.2.1) leads to

$$\begin{aligned} \|\mathbf{A}x - b\|_2^2 &= \|\Sigma \mathbf{V}^* x - \mathbf{U}^* b\|_2^2 = \left\| \begin{bmatrix} \mathbf{S} & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix} \begin{bmatrix} \mathbf{V}_{\mathcal{R}}^* \\ \mathbf{V}_{\mathcal{N}}^* \end{bmatrix} x - \begin{bmatrix} \mathbf{U}_{\mathcal{R}}^* \\ \mathbf{U}_{\mathcal{N}}^* \end{bmatrix} b \right\|_2^2 \\ &= \left\| \begin{bmatrix} \mathbf{S} \mathbf{V}_{\mathcal{R}}^* x \\ \mathbf{0} \end{bmatrix} - \begin{bmatrix} \mathbf{U}_{\mathcal{R}}^* b \\ \mathbf{U}_{\mathcal{N}}^* b \end{bmatrix} \right\|_2^2. \end{aligned}$$

The range space components are now untangled from the null space components.

Pseudoinverse solution

Using the Pythagorean theorem to isolate the range and null space components of the total error for the least squares problem

$$\|\mathbf{A}x - b\|_2^2 = \underbrace{\|\mathbf{S} \mathbf{V}_{\mathcal{R}}^* x - \mathbf{U}_{\mathcal{R}}^* b\|_2^2}_{\substack{x \text{ dependence} \\ \text{under control}}} + \underbrace{\|\mathbf{U}_{\mathcal{N}}^* b\|_2^2}_{\substack{\text{residual} \\ \text{no control}}}$$

There are now two terms; the first depends upon the solution vector x , the second does not. We only have control over the first term. To minimize the total error we must drive the first term to zero. Then the total error will be given by the residual error term. The error term that is controlled by the solution vector x is this

$$\mathbf{S}\mathbf{V}_{\mathcal{R}}^*x - \mathbf{U}_{\mathcal{R}}^*b. \quad (2.2)$$

Choosing the vector x which forces this term to zero leads to the SVD solution for the least squares problem:

$$x_{\dagger} = \mathbf{V}_{\mathcal{R}}\mathbf{S}^{-1}\mathbf{U}_{\mathcal{R}}^*b.$$

This is also the pseudoinverse solution

$$x_{\dagger} = \mathbf{A}^{\dagger}b$$

where the (thin) pseudoinverse is

$$\mathbf{A}^{\dagger} = \mathbf{V}_{\mathcal{R}}\mathbf{S}^{-1}\mathbf{U}_{\mathcal{R}}^*.$$

The error that can be controlled is forced to 0; but this leaves an error which cannot be removed, a residual error defined as

$$r^2 = \|\mathbf{U}_{\mathcal{N}}^*b\|_2^2.$$

The usually silent null space term can be heard as it pronounces the value of the total error.

To recap, the singular value decomposition leads immediately to the pseudoinverse solution and residual error.

In retrospect

We can reimagine the least squares problem as the challenge of resolving the data vector into range and null space components:

$$b = b_{\mathcal{R}} + b_{\mathcal{N}}.$$

Such an example is shown in (4.2).

$$\|\mathbf{A}x - b\|_2^2 = \|\mathbf{A}x - b_{\mathcal{R}} - b_{\mathcal{N}}\|_2^2 = \|b_{\mathcal{N}}\|_2^2$$

Because the vector $b_{\mathcal{R}} \in \mathcal{R}(\mathbf{A})$, there exists a vector x such that $\mathbf{A}x = b_{\mathcal{R}}$. Again, the error that cannot be removed is the residual error

$$\|b_{\mathcal{N}}\|_2^2$$

What we shown is that the vector x which minimizes the least squares error in (??) is exactly the same vector given by the SVD solution in equation (2.2.2). Using a unitary transform we were able to convert the general least squares problem into a form amenable to solution using the singular value decomposition.

For the overdetermined case as we have here the usually silent null space term can be heard as it pronounces the value of the total error

$$r^2 = \|\mathbf{U}_{\mathcal{N}}^*b\|_2^2 = (\mathbf{U}_{\mathcal{N}}^*b)^* (\mathbf{U}_{\mathcal{N}}^*b) = b^* (\mathbf{U}_{\mathcal{N}}\mathbf{U}_{\mathcal{N}}^*) b, \quad (2.3)$$

a restatement of (2.2.2).

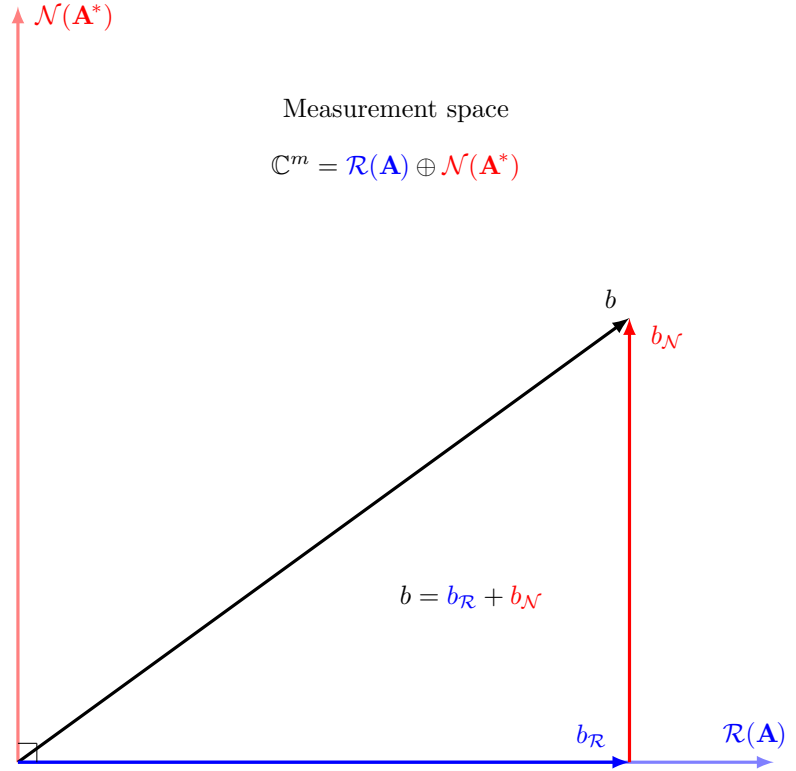


Figure 2.1. Decomposing the data vector.

2.3 Singular Value Decomposition - II

2.3.1 Σ Gymnastics

Success in manipulating the singular value decomposition includes builds upon success in manipulating the Σ matrices. Think of the Σ matrix as a sabot matrix for the matrix of singular values \mathbf{S} . The breakdown of block diagrams: the matrix Σ ($m \times n$) and the matrix Σ^T ($n \times m$) have different shapes, but equivalent block diagrams.

$$\Sigma = \begin{bmatrix} \mathbf{S} & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix},$$

$$\Sigma^T = \begin{bmatrix} \mathbf{S}^T & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix} = \begin{bmatrix} \mathbf{S} & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix},$$

$$\Sigma^\dagger = \begin{bmatrix} \mathbf{S}^{-1} & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix}.$$

$$\Sigma = \left[\begin{array}{c|c} \mathbf{S}_{\rho \times \rho} & \mathbf{0}_{\rho \times n-\rho} \\ \hline \mathbf{0}_{m-\rho \times \rho} & \mathbf{0}_{m-\rho \times n-\rho} \end{array} \right]$$

Notice the interchange of the indices m and n in the transpose:

$$\Sigma^T = \left[\begin{array}{c|c} \mathbf{S}_{\rho \times \rho} & \mathbf{0}_{\rho \times m-\rho} \\ \hline \mathbf{0}_{n-\rho \times \rho} & \mathbf{0}_{n-\rho \times m-\rho} \end{array} \right].$$

The pseudoinverse matrix has the same dimension as the parent matrix:

$$\Sigma = \left[\begin{array}{c|c} \mathbf{S}_{\rho \times \rho} & \mathbf{0}_{\rho \times n-\rho} \\ \hline \mathbf{0}_{m-\rho \times \rho} & \mathbf{0}_{m-\rho \times n-\rho} \end{array} \right]$$

$$\Sigma \Sigma^\dagger = \mathbb{I}_{m,\rho},$$

$$\Sigma^\dagger \Sigma = \mathbb{I}_{n,\rho}.$$

$$\Sigma = \begin{bmatrix} \sigma_1 & 0 \\ 0 & \sigma_2 \\ 0 & 0 \\ 0 & 0 \end{bmatrix}, \quad \Sigma^T = \begin{bmatrix} \sigma_1 & 0 & 0 & 0 \\ 0 & \sigma_2 & 0 & 0 \end{bmatrix},$$

$$\Sigma^\dagger = \begin{bmatrix} 1/\sigma_1 & 0 & 0 & 0 \\ 0 & 1/\sigma_2 & 0 & 0 \end{bmatrix}.$$

Stencil matrices

$$\Sigma \Sigma^\dagger = \mathbb{I}_{4,2} = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix} = \begin{bmatrix} \mathbf{I}_2 & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix},$$

$$\Sigma^\dagger \Sigma = \mathbb{I}_{2,2} = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} = \mathbf{I}_2.$$

Product matrices with the transpose

$$\Sigma \Sigma^T = \begin{bmatrix} \sigma_1^2 & 0 & 0 & 0 \\ 0 & \sigma_2^2 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix} = \begin{bmatrix} \mathbf{S}^2 & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix},$$

$$\Sigma^T \Sigma = \begin{bmatrix} \sigma_1^2 & 0 \\ 0 & \sigma_2^2 \end{bmatrix} = \mathbf{S}^2.$$

Product matrices with the pseudoinverse

$$\begin{aligned}\Sigma \Sigma^\dagger &= \mathbb{I}_{2,2} = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} = \mathbf{I}_2, \\ \Sigma \Sigma^T &= \begin{bmatrix} \sigma_1^2 & 0 \\ 0 & \sigma_2^2 \end{bmatrix} = \mathbf{S}^2, \\ \Sigma^\dagger \Sigma &= \mathbb{I}_{4,2} = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix} = \begin{bmatrix} \mathbf{I}_2 & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix}, \\ \Sigma^T \Sigma &= \begin{bmatrix} \sigma_1^2 & 0 & 0 & 0 \\ 0 & \sigma_2^2 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix} = \begin{bmatrix} \mathbf{S}^2 & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix},\end{aligned}$$

2.3.2 Fundamental Projectors

Given a matrix $\mathbf{A} \in \mathbb{C}_\rho^{m \times n}$, a matrix with complex entries with m rows, n columns, and matrix rank $0 < \rho \leq \min(m, n)$, then there exists a

Table 2.5. *Fundamental Projectors using the pseudoinverse.*

	Range space		Null space	
domain	$\mathbf{P}_{\mathcal{R}(\mathbf{A}^*)}$	$= \mathbf{A}^\dagger \mathbf{A}$	$\mathbf{P}_{\mathcal{N}(\mathbf{A})}$	$= \mathbf{I}_n - \mathbf{A}^\dagger \mathbf{A}$
codomain	$\mathbf{P}_{\mathcal{R}(\mathbf{A})}$	$= \mathbf{A} \mathbf{A}^\dagger$	$\mathbf{P}_{\mathcal{N}(\mathbf{A}^*)}$	$= \mathbf{I}_m - \mathbf{A} \mathbf{A}^\dagger$

Table 2.6. *Fundamental Projectors using domain matrices.*

	Range space		Null space	
domain	$\mathbf{P}_{\mathcal{R}(\mathbf{A}^*)}$	$= \mathbf{V}_{\mathcal{R}} \mathbb{I}_{n,\rho} \mathbf{V}_{\mathcal{R}}^*$	$\mathbf{P}_{\mathcal{N}(\mathbf{A})}$	$= \mathbf{I}_n - \mathbf{V}_{\mathcal{R}} \mathbb{I}_{n,\rho} \mathbf{V}_{\mathcal{R}}^*$
codomain	$\mathbf{P}_{\mathcal{R}(\mathbf{A})}$	$= \mathbf{U}_{\mathcal{R}} \mathbb{I}_{m,\rho} \mathbf{U}_{\mathcal{R}}^*$	$\mathbf{P}_{\mathcal{N}(\mathbf{A}^*)}$	$= \mathbf{I}_m - \mathbf{U}_{\mathcal{R}} \mathbb{I}_{m,\rho} \mathbf{U}_{\mathcal{R}}^*$

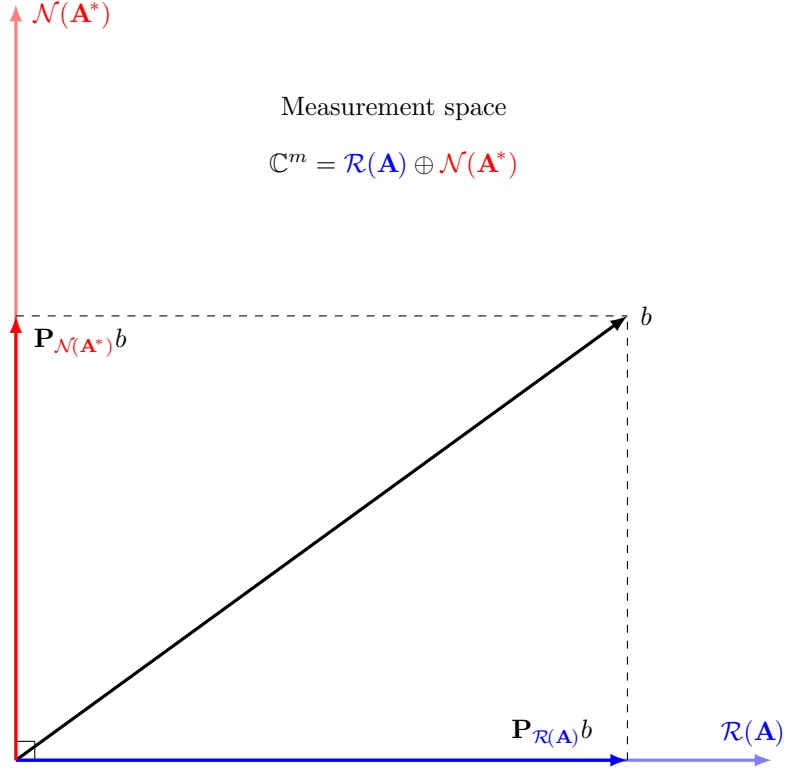


Figure 2.2. Projections of the data vector.

2.4 Least Squares Solution - Again

Let's revisit the canonical linear system in (1.1) the general solution in (1.10):

$$\begin{aligned} x_{LS} &= \mathbf{A}^\dagger b + (\mathbf{I}_n - \mathbf{A}^\dagger \mathbf{A}) y \\ &= \mathbf{A}^\dagger b + \mathbf{P}_{\mathcal{R}(\mathbf{A}^*)} y \end{aligned}$$

where the arbitrary vector $y \in \mathbb{C}^n$.

The projector onto the range space $\mathcal{R}(\mathbf{A}^*)$

$$\mathbf{A}^\dagger \mathbf{A} = \mathbf{V} \Sigma^\dagger \Sigma \mathbf{V}^* = \mathbf{V}_{\mathcal{R}} \mathbb{I}_{\rho, m} \mathbf{V}_{\mathcal{R}}^*$$

2.5 Reflection Invariance

Least squares solutions have a reflection invariance based upon the arithmetic fact

$$\sum r_2^2 = \sum (-r)^2.$$

Starting with the canonical linear system in (1.1), the least squares solution is x_{LS} and the residual errors may be defined as

$$r = \mathbf{A}x - b = r \quad \Rightarrow \quad r = \mathbf{A}^\dagger b - b = (\mathbf{A}^\dagger - \mathbf{I}_m) T.$$

Notice that $-r = b - \mathbf{A}^\dagger b$. To reflect the data points through the solution curve us

$$\tilde{b} = b + 2r$$

$$r = b - \mathbf{A}x$$

Two equivalent ways to define the residual vector

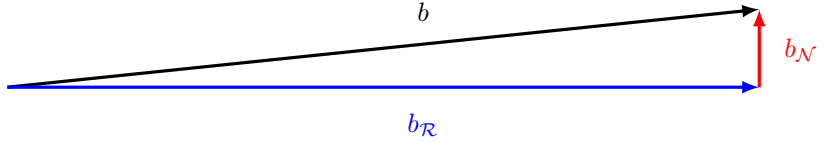


Figure 2.3. Data vector resolved as $b = b_{\mathcal{R}} + b_{\mathcal{N}}$.

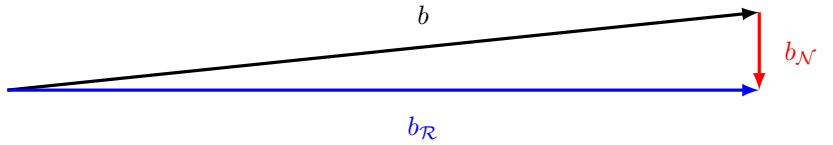


Figure 2.4. Data vector resolved as $b + b_{\mathcal{N}} = b_{\mathcal{R}}$.

$$x_{LS} = x_R$$

$$(\mathbf{A}^* \mathbf{A})^{-1} \mathbf{A}^* b = (\mathbf{A}^* \mathbf{A})^{-1} \mathbf{A}^* (b - 2r)$$

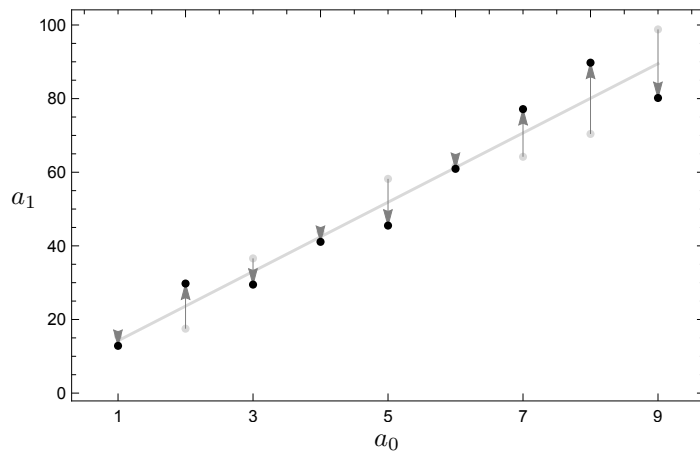
True iff $(\mathbf{A}^* \mathbf{A})^{-1} \mathbf{A}^* r = 0$:

$$\begin{aligned} (\mathbf{A}^* \mathbf{A})^{-1} \mathbf{A}^* r &= (\mathbf{A}^* \mathbf{A})^{-1} \mathbf{A}^* (\mathbf{A}x - b) \\ &= \left((\mathbf{A}^* \mathbf{A})^{-1} \mathbf{A}^* \mathbf{A} \right) x - (\mathbf{A}^* \mathbf{A})^{-1} \mathbf{A}^* b \\ &= x - x \\ &= \mathbf{0}. \end{aligned}$$

This test can be a filter as seen in §13.3.3

Table 2.7. *Reflecting the data for $\mathbf{A}x - T = r$.*

method	original data	reflected data
I	T	$T + 2r$
II	$\mathbf{A}^\dagger a - r$	$\mathbf{A}^\dagger a + r$

**Figure 2.5.** *The merit function for the learning curve showing the minimum and the value.*

Part II

Modal Example

Solving for the modes in a expansion basis. Here linear regression.

Chapter 3

Solution Via Calculus

3.1 Modal Approximation

The following example represents a problem in linear regression. A sequence of m data points (x_k, T_k) , $k = 1:m$ is recorded. The goal is to find the best approximation to a straight line. The *trial function* is

$$T(x) = a_0 + a_1x.$$

The residual errors are the difference between the measurements and predictions:

$$\text{residual error}_k = \underbrace{\text{measurement}_k}_{T_k} - \underbrace{\text{prediction}_k}_{T(x_k)}$$

Formally, the residual error is

$$r_k = T_k - T(x_k), \quad k = 1:m.$$

From this springs the *merit function*, the target of minimization,

$$\begin{aligned} M(a) &= \sum_{k=1}^m r_k^2 \\ &= \sum_{k=1}^m (\text{measurement}_k - \text{prediction}_k)^2 \\ &= \sum_{k=1}^m (T_k - T(x_k))^2 \\ &= \sum_{k=1}^m (T_k - a_0 - a_1x_k)^2 \end{aligned} \tag{3.1}$$

The least squares solution a_{LS} is defined as

$$a_{LS} = \left\{ a \in \mathbb{C}^2 : \|T(x_k) - a_0 - a_1x_k\|_2^2 \text{ is minimized} \right\}.$$

Colloquially, the least squares solution is the set a of complex pairs such that the square of the norm $\|T(x_k) - a_0 - a_1 x_k\|_2^2$ is minimized.

The solution set satisfies

$$\nabla M(a)|_{a_{LS}} = 0. \quad (3.2)$$

3.2 Bevington Example

To provide a common reference, see the example in Bevington [2, ch 6]. The data is summarized below in table 3.2. The problem involves temperature measurements T_k made at position x_k on a bar in contact with two heat baths (Dirichlet boundary conditions). A conceptualization is shown in figure 3.1. Arrowheads on the bottom show the nine locations where the temperature is measured.

In the ideal linear case, the temperature at the endpoints matches the temperature of the baths, $T(x = 0 \text{ cm}) = 0^\circ\text{C}$ and $T(x = 10 \text{ cm}) = 100^\circ\text{C}$ which describes a line with intercept $a_0 = 0^\circ\text{C}$ and slope $a_1 = 10^\circ\text{C/cm}$. Such an expectation is a crude quality measure, a “sniff test”.



Figure 3.1. *Measuring the temperature of a bar held between two constant-temperature heat baths.*

3.2.1 Problem Statement

Muddled conceptions are wellsprings for muddled execution. Success in dealing with complicated problems in least squares flows from being able to state the problem cleanly; a good practice is to start with a table specifying the problem of interest, such as table 13.4.

The first entry is the *trial function* which defines the functional form to be applied to the data. As the name indicates, this function is an initial guess. Whether or not Nature has chosen this model remains to be seen.

The *merit function* is created by inserting the trial function into (3.1). This function is the target of minimization and can provide a crude check on the solution. Given a candidate solution a_* , compute the value of $M(a_*)$. The least squares solution has the property that $M(a_*)$ has minimum value in the neighborhood of a_* . If the solution is perturbed, one must have $M(a_*) < M(a_* + \delta a)$. When you are developing and refining least squares algorithms, you may see that the computed solution a_* changes. For overdetermined problems, the solution is unique

and comparing the values of the merit function helps discriminate solutions. In a later section figure ?? will demonstrate this behavior.

The *measurements* define the quantities to be measured. It seems an obvious step, but more complex models may have ambiguities start here.

Results, or fit parameters, define the quantities to be computed using the least squares algorithm. Together with the trial function, and the measurements, we now have a clear idea of what will be measured and what will be computed.

The *residual error* specifies the difference between measurement and prediction at each point. A simple matter, apparent in the merit function, it is helpful to write it out, particularly for those who may be using the results and not intimate with the derivation.

The *system matrix* describes the measurement apparatus and contains the dependent variables. In this example we have $m = 9$ rows (measurements), $n = 2$ columns (fit parameters) and a matrix rank $\rho = 2$ (full column rank and overdetermined).

The *linear system* shows the application of the trial function to every measurement. It's a good idea to keep this image in mind.

The *ideal solution* is an infrequent visitor which helps provide a rough measure of quality. Caution is required, though. The ideal solution typically represents a concatenation of miracles which Nature may avoid. In this example, the ideal solution assumes magic barriers which absorb no heat, a bar of exact length, thermometers with exact measurements, heat baths at exact temperatures, no interaction with the local environment, etc. The hope is that systematic effects will be negligible and random effects will have 0 mean.

The next phase is to gather and record the data as shown in table 3.2. Discussion of significant digits in the input data is deferred.

3.2.2 Normal Equations via Calculus

In §6.4, Bevington solves the problem by applying calculus to the final form in (3.1), effectively solving (3.2). Introducing the notation

$$\partial_j M(a_0, a_1) = \frac{\partial M(a_0, a_1)}{\partial a_j}$$

the simultaneous equations to solve are

$$\begin{aligned}\partial_0 M(a_0, a_1) &= 0, \\ \partial_1 M(a_0, a_1) &= 0.\end{aligned}\tag{3.3}$$

These two differential equations spawn the linear system

$$\begin{aligned}-2 \sum_{k=1}^m (T_k - a_0 - a_1 x_k) &= 0, \\ -2 \sum_{k=1}^m (T_k - a_0 - a_1 x_k) x_k &= 0.\end{aligned}$$

Table 3.1. *Problem statement for linear regression.*

trial function	$T(x) = a_0 + a_1x$	$^{\circ}\text{C}$
residual error	$r_k = T_k - a_0 - a_1x$	$^{\circ}\text{C}$
merit function	$M(a) = \sum_{k=1}^m (T_k - a_0 - a_1x)^2$	$(^{\circ}\text{C})^2$
measurements	$x_k, k = 1:m$	position, cm
	$T_k, k = 1:m$	temperature, $^{\circ}\text{C}$
results	$a_0 \pm \epsilon_0$	intercept, $^{\circ}\text{C}$
	$a_1 \pm \epsilon_1$	slope, $^{\circ}\text{C} / \text{cm}$
# of measurements	$m = 9$	rows in A
# of parameters	$n = 2$	columns in A
system matrix	$\mathbf{A} \in \mathbb{R}_2^{9 \times 2}$	
linear system	$\begin{bmatrix} 1 & x_1 \\ \vdots & \vdots \\ 1 & x_m \end{bmatrix} \begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \begin{bmatrix} T_1 \\ \vdots \\ T_m \end{bmatrix}$	
ideal solution	$\begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \begin{bmatrix} 0 \\ 10 \end{bmatrix}$	
input data	table 3.2	

Distributing the summation operators creates a more revealing form

$$\begin{aligned} \sum_{k=1}^m T_k &= a_0 \sum 1 + a_1 \sum x_k, \\ \sum_{k=1}^m T_k x_k &= a_0 \sum x_k + a_1 \sum x_k^2, \end{aligned}$$

where summation from 1 to m is implied. (Therefore $\sum 1 = m$.) The minimization criteria are now recast as the linear system

$$\begin{bmatrix} \sum 1 & \sum x_k \\ \sum x_k & \sum x_k^2 \end{bmatrix} \begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \begin{bmatrix} \sum T_k \\ \sum T_k x_k \end{bmatrix}. \quad (3.4)$$

The solution can be written immediately. Defining the determinant

$$\Delta = m \sum x_k^2 - \left(\sum x_k \right)^2, \quad (3.5)$$

the matrix inverse is

$$\begin{bmatrix} m & \sum x_k \\ \sum x_k & \sum x_k^2 \end{bmatrix}^{-1} = \Delta^{-1} \begin{bmatrix} \sum x_k^2 & -\sum x_k \\ -\sum x_k & m \end{bmatrix}. \quad (3.6)$$

Table 3.2. *Raw data and results.*

k	Input		Output	
	(cm)	(°C)	(°C)	(°C)
	x_k	T_k	$T(x_k)$	r_k
1	1	15.6	14.2222	-1.37778
2	2	17.5	23.6306	6.13056
3	3	36.6	33.0389	-3.56111
4	4	43.8	42.4472	-1.35278
5	5	58.2	51.8556	-6.34444
6	6	61.6	61.2639	-0.336111
7	7	64.2	70.6722	6.47222
8	8	70.4	80.0806	9.68056
9	9	98.8	89.4889	-9.31111

The solution to equation (3.4) is the matrix product

$$\begin{aligned}
 \begin{bmatrix} a_0 \\ a_1 \end{bmatrix} &= \Delta^{-1} \begin{bmatrix} \sum x_k^2 & -\sum x_k \\ -\sum x_k & \sum 1 \end{bmatrix} \begin{bmatrix} \sum T_k \\ \sum T_k x_k \end{bmatrix}, \\
 a_0 &= \Delta^{-1} \left(\sum x_k^2 \sum T_k - \sum x_k \sum T_k x_k \right), \\
 a_1 &= \Delta^{-1} \left(m \sum T_k x_k - \sum x_k \sum T_k \right).
 \end{aligned} \tag{3.7}$$

The final results agree with Bevington's equations 6–17:

Bevington's §6–5 is a succinct explanation of error propagation. In short, measurements are inexact, therefore results will be inexact. The beauty of the method of least squares is that the error in the solution parameters can be expressed in terms of the error in the data. Measurements without uncertainties are incomplete measurements.

The computation chain begins with an estimate of the parent standard deviation which is based upon the total error:

$$s^2 \approx \frac{r^* r}{m - n}.$$

Error contributions for individual parameters are harvested from the diagonal elements of the matrix inverse in (3.6):

$$\begin{aligned}
 \epsilon_0^2 &= \frac{r^T r}{\Delta(m - n)} \sum x_k^2 \\
 \epsilon_1^2 &= \frac{r^T r}{\Delta(m - n)} \sum 1
 \end{aligned}$$

Table 3.3. *Results for linear regression.*

fit parameters	a_0	intercept, °C
	a_1	slope, °C / cm
solution function	$T(x) = a_0 + a_1x$	°C
solution error	$\epsilon_T^2(x) = \epsilon_0^2 + x^2\epsilon_1^2 + a_1^2\epsilon_x^2$	°C
computed solution	$\begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \begin{bmatrix} 4.8 \\ 9.41 \end{bmatrix} \pm \begin{bmatrix} 4.9 \\ 0.87 \end{bmatrix}$	
ideal solution	$\begin{bmatrix} \tilde{a}_0 \\ \tilde{a}_1 \end{bmatrix} = \begin{bmatrix} 0 \\ 10 \end{bmatrix}$	
r^*r	316.6	
curvature matrix $(\mathbf{A}^*\mathbf{A})^{-1}$	$\frac{1}{180} \begin{bmatrix} 95 & -15 \\ -15 & 3 \end{bmatrix}$	
problem statement	table 13.4	
input data	table 3.2	
plots	figure 3.2	1. data & solution
	figure 3.3	2. residual errors
	figure ??	3. merit function

3.3 Numerical Results

Results are stated in two forms. The first is an integer form free of numerical errors inherent in binary representations with finite length. This liberates one from trying to determine if errors are in the algorithm or in machine arithmetic. To aid debugging, intermediate results are also provided.

The second form represents the answer which would be provided to a customer: the fit parameters and associated errors quoted with the proper amount of significant digits.

3.3.1 Exact Form

Exact results for the fit parameters are error follow. The product matrix in (3.4) is

$$\begin{bmatrix} m & \sum x_k \\ \sum x_k & \sum x_k^2 \end{bmatrix} = \begin{bmatrix} 9 & 45 \\ 45 & 285 \end{bmatrix},$$

with determinant

$$\Delta = 540.$$

The inverse of this matrix, (3.6), is

$$\begin{bmatrix} m & \sum x_k \\ \sum x_k & \sum x_k^2 \end{bmatrix}^{-1} = \Delta^{-1} \begin{bmatrix} 285 & -45 \\ -45 & 9 \end{bmatrix}.$$

The solution vector, (??), is

$$a = \begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \frac{1}{360} \begin{bmatrix} 1733 \\ 3387 \end{bmatrix}.$$

The residual error vector is

$$r = \frac{1}{360} \begin{bmatrix} -496 \\ 2207 \\ -1282 \\ -487 \\ -2284 \\ -121 \\ 2330 \\ 3485 \\ -3352 \end{bmatrix},$$

making the total error

$$r^*r = \frac{1\,139\,969}{3600}.$$

The uncertainties are then

$$\epsilon = \begin{bmatrix} \epsilon_0 \\ \epsilon_1 \end{bmatrix} = \left(360\sqrt{35}\right)^{-1} \begin{bmatrix} 108\,297\,055 \\ 3\,419\,907 \end{bmatrix}.$$

3.3.2 Computed Form

The results in the previous section are shown in integer form as debugging tool: readers can check answer to arbitrary precision. This section deals with formats appropriate for formal reporting. One way to quote numbers with uncertainties is using the \pm (plus – minus) notation:

$$\begin{aligned} a_0 &= 4.8 \pm 4.9 \quad \text{intercept} \quad ^\circ\text{C}, \\ a_1 &= 9.41 \pm 0.87 \quad \text{slope} \quad ^\circ\text{C} / \text{cm}. \end{aligned}$$

An alternative presentation uses parentheses:

$$\begin{aligned} a_0 &= 4.8(4.9) \quad \text{intercept} \quad ^\circ\text{C}, \\ a_1 &= 9.41(0.87) \quad \text{slope} \quad ^\circ\text{C} / \text{cm}. \end{aligned}$$

Quote the total error as $r^*r \approx 320$.

The uncertainty determines the number of significant digits. Common practice quotes the first two digits in the uncertainty (blue); the location of these two digits determines the number of significant digits in the solution parameter. The double precision computations are

$$\begin{aligned}\epsilon_0 &= 4.886206312183354 && \text{rounded to 4.9,} \\ a_0 &= 4.81388888888889 && \text{rounded to 4.8;} \\ \epsilon_1 &= 0.8633016476563611 && \text{rounded to 0.87,} \\ a_1 &= 9.408333333333333 && \text{rounded to 9.41.}\end{aligned}$$

At this point the model can be explored and evaluated. If the model is not acceptable, another trial function can be posed. Otherwise, the trial function becomes the solution function and is stated with error:

$$\begin{aligned}T(x) &= a_0 + a_1x, \\ \epsilon_T^2(x) &= \epsilon_0^2 + x^2\epsilon_1^2 + a_1^2\epsilon_x^2,\end{aligned}$$

which allows for interpolation and extrapolation. What happens when the solution is extrapolated to the heat baths? The expected answers are 0°C at 0 cm, and 100°C at 10 cm:

$$\begin{aligned}T(0) &= (4.8 \pm 4.9)^\circ\text{C}, \\ T(10) &= (99. \pm 10.)^\circ\text{C}.\end{aligned}$$

One final thought. The method of least squares minimizes the sums of the squares of the residual errors. And in linear regression, the sum of these residuals must be 0. That is,

$$\sum_{k=1}^m r_k = 0.$$

This can be a quick method for evaluating solutions and data sets. Given the data and the solution parameters a a quick Python or Mathematica script can compute and sum the residuals. If a data point is omitted, the sum will not be 0. If the parameters are misquoted, the sum will not be 0. If a data point is corrupted, the sum will not be 0. Or if the solutions are for another data set, the sum will not be 0. The 0 test is simple and powerful.

Thanks to integer arithmetic, it is easy to verify that the sum of the residuals in this problem is exactly 0.

3.4 Visualization

3.4.1 Seeing the Solution

Numbers tell a story and plots bring a story to life. Besides a powerful and immediate impact, the plots are often a defense against a wide host of problems.

The first plot is the solution curve plotted against the measurements as shown in figure 3.3. The residual errors are shown as red segments which are the actual components of the residual error vector in \mathbb{C}^9 . That is, the length of the red bars are given by the coordinates of r in the null space $\mathcal{N}(\mathbf{A}^*)$. But what is the basis for which these coordinates correspond? That will be addressed in §4.2.

Examination of this first plot is a first step, a crude tool which shows that the model was correctly applied. It reveals that there are no gross systematic errors; it says little about the quality of the fit. For that, more refinement is needed.

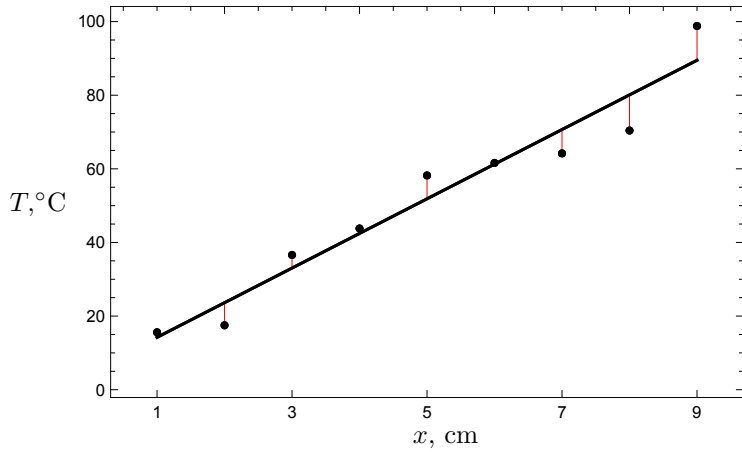


Figure 3.2. Solution plotted against data with residual errors shown in red.

Next, examine the residual errors as in figure 3.3. This is a more detailed examination of fit quality. The resolution increases by roughly a factor of 10 here. The dark line shows the mean of the residuals; exactly 0 in this case. The gray band represents the first standard deviation.

If data points need to be excluded, the residuals plot buttresses the case. The lone point that is five standard deviations away from the mean is quite obvious. Yes, it will be apparent on the solution plot, but that does not show rejection criteria in terms of the standard deviation.

Ideally, the residuals should be uncorrelated and scattered. In this case, there is a hint of correlation; a hint that the model may be extended. Or we could be a victim of low statistics – the behavior is random, but the sampling is too limited. To help think about the issue of correlation, just connect the dots as in figure 3.4. The result should be a jittery line with an oscillating appearance, not a trend line. As in this case.

The third and final diagnostic plot, figure 3.3, puts the solution on a map of the merit function. This addresses the question of whether the solution is a true minimum and applies whether the problem is linear or nonlinear. Complicated codes can produce answers which appear valid and this plot helps to filter the good from the bad.

The plot shows a level sets of the merit function, loci where $M(a) = \text{const.}$ The least squares solution is plotted in the middle. The three concentric circles

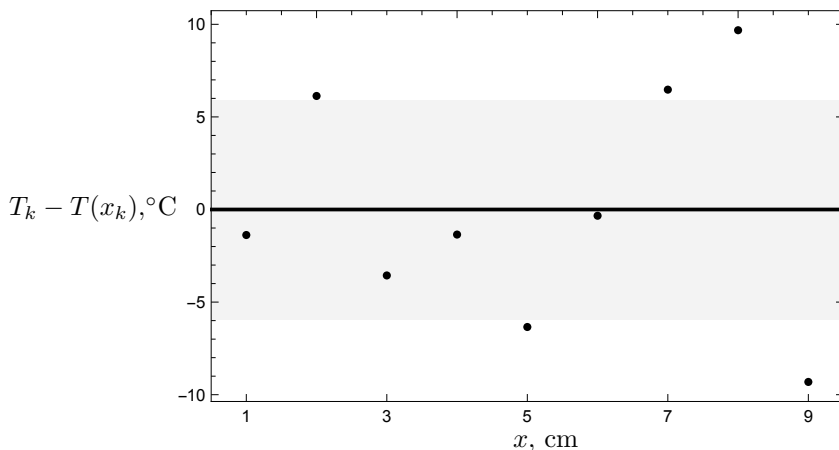


Figure 3.3. Scatter plot of residual errors showing mean (thick line) and one standard deviation (gray box).

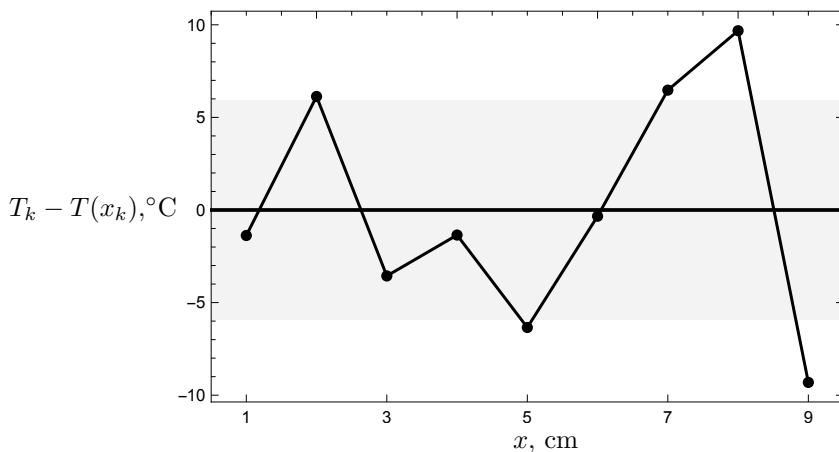


Figure 3.4. Scatter plot of residual errors with data points connected.

represent 1, 2, and 3 standard deviations in the fit parameters. A modified version of this plot is shown in figure 3.4.1.

The modified version explicitly labels the radii of the error ellipse, ϵ_0 , and ϵ_1 . The contour values are marked and create a feeling of a paraboloid with a minimum at $r^2 \approx 317$. The large arrowheads show the location of the solution precisely along the axes. The smaller escort arrowheads show the extent of the first standard deviation.

3.4.2 Seeing the Uncertainty

The interpretation of the variation of the parameters a is straightforward. A nice visualization is the “whisker plot” in figure 3.7. Given the mean a and the standard



Figure 3.5. Contour plot of merit function showing the solution (white cross) and three concentric error bands.

deviation ϵ , solutions can be sampled as a population. For the plot, 250 solutions a_μ were sampled and drawn against the data.

$$T_\mu(x) = a_{0_\mu} + a_{1_\mu}x, \quad \mu = 1:250.$$

Each solution represents a line. All of the lines provide a qualitative feel for the simultaneous variation of slope and intercept prescribed by the solution. Notice the white dots representing the ideal solution.

3.4.3 Digging Deeper

The whisker plot is a nice tool to quickly communicate the variation of the solution. This section describes the nuts and bolts of creating the representation. The primary tool is the assumed distribution of errors, here the normal distribution:

$$f(x; \mu, \sigma) = \frac{1}{\sigma\sqrt{2\pi}} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$$

The curve is normalized to unity, that is,

$$\int_{-\infty}^{\infty} f(x; \mu, \sigma) dx = 1.$$

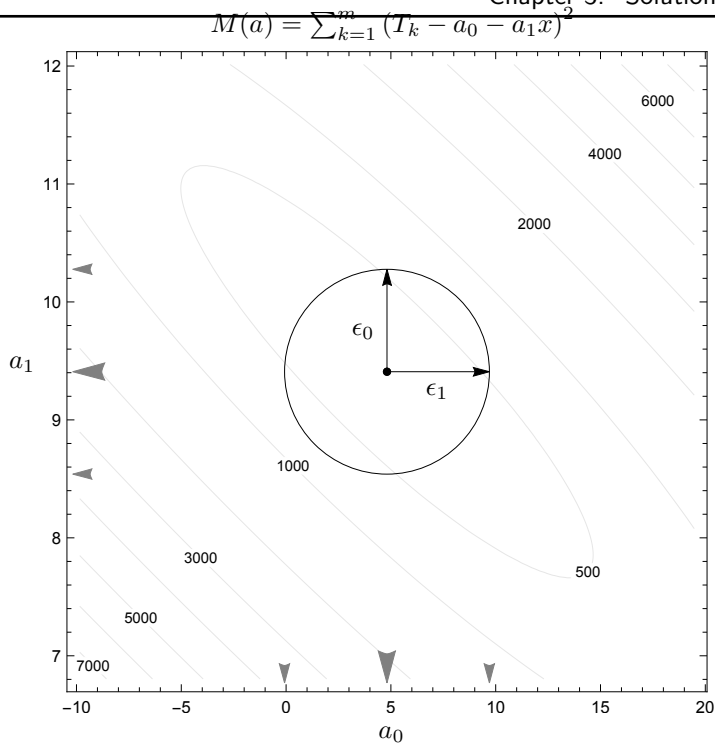


Figure 3.6. Another look at the merit function showing the primary error ellipse.

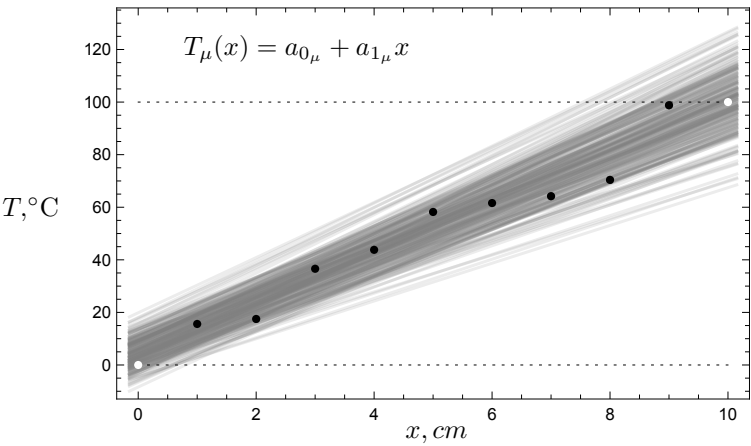
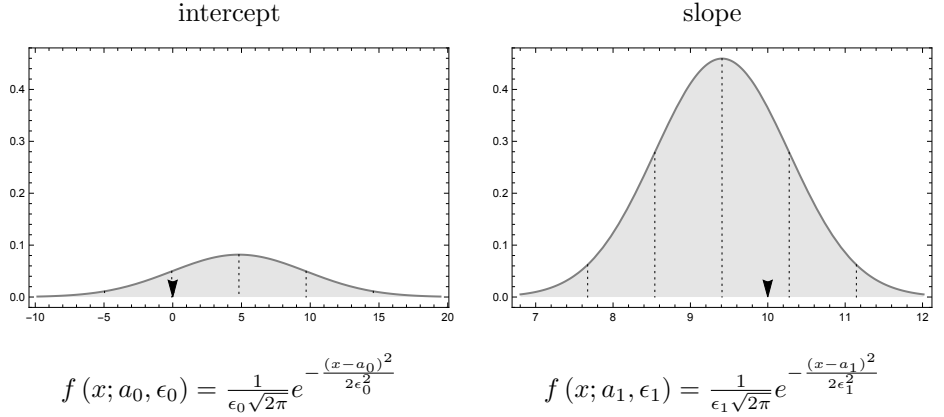


Figure 3.7. Whisker plot showing 250 randomly sampled solutions.

The mean value is μ , and the standard deviation is σ . For the data analysis, the mean represents the fit parameters a , and the standard deviation represents the error ϵ . Figure 3.4 shows these connections explicitly and provides a qualitative representation for the quality of the parameters.

The numbers on the vertical axis are devoid of meaning and are included only

Table 3.4. *The solution parameters expressed as normal distributions.*

to emphasize the point that the two distributions were plotted on a common scale. Knowing that, it is apparent that the slope measurement on the right is much more precise than the intercept measurement on the left.

Arrowheads locate the ideal solution and dotted lines demarcate the first three standard deviations. This provides a visual gauge of the accuracy and precision of the calculation.

Figure 3.8 shows the actual solutions used in the whisker plot as 250 points in the solution space $\mathcal{R}(\mathbf{A}^*)$. The symbol “+” marks the ideal solution.

A crude, but useful tool, is the counting of the number of points within the error bands which allows for a direct comparison with continuum theory. For example, the innermost circle should contain, in the continuum limit, 68.27% of the data points. The second band should envelope 95.45% of the points, and so on. Table 3.5 displays the actual census and compares to the continuum results.

The first column identifies the annular region in which the points were counted, and the census for that region. For example, the fourth annulus contains just 6 points. The area column displays the relative area of the annular regions computed using the area formula

$$\text{area}_k = \pi \left(k^2 - (k-1)^2 \right), \quad k = 1:5.$$

The relative area removes the constant π ; this does not affect the final answer and it simplifies the cross check. The count density divides the counts in a zone by the relative area of that zone, e.g. $115/3 = 38.3$. The normalized density is the ratio of the count density to the sum of the count densities, e.g. $38.3/135.7 = 28.24$. The cumulative density accumulates the normalized densities. These normalized

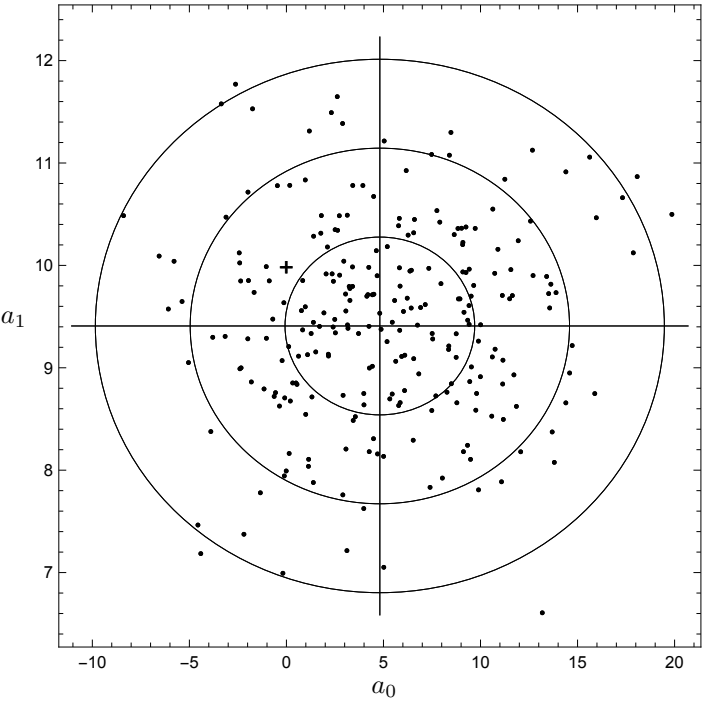


Figure 3.8. Scatter plot showing sampling of 250 solution points in $\mathcal{R}(\mathbf{A}^*)$.

Table 3.5. Comparing samples to ideal normal distribution.

annulus	count	rel. area	count density	normalized density	cumulative density	continuum limit
1	88	1	88.0	64.83%	64.83%	68.27%
2	115	3	38.3	28.24%	93.07%	95.45%
3	41	5	8.2	6.41%	99.16%	99.73%
4	6	5	1.2	0.88%	100.0%	99.99%
250			135.7	100.00%		

densities approximate the final column, the continuum limit given by

$$\int_{-k\sigma}^{k\sigma} f(x; \mu = 1, \sigma = 1) dx.$$

Chapter 4

Solution via SVD

The singular value decomposition is an x -ray revealing the structure of the fundamental spaces.

4.1 Computing the SVD

Solution steps

1. Compute $\lambda(\mathbf{A}^* \mathbf{A})$.
2. Educated guess at domain matrix \mathbf{V} .
3. Compute codomain matrix \mathbf{U} .

The least squares problem delivers a singular value decomposition (SVD) without the muss and fuss of solving an eigensystem. The SVD is given by the matrix product

$$\mathbf{A} = \mathbf{U} \mathbf{\Sigma} \mathbf{V}^*.$$

For the full column rank problem we have we can expand in the following block decomposition

$$\mathbf{A} = \left[\begin{array}{c|c} \mathbf{U}_{\mathcal{R}} & \mathbf{U}_{\mathcal{N}} \end{array} \right] \left[\begin{array}{c} \mathbf{S} \\ \mathbf{0} \end{array} \right] \left[\begin{array}{c} \mathbf{V}_{\mathcal{R}}^* \end{array} \right]$$

The \mathbf{S} matrix contains the singular values

$$\mathbf{S} = \begin{bmatrix} \sigma_1 & 0 \\ 0 & \sigma_2 \end{bmatrix}.$$

The column vectors of the matrix $\mathbf{V}_{\mathcal{R}}$ represent an orthonormal basis for the row space (domain). The column vectors of the matrix $\mathbf{U}_{\mathcal{R}}$ represent two of the nine vectors in an orthonormal basis for the column space (codomain).

4.1.1 Singular values

The singular value spectrum of the matrix \mathbf{A} is the square root of the (non-zero) eigenvalues of the product matrix $\mathbf{A}^* \mathbf{A}$

$$\sigma(\mathbf{A}) = \sqrt{\lambda(\mathbf{A}^* \mathbf{A})}.$$

The eigenvalues of the product matrix are the roots of the characteristic polynomial $p(\lambda)$ for said matrix.

$$p(\lambda) = \lambda^2 - \lambda \operatorname{tr}(\mathbf{A}^* \mathbf{A}) + \det(\mathbf{A}^* \mathbf{A})$$

The determinant appeared in equations (9.3) and (5.2); the trace is $\operatorname{tr}(\mathbf{A}^* \mathbf{A}) = \mathbf{1}^T \mathbf{1} + x^T x$. The singular values are then

$$\sigma = \sqrt{\frac{1}{2} \left(\mathbf{1}^T \mathbf{1} + x^T x \pm \sqrt{4(\mathbf{1}^T x)^2 - (\mathbf{1}^T \mathbf{1} - x^T x)^2} \right)}.$$

(The astute reader will notice that the discriminant does not seem to have the familiar form of $b^2 - 4ac$. The earnest reader will discover why this is so.) The singular value spectrum for these data is

$$\sigma = \sqrt{3 \left(49 \pm \sqrt{2341} \right)} \approx (17.0924, 1.35954).$$

We now have the sigma matrix and the matrix of singular values \mathbf{S} :

$$\Sigma = \begin{bmatrix} \mathbf{S} \\ \mathbf{0} \end{bmatrix} = \begin{bmatrix} \sigma_1 & 0 \\ 0 & \sigma_2 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \end{bmatrix} \approx \begin{bmatrix} 17.0924 & 0 \\ 0 & 1.35954 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \end{bmatrix}.$$

4.1.2 Domain matrix

We can skip the eigenvector problem. To find the domain matrix we exploit the singular value decomposition of the product matrix

$$\Sigma^T \Sigma = \begin{bmatrix} \mathbf{S} & \mathbf{0} \end{bmatrix} \begin{bmatrix} \mathbf{S} \\ \mathbf{0} \end{bmatrix} = \mathbf{S}^2$$

$$\mathbf{V}_{\mathcal{R}} \mathbf{S}^2 \mathbf{V}_{\mathcal{R}}^* = \mathbf{A}^* \mathbf{A}. \quad (4.1)$$

By the singular value theorem the matrix is unitary and will be a rotation matrix, a reflection matrix or a convolution. We begin by trying a rotation matrix

$$\mathbf{V}_{\mathcal{R}} = \begin{bmatrix} \cos \theta & -\sin \theta \\ \sin \theta & \cos \theta \end{bmatrix} \quad (4.2)$$

which is colored blue because the column vectors belong to $\mathcal{R}(\mathbf{A}^*)$. The objective is to find the angle θ . The immediate result of equations (??), (4.1), and (4.2) is

$$\begin{bmatrix} \cos \theta & -\sin \theta \\ \sin \theta & \cos \theta \end{bmatrix} \mathbf{S}^2 \begin{bmatrix} \cos \theta & -\sin \theta \\ \sin \theta & \cos \theta \end{bmatrix}^T = \mathbf{A}^* \mathbf{A},$$

$$\begin{bmatrix} \sigma_1^2 \cos^2 \theta + \sigma_2^2 \sin^2 \theta & (\sigma_1^2 - \sigma_2^2) \cos \theta \sin \theta \\ (\sigma_1^2 - \sigma_2^2) \cos \theta \sin \theta & \sigma_2^2 \cos^2 \theta + \sigma_1^2 \sin^2 \theta \end{bmatrix} = \begin{bmatrix} \mathbf{1}^T \mathbf{1} & \mathbf{1}^T x \\ x^T \mathbf{1} & x^T x \end{bmatrix}.$$

which presents multiple solution paths for the angle θ . For example

$$\cos \theta = \sqrt{\frac{\sigma_2^2 - \mathbf{1}^T \mathbf{1}}{\sigma_2^2 - \sigma_1^2}} = \sqrt{\frac{x^T x - \sigma_1^2}{\sigma_2^2 - \sigma_1^2}}. \quad (4.3)$$

This implies

$$\sin \theta =$$

The domain matrix is now

$$\mathbf{V} = \begin{bmatrix} \mathbf{V}_{\mathcal{R}} \end{bmatrix} = (\sigma_2^2 - \sigma_1^2)^{-\frac{1}{2}} \begin{bmatrix} \sqrt{\sigma_2^2 - \mathbf{1}^T \mathbf{1}} & -\sqrt{\mathbf{1}^T \mathbf{1} - \sigma_1^2} \\ \sqrt{\mathbf{1}^T \mathbf{1} - \sigma_1^2} & \sqrt{\sigma_2^2 - \mathbf{1}^T \mathbf{1}} \end{bmatrix}.$$

Using the data set at hand

$$\begin{bmatrix} \mathbf{V}_{\mathcal{R}} \end{bmatrix} = \begin{bmatrix} \sqrt{\frac{1}{2} - \frac{23}{\sqrt{2341}}} & -\sqrt{\frac{1}{2} + \frac{23}{\sqrt{2341}}} \\ \sqrt{\frac{1}{2} + \frac{23}{\sqrt{2341}}} & \sqrt{\frac{1}{2} - \frac{23}{\sqrt{2341}}} \end{bmatrix} \approx \begin{bmatrix} 0.156956 & -0.987606 \\ 0.987606 & 0.156956 \end{bmatrix}$$

4.1.3 Codomain matrix

The final component is of course the codomain matrix. Knowing the decomposition for the adjoint matrix \mathbf{A}^* and that the linear system is overdetermined we can write

$$\mathbf{U}_{\mathcal{R}}^* = \mathbf{S}^{-1} \mathbf{V}_{\mathcal{R}}^* \mathbf{A}^*.$$

The k th column vector of this matrix has the compact form

$$[\mathbf{U}_{\mathcal{R}}^*]_k = (\sigma_1^2 - \sigma_2^2)^{-\frac{1}{2}} \begin{bmatrix} \sigma_1^{-2} \left(\sqrt{\sigma_2^2 - \mathbf{1}^T \mathbf{1}} - x \sqrt{\mathbf{1}^T \mathbf{1} - \sigma_1^2} \right) \\ \sigma_2^{-2} \left(\sqrt{\sigma_2^2 - \mathbf{1}^T \mathbf{1}} + x \sqrt{\mathbf{1}^T \mathbf{1} - \sigma_1^2} \right) \end{bmatrix}.$$

Using the following shorthand,

$$f(x, y) = \sqrt{x + y/\sqrt{2341}},$$

the range space component of the codomain matrix can be written as

$$\mathbf{U}_{\mathcal{R}} = \left(6\sqrt{10}\right)^{-1} \begin{bmatrix} f(68, -3212) & -f(68, 3212) \\ f(47, -2003) & -f(47, 2003) \\ f(32, -968) & -f(32, 968) \\ f(23, -107) & -f(23, 107) \\ f(50, 580) & -f(50, -580) \\ f(23, 1093) & -f(23, -1093) \\ f(32, 1432) & f(32, -1432) \\ f(47, 1597) & f(47, -1597) \\ f(68, 1588) & f(68, -1588) \end{bmatrix}$$

If one wishes to complete the codomain matrix, use the Gram-Schmidt orthonormalization process on the matrix

$$\mathbf{U} = \begin{bmatrix} [\mathbf{U}_{\mathcal{R}}]_1 & [\mathbf{U}_{\mathcal{R}}]_2 & \mathbf{e}_{1,9} & \mathbf{e}_{2,9} & \mathbf{e}_{3,9} & \mathbf{e}_{4,9} & \mathbf{e}_{5,9} & \mathbf{e}_{6,9} & \mathbf{e}_{7,9} \end{bmatrix}$$

starting with column three.

Completing the codomain matrix with an orthonormal span of the null space is optional and can be done by feeding the range space components and a complementary set of unit vectors into a Gram-Schmidt algorithm.

4.1.4 Error terms

$$\epsilon_k^2 = \frac{(\mathbf{A}\alpha - T)^* (\mathbf{A}\alpha - T)}{m - n} \left[(\mathbf{A}^* \mathbf{A})^{-1} \right]_{kk}$$

The error terms can be computed after this step. Given that the product matrix decomposition in (4.1) is a singular value decomposition we can trivially write the inverse matrix as

$$(\mathbf{A}^* \mathbf{A})^{-1} = \mathbf{V}_{\mathcal{R}} \mathbf{S}^{-2} \mathbf{V}_{\mathcal{R}}^* = \frac{1}{180} \begin{bmatrix} 95 & -15 \\ -15 & 3 \end{bmatrix}$$

The error terms in (??) become

$$\begin{aligned} \epsilon &= \sqrt{\frac{r^* r}{\mathbf{1}^T \mathbf{1} - n}} \sqrt{\frac{1}{\sigma_1 \sigma_2}} \sqrt{\left[\begin{array}{c} \frac{\cos^2 \theta}{\sigma_2^2} + \frac{\sin^2 \theta}{\sigma_1^2} \\ \frac{\cos^2 \theta}{\sigma_1^2} + \frac{\sin^2 \theta}{\sigma_2^2} \end{array} \right]}, \\ &= \sqrt{\frac{r^* r}{m - n}} \sqrt{\left[\begin{array}{c} \sigma_1^2 - \sqrt{(\sigma_2^2 - \mathbf{1}^T \mathbf{1})(\sigma_2^2 - \sigma_1^2)} \\ \sigma_2^2 + \sqrt{(\sigma_2^2 - \mathbf{1}^T \mathbf{1})(\sigma_2^2 - \sigma_1^2)} \end{array} \right]}. \end{aligned}$$

To close this section, look at a numeric representation of the \mathbf{U} matrix:

$$\mathbf{U} = \begin{bmatrix} 0.0670 & -0.611 & 0.656 & -0.0132 & -0.0780 & -0.137 & -0.193 & -0.240 & -0.270 \\ 0.125 & -0.496 & -0.749 & -0.0927 & -0.113 & -0.147 & -0.184 & -0.215 & -0.233 \\ 0.183 & -0.380 & 0 & 0 & 0 & 0 & 0 & 0 & 0.907 \\ 0.240 & -0.265 & 0 & 0 & 0 & 0 & 0 & 0.920 & -0.159 \\ 0.298 & -0.149 & 0 & 0 & 0 & 0 & 0.924 & -0.142 & -0.123 \\ 0.356 & -0.0337 & 0 & 0 & 0 & 0.910 & -0.150 & -0.117 & -0.0858 \\ 0.414 & 0.0817 & 0 & 0 & 0.867 & -0.198 & -0.141 & -0.0930 & -0.0490 \\ 0.471 & 0.197 & 0 & 0.755 & -0.321 & -0.209 & -0.132 & -0.0685 & -0.0123 \\ 0.529 & 0.313 & 0.0937 & -0.649 & -0.356 & -0.219 & -0.124 & -0.0441 & 0.0245 \end{bmatrix}$$

4.2 Visualization

With the singular value decomposition in hand, the domain space plots become more concrete and we do so below beginning in figure (4.2). The black vector represents the measurements

$$T = \mathbf{A}a - \mathbf{R}$$

$$y \in \mathbb{C}^9, \quad \mathbf{A}a \in \mathcal{R}(\mathbf{A}) \subseteq \mathbb{C}^2, \quad r \in \mathcal{N}(\mathbf{A}^*) \subseteq \mathbb{C}^7$$

$$T = \frac{1}{10} \begin{bmatrix} 156 \\ 175 \\ 366 \\ 438 \\ 582 \\ 616 \\ 642 \\ 704 \\ 988 \end{bmatrix}$$

The closest point to the data vector in the range $\mathcal{R}(\mathbf{A})$ is

$$\mathbf{A}a = \frac{1}{360} \begin{bmatrix} 5120 \\ 8507 \\ 11894 \\ 15281 \\ 18668 \\ 22055 \\ 25442 \\ 28829 \\ 32216 \end{bmatrix} = \alpha_1 [\mathbf{U}_{\mathcal{R}}]_1 + \alpha_2 [\mathbf{U}_{\mathcal{R}}]_2 \in \mathcal{R}(\mathbf{A})$$

where the coordinates are

$$\begin{aligned} \begin{bmatrix} \alpha_1 \\ \alpha_2 \end{bmatrix} &= \left(30 \left(\sqrt{2341} - 31 \right) \sqrt{4682} + 58\sqrt{2341} \right)^{-1} \times \\ &\quad \begin{bmatrix} 4\,104\,889 + 75\,341\sqrt{2341} \\ 3\sqrt{15} \left(753\,593 - 15\,933\sqrt{2341} \right) \end{bmatrix} \\ &\approx \begin{bmatrix} 171.733 \\ -4.45594 \end{bmatrix} \\ T_{\mathcal{R}} &\approx 171.733u_1 - 4.45594u_2 \end{aligned} \quad (4.4)$$

Another way to think of the geometry is to pose the merit function in terms of the

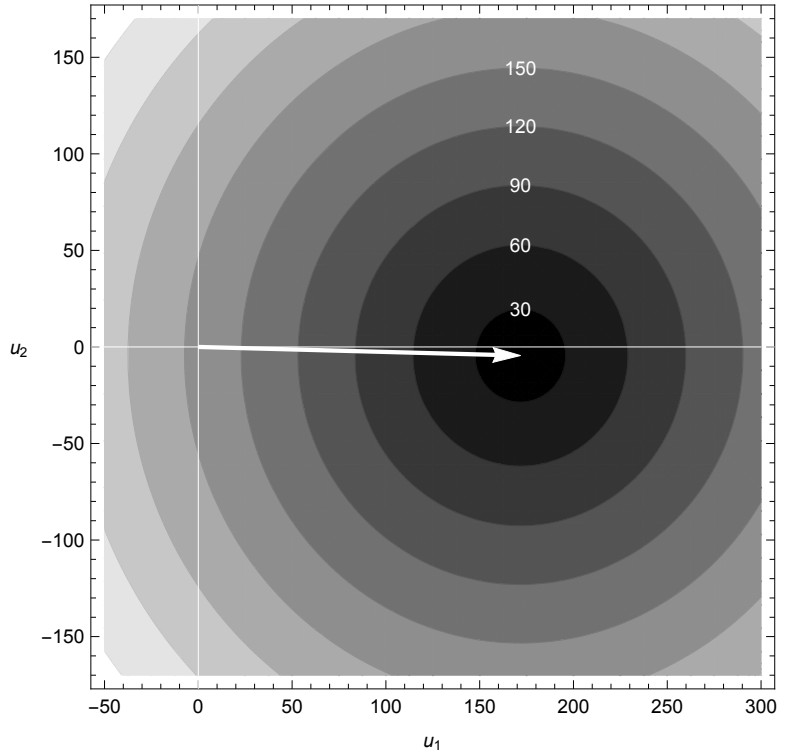


Figure 4.1. The solution vector in (4.4) (white arrow) is the mixture of u_1 and u_2 which eliminates the error.

range space vectors

$$M(\alpha) = \|T - \alpha_1 u_1 - \alpha_2 u_2\|_2^2.$$

The minimizer is given by (4.2) and displayed in figure 4.1.

$$M(171.733, -4.45594) \approx 17.7949.$$

The residual error vector lies entirely in the null space $\mathcal{N}(\mathbf{A}^*)$

$$r = -\frac{1}{360} \begin{bmatrix} 496 \\ -2207 \\ 1282 \\ 487 \\ 2284 \\ 121 \\ -2330 \\ -3485 \\ 3352 \end{bmatrix}$$

$$= \alpha_3[\mathbf{U}_{\mathcal{N}}]_1 + \alpha_4[\mathbf{U}_{\mathcal{N}}]_2 + \alpha_5[\mathbf{U}_{\mathcal{N}}]_3 + \alpha_6[\mathbf{U}_{\mathcal{N}}]_4 + \alpha_7[\mathbf{U}_{\mathcal{N}}]_5 + \alpha_8[\mathbf{U}_{\mathcal{N}}]_6 + \alpha_9[\mathbf{U}_{\mathcal{N}}]_7$$

$$\in \mathcal{N}(\mathbf{A}^*)$$

The coordinates are now

$$\begin{bmatrix} \alpha_3 \\ \alpha_4 \\ \alpha_5 \\ \alpha_6 \\ \alpha_7 \\ \alpha_8 \\ \alpha_9 \end{bmatrix} = \left(60\sqrt{24\,747\,709}\right)^{-1} \begin{bmatrix} 680\sqrt{7\,815\,066} \\ -1933\sqrt{3\,907\,533} \\ -3621\sqrt{186\,073} \\ 6679\sqrt{10\,434} \\ 13\,406\sqrt{29\,526} \\ 2196\sqrt{85\,386} \\ 641\sqrt{3\,344\,285} \end{bmatrix}$$

$$T = \mathbf{U}\alpha = \mathbf{A}a - r$$

The decomposition shown in figure 2.1.

$$\frac{1}{360} \begin{bmatrix} 5616 \\ 6300 \\ 13\,176 \\ 15\,768 \\ 20\,952 \\ 22\,176 \\ 23\,112 \\ 25\,344 \\ 35\,568 \end{bmatrix} = \frac{1}{360} \begin{bmatrix} 5120 \\ 8507 \\ 11\,894 \\ 15\,281 \\ 18\,668 \\ 22\,055 \\ 25\,442 \\ 28\,829 \\ 32\,216 \end{bmatrix} + \frac{1}{360} \begin{bmatrix} 496 \\ -2207 \\ 1282 \\ 487 \\ 2284 \\ 121 \\ -2330 \\ -3485 \\ 3352 \end{bmatrix}$$

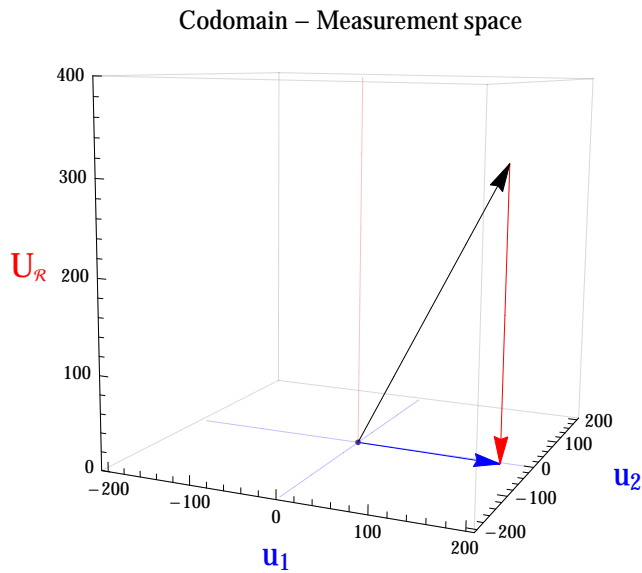


Figure 4.2. Measurement space $\mathcal{R}(\mathbf{A})$ for the Bevington example showing resolution of the data vector (black) into range (blue) and null space (red) components.

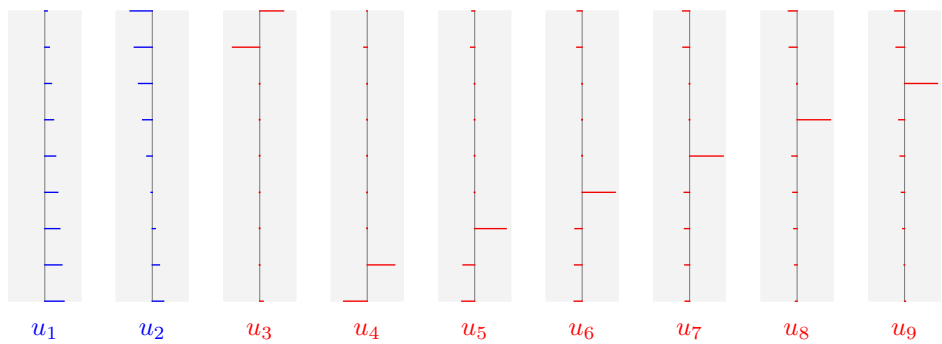


Table 4.1. The column vectors of \mathbf{U} . The gray box represents the maximum length an element may have: $[-1, 1]$.

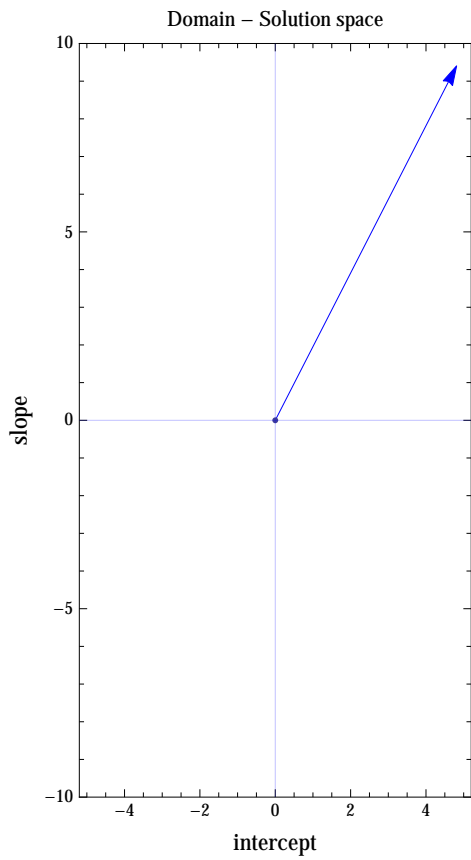


Figure 4.3. *Minimization occurs in the codomain.*

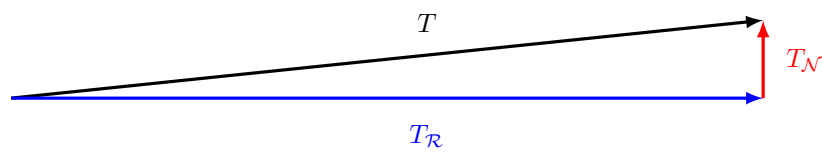


Figure 4.4. *Data vector resolved into range and null space components.*



Figure 4.5. *Decomposing $\|r = T_N\|_2^2$ into residual error terms r_k^2 of table 4.3.*

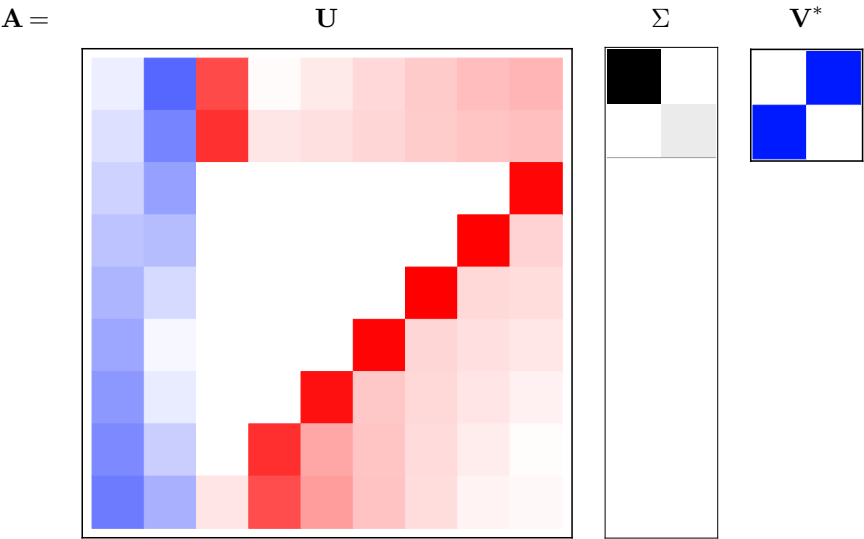


Table 4.2. Singular value decomposition for the system matrix A in (5.1) showing range and null space components.

Table 4.3. A summary of the residual errors and their contributions to $\|r\|_2$.

k	r_k	r_k^2
1	-1.4	1.9
2	6.1	37.6
3	-3.6	12.7
4	-1.4	1.8
5	-6.3	40.3
6	-0.3	0.1
7	6.5	41.9
8	9.7	93.7
9	-9.3	86.7

Chapter 5

Solution Via Other Methods

For comparison, other solution methods are applied to the Bevington data. These methods will reinforce the geometry of least squares.

5.1 Normal Equations from Vectors

In §3.2.2, the normal equations appeared from the linear system resulting from the two partial differential equations in (3.3). Another approach starts with the linear equation $\mathbf{A}x = b$ and uses column vectors. The data is posed in terms of the column vectors in table 3.2. These, plus a constant vector, are the elements of composition:

$$\mathbf{1} = \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix}, \quad x = \begin{bmatrix} 1 \\ 2 \\ 3 \\ 4 \\ 5 \\ 6 \\ 7 \\ 8 \\ 9 \end{bmatrix}, \quad T = \frac{1}{10} \begin{bmatrix} 156 \\ 175 \\ 366 \\ 438 \\ 582 \\ 616 \\ 642 \\ 704 \\ 988 \end{bmatrix}.$$

There are two column vectors in the system matrix \mathbf{A} :

$$\mathbf{A} = \left[\mathbf{1} \mid x \right]$$

In the service of a crisp mental image, the linear system is written explicitly:

$$\begin{array}{ccc} \mathbf{A} & a & = T \\ \left[\begin{array}{cc} 1 & 1 \\ 1 & 2 \\ 1 & 3 \\ 1 & 4 \\ 1 & 5 \\ 1 & 6 \\ 1 & 7 \\ 1 & 8 \\ 1 & 9 \end{array} \right] & \left[\begin{array}{c} a_0 \\ a_1 \end{array} \right] = \frac{1}{10} & \left[\begin{array}{c} 156 \\ 175 \\ 366 \\ 438 \\ 582 \\ 616 \\ 642 \\ 704 \\ 988 \end{array} \right]. \end{array} \quad (5.1)$$

5.1.1 Composing the normal equations

The system encodes nine linear problems:

$$\begin{aligned} a_0 + a_1 x_1 &= T_1, \\ &\vdots \\ a_0 + a_1 x_9 &= T_9. \end{aligned}$$

But there are no parameters a_0 and a_1 which solve all nine equations. More formally, the data vector T is not in the column space of \mathbf{A} . There are no parameters a_0 and a_1 such that

$$a_0 \mathbf{1} + a_1 x = T.$$

One strategy is to compose a solution does have a solution

$$\mathbf{A}^* \mathbf{A} a = \mathbf{A}^* T.$$

Certainly the vector $\mathbf{A}^* T$ is in the column space of \mathbf{A}^* – the coordinates are T ! The solution is

$$\left[\begin{array}{c} a_0 \\ a_1 \end{array} \right] = (\mathbf{A}^* \mathbf{A})^{-1} \mathbf{A}^* T.$$

For the Bevington data set the dot products are,

$$\begin{aligned} \mathbf{1}^T \mathbf{1} &= m = 9, \\ \mathbf{1}^T x &= x^T \mathbf{1} = 45, \\ x^T x &= 285, \\ \mathbf{1}^T T &= \frac{4667}{10}, \\ x^T T &= 2898. \end{aligned}$$

Other intermediate products include the product matrix,

$$\mathbf{A}^* \mathbf{A} = \begin{bmatrix} \mathbf{1}^T \mathbf{1} & \mathbf{1}^T x \\ x^T \mathbf{1} & x^T x \end{bmatrix} = \begin{bmatrix} 9 & 45 \\ 45 & 285 \end{bmatrix},$$

its determinant,

$$\Delta = (\mathbf{1}^T \mathbf{1}) (x^T x) - (\mathbf{1}^T x)^2, \quad (5.2)$$

inverse,

$$(\mathbf{A}^* \mathbf{A})^{-1} = \Delta^{-1} \begin{bmatrix} x^T x & -\mathbf{1}^T \mathbf{1} \\ -\mathbf{1}^T \mathbf{1} & \mathbf{1}^T x \end{bmatrix}$$

and the vector,

$$\mathbf{A}^* T = \begin{bmatrix} \mathbf{1}^T T \\ x^T T \end{bmatrix} = \frac{1}{10} \begin{bmatrix} 4667 \\ 28980 \end{bmatrix}.$$

The solution is again

$$a = \Delta^{-1} \begin{bmatrix} x^T x & -\mathbf{1}^T \mathbf{1} \\ -\mathbf{1}^T \mathbf{1} & \mathbf{1}^T x \end{bmatrix} \begin{bmatrix} \mathbf{1}^T T \\ x^T T \end{bmatrix}$$

$$\begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \Delta^{-1} \begin{bmatrix} (x^T x) (\mathbf{1}^T T) - (\mathbf{1}^T x) (x^T T) \\ (\mathbf{1}^T \mathbf{1}) (x^T T) - (\mathbf{1}^T x) (\mathbf{1}^T T) \end{bmatrix} \quad (5.3)$$

leading to the same solution presented in §3.3.1.

5.2 QR Decomposition

Resolve the range space $\mathcal{R}(\mathbf{A})$ into an orthonormal basis.

$$\mathbf{A} = \mathbf{Q} \mathbf{R}$$

$$a = \mathbf{R}^{-1} \mathbf{Q}^* T$$

5.2.1 Computing the QR Decomposition

$$\mathbf{R} \in \mathbb{R}^{\rho \times \rho}$$

$$\mathbf{R} = \begin{bmatrix} r_{11} & r_{12} \\ 0 & r_{22} \end{bmatrix} = \begin{bmatrix} \nu_1 & q_1^2 a_2 \\ 0 & \nu_2 \end{bmatrix}$$

The norm of the first column vector:

$$r_{11} = \|a_1\|_2 = 3 \quad (5.4)$$

Normalized column vector

$$q_1 = \frac{a_1}{r_{11}} = \frac{1}{3} \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix} \tag{5.5}$$

Next scale factor in row 1:

$$r_{12} = q_1^* a_2 = 15 \tag{5.6}$$

$$q_2 = a_2 - r_{12} q_1 \tag{5.7}$$

$$\mathbf{Q} = \begin{bmatrix} \frac{1}{3} \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix} \quad \frac{1}{2\sqrt{15}} \begin{bmatrix} -4 \\ -3 \\ -2 \\ -1 \\ 0 \\ 1 \\ 2 \\ 3 \\ 4 \end{bmatrix} \end{bmatrix}$$

$$\begin{matrix} \mathbf{A} & = & \mathbf{Q} & \mathbf{R}, \\ \begin{bmatrix} 1 & 1 \\ 1 & 2 \\ 1 & 3 \\ 1 & 4 \\ 1 & 5 \\ 1 & 6 \\ 1 & 7 \\ 1 & 8 \\ 1 & 9 \end{bmatrix} & = & \begin{bmatrix} \frac{1}{3} \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix} \quad \frac{1}{2\sqrt{15}} \begin{bmatrix} -4 \\ -3 \\ -2 \\ -1 \\ 0 \\ 1 \\ 2 \\ 3 \\ 4 \end{bmatrix} \end{bmatrix} & \begin{bmatrix} 3 & 15 \\ 0 & 2\sqrt{15} \end{bmatrix} \end{matrix}$$

Chapter 6

Finer Points

A deeper look at the example.

6.1 Invariances

Invariance is a touchstone which allows for elegant discrimination of theoretical and numerical analysis. The critical insight is this: complicated problems have complicated solutions, but they retain simple invariances. Simple properties afford simple tools for checking complicated problems.

6.1.1 Translation Invariance

Intuitively, a translation of the coordinate system will not change the temperature measurements, which means the gradient of the temperature is invariant under translation of the coordinate system. For the example in figure 3.1, the temperature gradient was measured over the length of a 10 cm bar. For convenience, the end of the bar in contact with the ice bath was taken as the origin. But just as readily, that point could have been taken as $x = 5$ cm; the end of the bar touching the boiling water bath is then 15 cm. The physics is unchanged. For an arbitrary shift of x_* the gradient is invariant:

$$\nabla T = \frac{T_1 - T_0}{(x_1 + x_*) - (x_0 + x_*)} = \frac{T_1 - T_0}{x_1 - x_0}$$

How should the solution change? We insist the slope (gradient) remain unchanged. The origin shifts from $x_0 = 0$ to $x_0 = -x_*$, the intercept shifts according to

$$y(-x_0) = a_0 - a_1 x,$$

that is, $a_0 \rightarrow a_0 - a_1 x_*$.

Translation

Formally evaluate how the least squares solution in (??) transforms under the translation

$$x_k \rightarrow x_k + x_*$$

The position vector x can be written as the sum of the original vector plus the translation:

$$\begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_m \end{bmatrix} \rightarrow \begin{bmatrix} x_1 + x_* \\ x_2 + x_* \\ \vdots \\ x_m + x_* \end{bmatrix} = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_m \end{bmatrix} + \begin{bmatrix} x_* \\ x_* \\ \vdots \\ x_* \end{bmatrix}.$$

In vector notation this is

$$x \rightarrow x + \mathbf{1}x_*. \quad (6.1)$$

Demonstrations

First compute the transformation rules on the inner products using (6.1):

$$\begin{aligned} \mathbf{1}^T \mathbf{1} &\rightarrow \mathbf{1}^T \mathbf{1}, \\ \mathbf{1}^T x &\rightarrow \mathbf{1}^T x + x_* (\mathbf{1}^T \mathbf{1}), \\ x^T x &\rightarrow x^T x + 2x_* (\mathbf{1}^T x) + x_*^2 (\mathbf{1}^T \mathbf{1}), \\ x^T T &\rightarrow x^T T + x_* (\mathbf{1}^T T). \end{aligned} \quad (6.2)$$

Computation of the transformation of the intercept a_0 and slope a_1 requires computation of the transformation of the determinant in (??): (5.2):

$$\begin{aligned} \Delta &= (\mathbf{1}^T \mathbf{1}) (x^T x) - (\mathbf{1}^T x)^2, \\ &\rightarrow (\mathbf{1}^T \mathbf{1}) (x^T x + 2x_* (\mathbf{1}^T x) + x_*^2 (\mathbf{1}^T \mathbf{1})) - (\mathbf{1}^T x + x_* (\mathbf{1}^T \mathbf{1}))^2, \\ &= \Delta + (\mathbf{1}^T \mathbf{1}) (2x_* (\mathbf{1}^T x) + x_*^2 (\mathbf{1}^T \mathbf{1})) - (2x_* (\mathbf{1}^T \mathbf{1}) (\mathbf{1}^T x)) - (x_* (\mathbf{1}^T \mathbf{1}))^2, \\ &= \Delta. \end{aligned}$$

Thus the determinant is invariant under translation. Using the substitutions in (6.2) one can show (exercise ???) the solution parameters transform as expected:

$$\begin{aligned} a_0 &\rightarrow a_0 - a_1 x_*, \\ a_1 &\rightarrow a_1. \end{aligned}$$

Homework: C to F, affine transform $T \rightarrow \alpha T + \beta$

Computations

The slope is invariant, the intercept:

$$\begin{aligned} a_1 &\rightarrow a_1, \\ a_0 &\rightarrow a_0 - x_* a_1. \end{aligned}$$

Predictions for $x_* = 1$.

$$\begin{aligned} a_0 &\rightarrow -\frac{827}{180}, \\ a_1 &\rightarrow \frac{1129}{120}. \end{aligned} \tag{6.3}$$

$$\begin{aligned} \begin{bmatrix} a_0 \\ a_1 \end{bmatrix} &= \Delta^{-1} \begin{bmatrix} \sum x_k^2 & -\sum x_k \\ -\sum x_k & \sum 1 \end{bmatrix} \begin{bmatrix} \sum T_k \\ \sum T_k x_k \end{bmatrix}, \\ &= 540^{-1} \begin{bmatrix} 384 & -54 \\ -54 & 9 \end{bmatrix} \frac{1}{10} \begin{bmatrix} 4667 \\ 33647 \end{bmatrix}, \\ &= \frac{1}{360} \begin{bmatrix} 1733 \\ 3387 \end{bmatrix}. \end{aligned}$$

In agreement with (6.3).

Visuals

Compare to figure 3.2

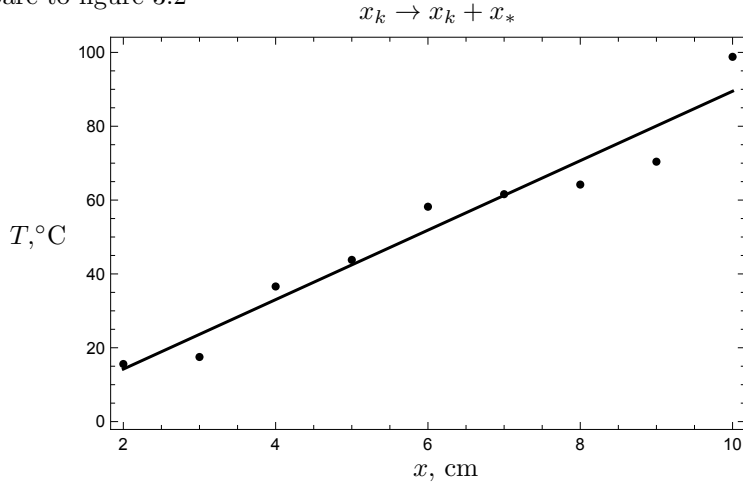


Figure 6.1. *Solution after translation along x axis: the slope is invariant.*

Compare to 3.3

Compare to figure 3.4.1

6.1.2 Reflection Invariance

A demonstration of the reflection method of §2.5.

$$\sum_{k=1}^m r^2 = \sum_{k=1}^m (-r)^2$$

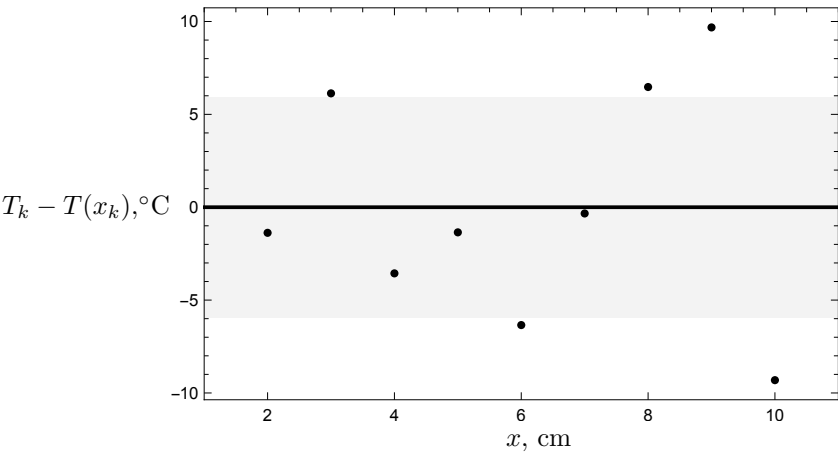


Figure 6.2. Scatter plot of residual errors after translation along x axis: the residuals are invariant.

table 13.6

6.2 Fitting To Higher Orders

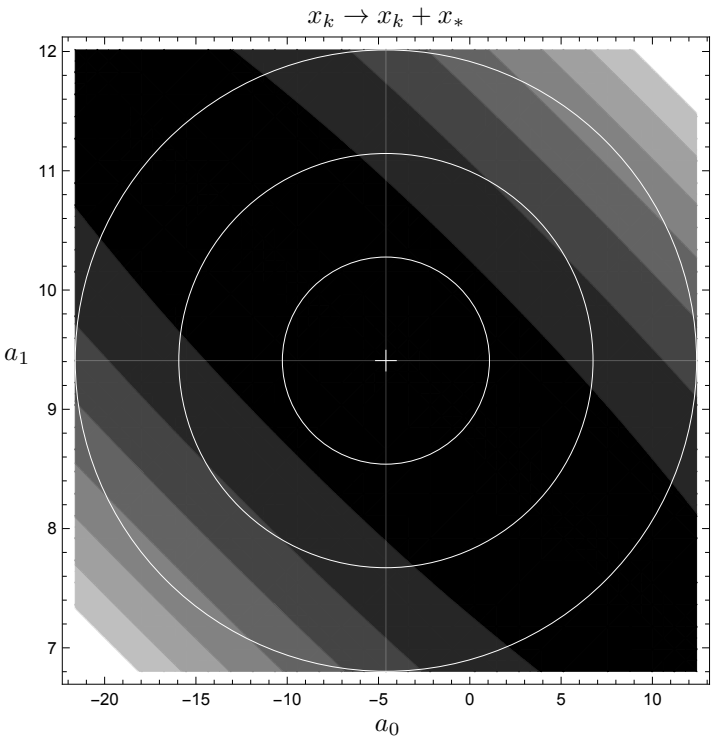


Figure 6.3. Contour plot of merit function after translation along x axis: a_1 is invariant.

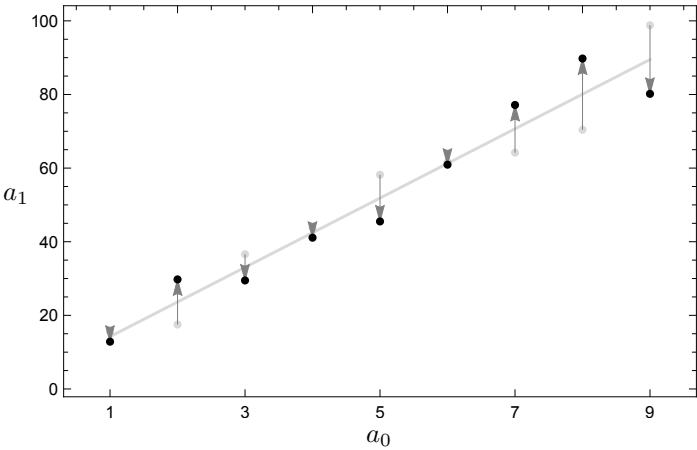


Figure 6.4. The merit function for the learning curve showing the minimum and the value.

Chapter 7

Tricks

7.1 Where Do Parallel Lines Cross?

The provocative question which opens this section has an obvious answer in Euclidean space. There is no such point; parallel lines never cross. Yet, if we input these lines as a linear system, we compute a least squares solution. What is the significance of the least squares solution?

Consider this tantalizing example. The two lines,

$$\begin{aligned}y(x) &= \frac{1}{2} - \frac{1}{2}x, \\y(x) &= 1 - \frac{1}{2}x,\end{aligned}$$

are plotted in figure 7.1 along with the least squares solution is

$$x_{LS} = \mathbf{A}^\dagger \mathbf{b} = \frac{1}{10} \begin{bmatrix} 3 \\ 6 \end{bmatrix}.$$

What is so special about this point?

7.1.1 Intersecting Lines

$$\begin{aligned}y &= 1, \\y &= x.\end{aligned}$$

$$\begin{bmatrix} 0 & 1 \\ -1 & 1 \end{bmatrix} \begin{bmatrix} x \\ y \end{bmatrix} = \begin{bmatrix} 1 \\ 0 \end{bmatrix}$$

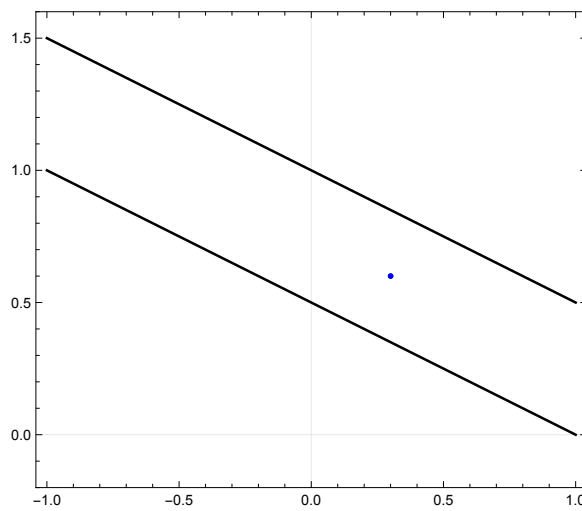


Figure 7.1. *Parallel lines and the least squares solution.*

Part III

Zonal Example

Chapter 8

Zonal Example

8.1 Problem

8.1.1 Zonal Subsection

Chapter 9

Lines

9.1 Face-centered cubic lattice

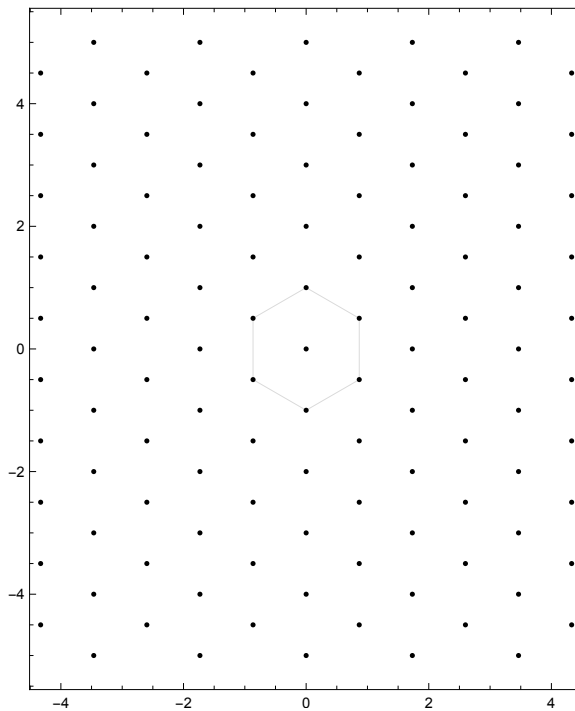


Figure 9.1. *A slice of a face-centered cubic lattice showing a single crystal.*

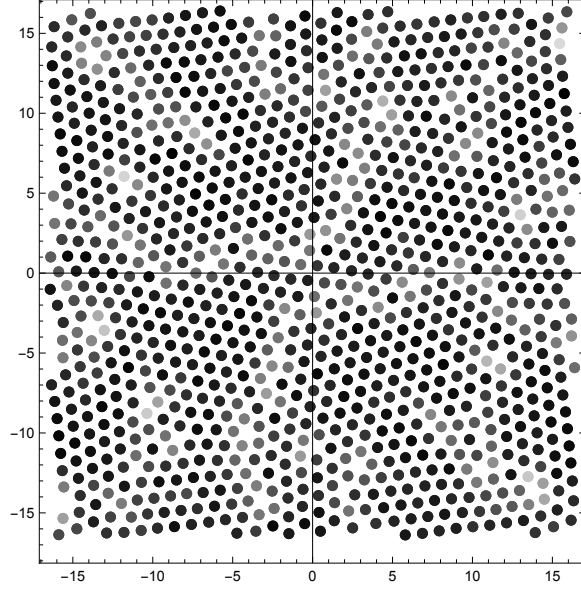


Figure 9.2. *Simulation output showing atomic shades shaded by potential energy.*

9.2 Model

$$y_{(\mu)}(x) = \mu\alpha_* + \alpha_0 + \alpha_1 x, \quad \mu = 0, 1, 2, \dots, M-1. \quad (9.1)$$

$$\begin{cases} 0 \cdot \alpha_* + \alpha_0 + \alpha_1 x_{1_1} = y_{1_1} \\ \vdots \\ 0 \cdot \alpha_* + \alpha_0 + \alpha_1 x_{m_1} = y_{m_1} \end{cases} \quad \text{row 1}$$

$$\begin{cases} 1 \cdot \alpha_* + \alpha_0 + \alpha_1 x_{1_2} = y_{1_2} \\ \vdots \\ 1 \cdot \alpha_* + \alpha_0 + \alpha_1 x_{m_2} = y_{m_2} \end{cases} \quad \text{row 2} \quad (9.2)$$

$$\begin{cases} 2 \cdot \alpha_* + \alpha_0 + \alpha_1 x_{1_3} = y_{1_3} \\ \vdots \\ 2 \cdot \alpha_* + \alpha_0 + \alpha_1 x_{m_3} = y_{m_3} \end{cases} \quad \text{row 3}$$

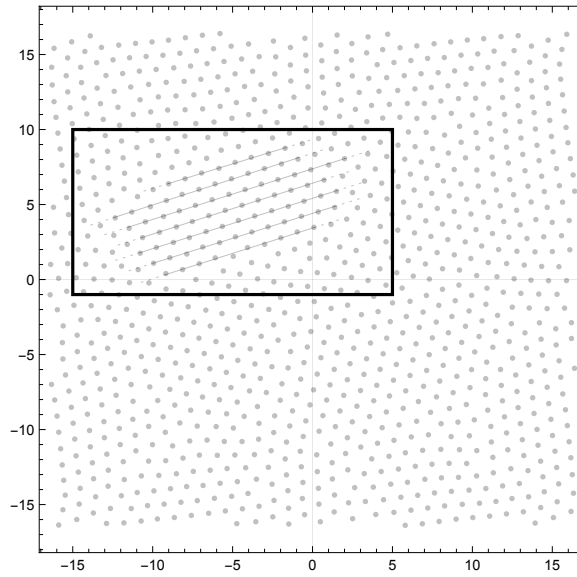


Figure 9.3. *Full data set showing inset.*

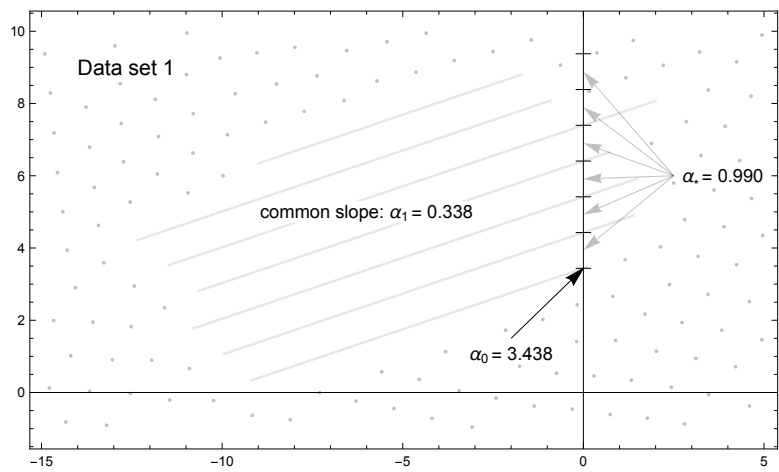
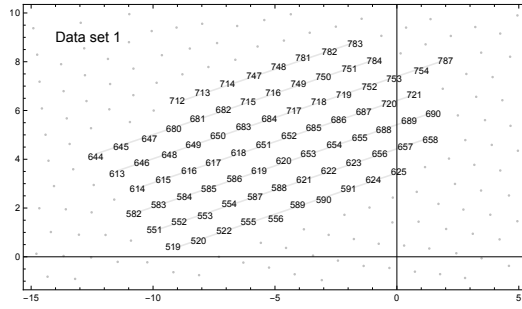
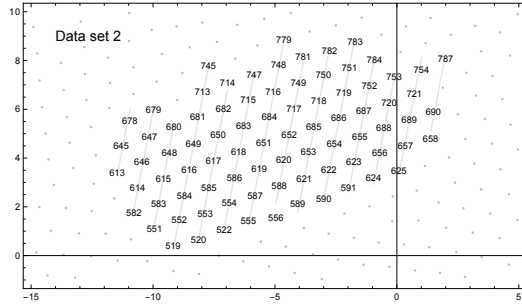


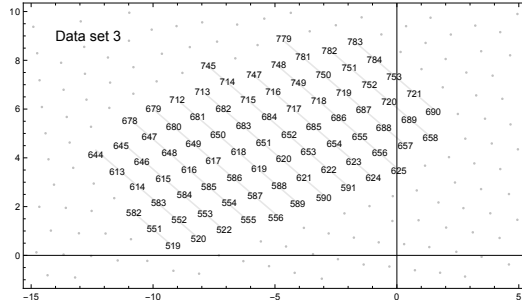
Figure 9.4. *Sample data set showing fit parameters.*

Table 9.1. *Data sets and basic results*

$$\begin{aligned}
 \alpha_* &= 0.9899 \pm 0.0032 \\
 \alpha_0 &= 3.438 \pm 0.013 \\
 \alpha_1 &= 0.3376 \pm 0.0017 \\
 \sqrt{\langle r^2 \rangle} &= 0.052
 \end{aligned}$$



$$\begin{aligned}
 \alpha_* &= 4.974 \pm 0.052 \\
 \alpha_0 &= -2.075 \pm 0.093 \\
 \alpha_1 &= 5.168 \pm 0.052 \\
 \sqrt{\langle r^2 \rangle} &= 0.18
 \end{aligned}$$



$$\begin{aligned}
 \alpha_* &= 1.2322 \pm 0.0039 \\
 \alpha_0 &= -7.505 \pm 0.043 \\
 \alpha_1 &= -0.8576 \pm 0.0038 \\
 \sqrt{\langle r^2 \rangle} &= 0.054
 \end{aligned}$$

$$\begin{array}{c}
\mathbf{A} \\
\left[\begin{array}{ccc}
0 & 1 & x_{1_1} \\
\vdots & \vdots & \vdots \\
0 & 1 & x_{\mu_1} \\
1 & 1 & x_{1_2} \\
\vdots & \vdots & \vdots \\
1 & 1 & x_{\mu_2} \\
2 & 1 & x_{1_3} \\
\vdots & \vdots & \vdots \\
2 & 1 & x_{\mu_3} \\
\vdots & \vdots & \vdots \\
M-1 & 1 & x_{1_M} \\
\vdots & \vdots & \vdots \\
M-1 & 1 & x_{m_M}
\end{array} \right]
\end{array}
\begin{array}{c}
\alpha \\
\left[\begin{array}{c}
\alpha_* \\
\alpha_0 \\
\alpha_1
\end{array} \right]
\end{array}
=
\begin{array}{c}
y \\
\left[\begin{array}{c}
y_{1_1} \\
\vdots \\
y_{\mu_1} \\
y_{1_2} \\
\vdots \\
y_{\mu_2} \\
y_{1_3} \\
\vdots \\
y_{\mu_3} \\
\vdots \\
y_{1_M} \\
\vdots \\
y_{\mu_M}
\end{array} \right]
\end{array}
\quad (9.3)$$

9.3 Solution

Once again the normal equations offer the easy path to solution as in (??). The first step is to compute the inverse of the product matrix. Recall that the dot product is a commutative operator; therefore only six of the nine matrix entries are unique:

$$\mathbf{A}^T \mathbf{A} = \begin{bmatrix} \mathbf{J} \cdot \mathbf{J} & \mathbf{J} \cdot \mathbf{1} & \mathbf{J} \cdot x \\ \mathbf{1} \cdot \mathbf{J} & \mathbf{1} \cdot \mathbf{1} & \mathbf{1} \cdot x \\ x \cdot \mathbf{J} & x \cdot \mathbf{1} & x \cdot x \end{bmatrix} = \begin{bmatrix} a & b & c \\ b & d & e \\ c & e & f \end{bmatrix}.$$

For clarity, the unique elements are specified:

$$\begin{array}{lll}
a = \mathbf{J} \cdot \mathbf{J} & b = \mathbf{J} \cdot \mathbf{1} & c = \mathbf{J} \cdot x \\
& d = \mathbf{1} \cdot \mathbf{J} & e = \mathbf{1} \cdot x \\
& & f = x \cdot x
\end{array}$$

In advance of the computing the inverse, first compute the determinant

$$\det(\mathbf{A}^T \mathbf{A}) = \Delta = 2bce + adf - ae^2 - c^2d - fb^2.$$

Using (??) the inverse is

$$(\mathbf{A}^T \mathbf{A})^{-1} = \Delta^{-1} \begin{bmatrix} df - e^2 & ce - bf & be - cd \\ \cdot & af - c^2 & bc - ae \\ \cdot & \cdot & ad - b^2 \end{bmatrix}.$$

The right-hand side in (9.3) is

$$\mathbf{A}^T y = \beta = \begin{bmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \end{bmatrix} = \begin{bmatrix} \mathbf{J} \cdot y \\ \mathbf{1} \cdot y \\ x \cdot y \end{bmatrix}.$$

The least squares solution is provided as

$$\begin{bmatrix} \alpha_0 \\ \alpha_* \\ \alpha_1 \end{bmatrix} = \left(\mathbf{A}^T \mathbf{A} \right)^{-1} \mathbf{A}^T y$$

which distills down to

$$\alpha = \Delta^{-1} \begin{bmatrix} \beta_1 (df - e^2) + \beta_2 (ce - bf) + \beta_3 (be - cd) \\ \beta_1 (ce - bf) + \beta_2 (af - c^2) + \beta_3 (bc - ae) \\ \beta_1 (be - cd) + \beta_2 (bc - ae) + \beta_3 (ad - b^2) \end{bmatrix}.$$

The errors associate with the fit parameters are

$$\begin{bmatrix} \sigma_* \\ \sigma_0 \\ \sigma_1 \end{bmatrix} = \sqrt{\frac{r^T}{(m-n)\Delta}} \sqrt{\begin{bmatrix} df - e^2 \\ af - c^2 \\ ad - b^2 \end{bmatrix}}.$$

The solutions are expressed in terms of dot products readily available in Fortran.

9.4 Problem Statement

9.5 Data

9.6 Results

9.6.1 Least Squares Results

9.6.2 Apex Angles

9.6.3 Qualitative Results

Table 9.3. *Point membership in data sets shown in figure 9.1.*

set	row	1	2	3	4	5	6	7	8	9	10	11	12
1	1	519	520	522	555	556	589	590	591	624	625		
1	2	551	552	553	554	587	588	621	622	623	656	657	658
1	3	582	583	584	585	586	619	620	653	654	655	688	689
1	4	614	615	616	617	618	651	652	685	686	687	720	721
1	5	613	646	648	649	650	683	684	717	718	719	752	753
1	6	644	645	647	680	681	682	682	715	716	749	750	751
1	7	712	713	714	747	748	781	782	783				
2	1	658	690	787									
2	2	625	657	689	721	754							
2	3	624	656	688	720	753							
2	4	591	623	655	687	752	784						
2	5	590	622	654	686	719	751	783					
2	6	589	621	653	685	718	750	782					
2	7	556	588	620	652	717	749	781					
2	8	555	587	619	651	684	716	748	779				
2	9	522	554	586	618	683	715	747					
2	10	520	553	585	617	650	682	714					
2	11	519	552	584	616	649	681	713	745				
2	12	551	583	615	648	680							
2	13	582	614	646	647	679							
2	14	613	645	678									
3	1	582	551	519									
3	2	644	613	614	583	552	520						
3	3	645	646	615	584	553	522						
3	4	647	648	616	585	554	555	678					
3	5	679	680	649	617	586	587	556					
3	6	712	681	650	618	619	588	589					
3	7	713	682	683	651	620	621	590					
3	8	745	714	715	684	652	653	622	591				
3	9	747	716	717	685	654	623	624					
3	10	748	749	718	686	655	656	625					
3	11	779	781	750	719	687	688	657					
3	12	782	751	752	720	689	658						
3	13	783	784	753	721	690							

Table 9.4. *Excerpted data set.*

	k	x_k	y_k	ϕ_k
	1	-15.879001	-16.365496	-2.597531
	2	-14.749446	-15.995488	-2.613017
	3	-13.905339	16.242941	-2.557543
	\vdots			
	1 022	13.927362	-16.235010	-2.780323
	1 023	14.741765	15.957687	-2.687929
	1 024	15.905518	16.346979	-2.599001

Table 9.5. *Least squares results for three axes.*

axis	gap			intercept			slope			$\sqrt{\langle r^2 \rangle}$
1	0.9899	\pm	0.0032	3.438	\pm	0.013	0.3376	\pm	0.0017	0.052
2	4.974	\pm	0.052	-2.075	\pm	0.093	5.168	\pm	0.052	0.18
3	1.2322	\pm	0.0039	-7.505	\pm	0.043	-0.8576	\pm	0.0038	0.054

Table 9.6. *Intermediate results: angles for the axes.*

axis	θ	\pm	σ_θ							
1	0.3256	\pm	0.0015	=	(18.655	\pm	0.086) $^\circ$			
2	1.380	\pm	0.018	=	(79.0	\pm	1.0) $^\circ$			
3	-0.7089	\pm	0.0025	=	(-40.62	\pm	0.14) $^\circ$			

Table 9.7. *Final results: apex angle measurements*

	θ	\pm	σ_θ							
α	1.040	\pm	0.018	=	(59.6	\pm	1.0) $^\circ$			
β	1.0345	\pm	0.0029	=	(59.27	\pm	0.17) $^\circ$			
γ	1.041	\pm	0.018	=	(59.7	\pm	1.0) $^\circ$			
total	3.116	\pm	0.026	=	(178.5	\pm	1.7) $^\circ$			

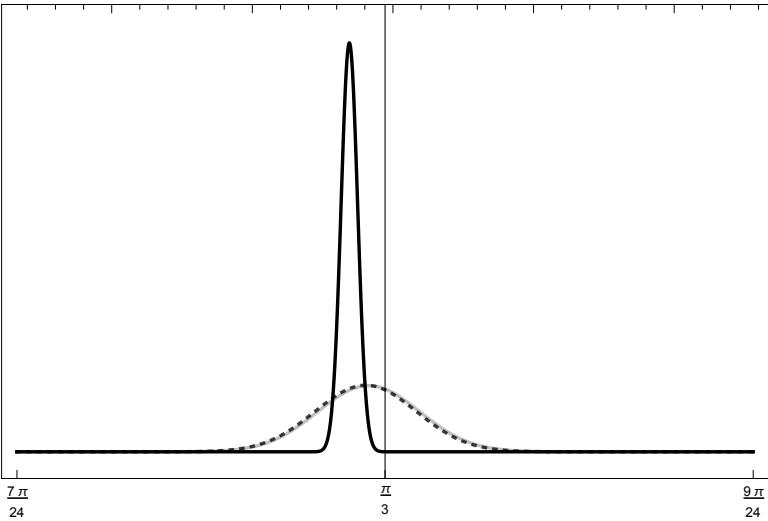


Figure 9.6. *Apex angles displayed in table 9.7.*

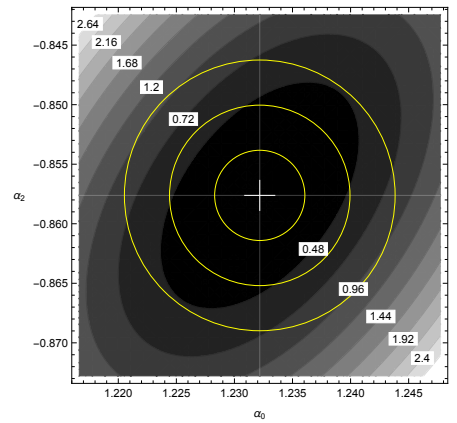
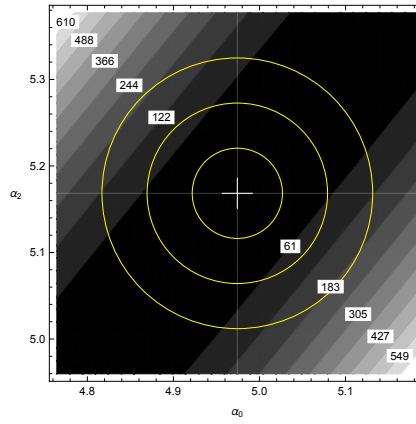
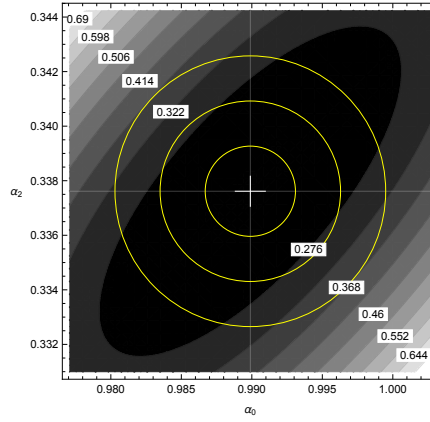


Figure 9.7. Merit functions for the three data sets.

Chapter 10

Crystals

In the previous model, the rows of atoms were treated independently. In this section the basic unit is not a row, it is instead a crystal. Mathematically, the process will imitate Nature: a seed crystal is picked, and other crystals will be identified from that.

Chapter 11

Stitching Local Maps

11.1 What is stitching?

Stitching is the process of combining local maps to create a global map.

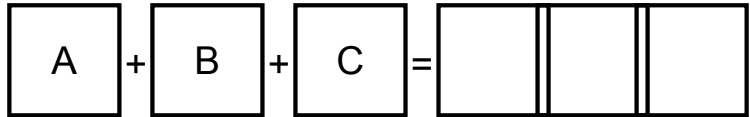


Figure 11.1. *Stitching local maps together to form a global map.*

1. ϕ
2. $\nabla\phi$
3. ϕ and $\nabla\phi$

11.2 Stitch ϕ

11.2.1 Genesis

$$\phi(x) = \exp\left(-\frac{x}{5}\right) \sin(\pi x)$$

11.2.2 Data

The central idea is simple; the mathematical expression is a tedious exercise in index gymnastics.

$$\zeta = 3$$

Table 11.1. *The input data in continuous and discrete form.*

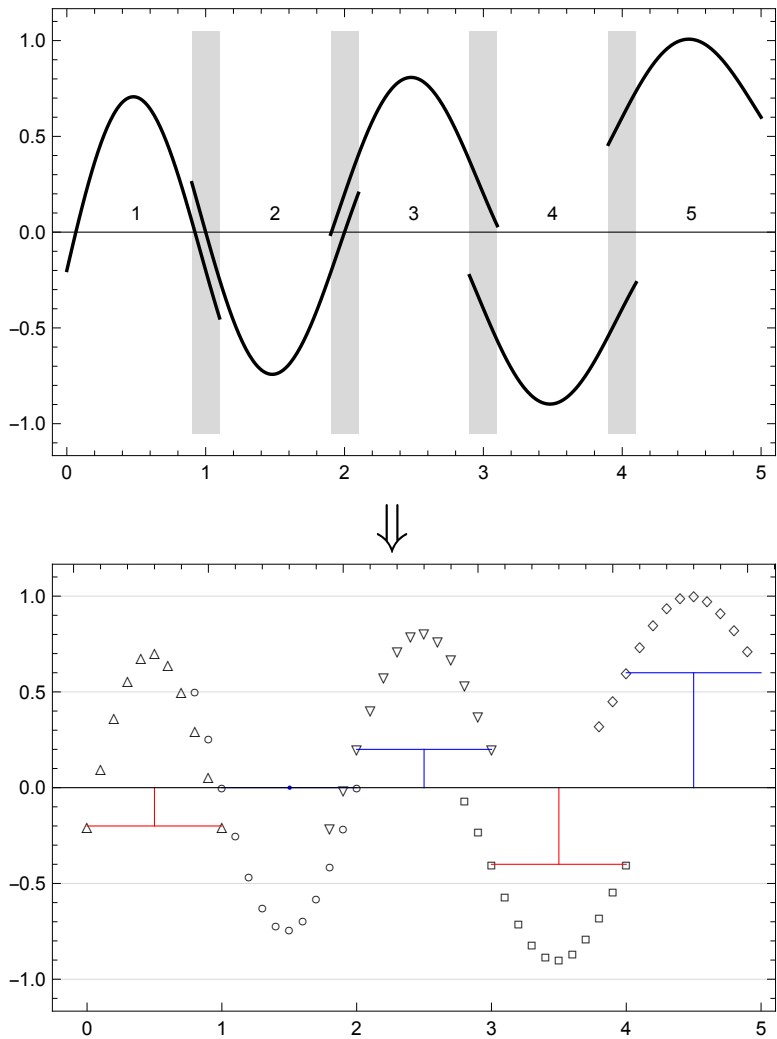


Table 11.2. *Sample showing an overlap of $\zeta = 3$ between the first two zones.*

$\phi_{1,1}$	$\phi_{2,1}$	\cdots	$\phi_{\lambda_1-2,1}$	$\phi_{\lambda_1-1,1}$	$\phi_{\lambda_1,1}$				
			$\phi_{1,2}$	$\phi_{2,2}$	$\phi_{3,2}$	\cdots	$\phi_{\lambda_1-2,2}$	$\phi_{\lambda_2-1,2}$	$\phi_{\lambda_2,2}$

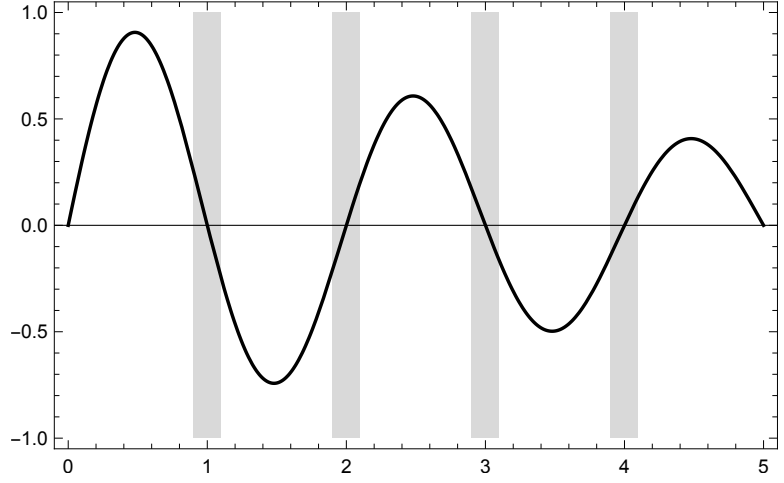


Figure 11.2. The ideal potential function showing five measurement zones and four overlap bands.

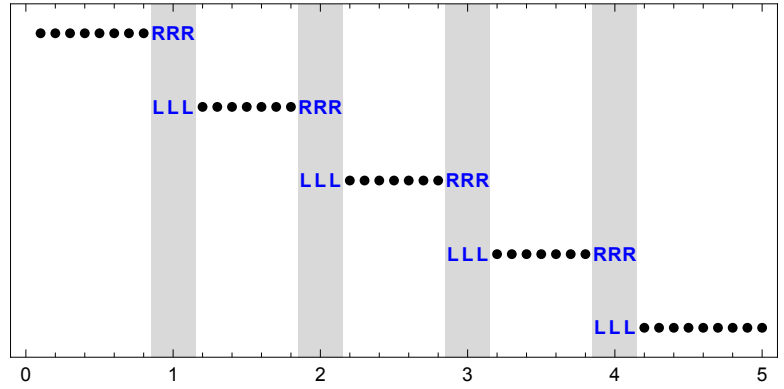


Figure 11.3. Waterfall diagram showing discretization within measurement zones with left and right zone overlaps.

$$\Delta_{12} = \zeta^{-1} \left(\underbrace{(\phi_{\lambda_1-2,1} + \phi_{\lambda_1-1,1} + \phi_{\lambda_1,1})}_{\text{zone 1}} - \underbrace{(\phi_{1,2} + \phi_{2,2} + \phi_{3,2})}_{\text{zone 2}} \right)$$

$$\Delta_{12} = \zeta^{-1} \left(\underbrace{(\phi_{\lambda_1-2,1} - \phi_{1,2})}_{\text{pair 1}} + \underbrace{(\phi_{\lambda_1-1,1} - \phi_{2,2})}_{\text{pair 2}} + \underbrace{(\phi_{\lambda_1,1} - \phi_{3,2})}_{\text{pair 3}} \right)$$

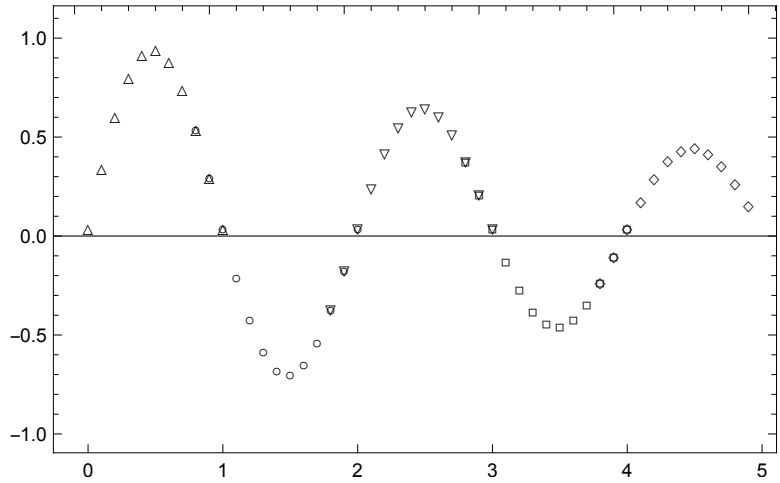


Figure 11.4. *Stitching unifies the data.*

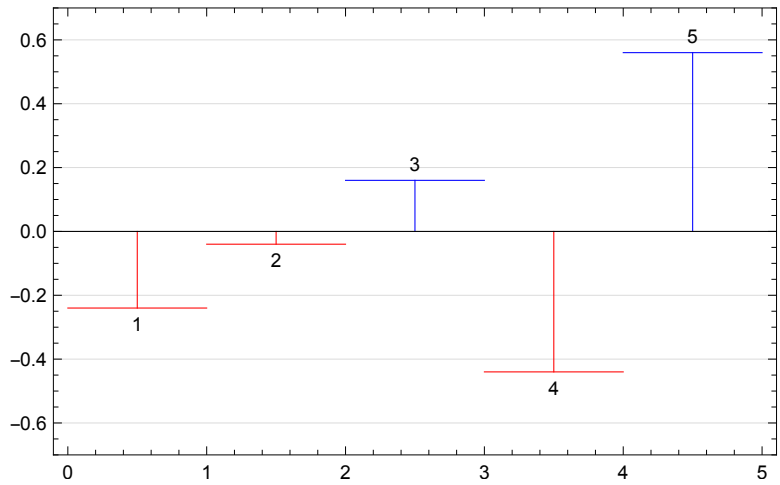


Figure 11.5. *A set of piston adjustments which restores continuity across the domain.*

Mean value of the differences.

$$\Delta_{j,j+1} = \zeta^{-1} \sum_{k=1}^{\zeta} p_{j,\lambda_j - \zeta + k} - p_{j+1,k}$$

k	y_1	y_2	y_3	y_4	y_5
1	-0.2	0.500 878	-0.210 084	-0.064 251 7	0.325 113
2	0.102 898	0.258 113	-0.011 324 8	-0.226 982	0.458 345
3	0.364 738	0.0	0.2	-0.4	0.6
4	0.561 904	-0.247 992	0.403 039	-0.566 234	0.736 101
5	0.677 936	-0.462 368	0.578 555	-0.709 935	0.853 753
6	0.704 837	-0.623 794	0.710 719	-0.818 142	0.942 345
7	0.643 511	-0.718 793	0.788 498	-0.881 821	0.994 482
8	0.503 326	-0.740 818	0.806 531	-0.896 585	1.006 57
9	0.300 878	-0.690 609	0.765 423	-0.862 929	0.979 014
10	0.058 112 7	-0.575 834	0.671 453	-0.785 993	0.916 025
11	-0.2	-0.410 084	0.535 748	-0.674 887	0.825 059
12		-0.211 325	0.373 018	-0.541 655	0.715 978
13		0.0	0.2	-0.4	

Table 11.3. *Measurements displaying the connection between overlap bands in figure 11.3.*

First overlap region.

$$\Delta_{12} = \zeta^{-1} \left(\underbrace{(\phi_{9,1} + \phi_{10,1} + \phi_{11,1})}_{\text{last 3 elements of zone 1}} - \underbrace{(\phi_{1,2} + \phi_{2,2} + \phi_{3,2})}_{\text{first 3 elements of zone 2}} \right)$$

11.2.3 Data and results

Table 11.4. *Computation of the zone shift values.*

$$\begin{aligned} \Delta_{12} &= \frac{1}{3} ((\phi_{9,1} + \phi_{10,1} + \phi_{11,1}) - (\phi_{1,2} + \phi_{2,2} + \phi_{3,2})) \\ \Delta_{23} &= \frac{1}{3} ((\phi_{11,2} + \phi_{12,2} + \phi_{13,2}) - (\phi_{1,3} + \phi_{2,3} + \phi_{3,3})) \\ \Delta_{34} &= \frac{1}{3} ((\phi_{11,3} + \phi_{12,3} + \phi_{13,3}) - (\phi_{1,4} + \phi_{2,4} + \phi_{3,4})) \\ \Delta_{45} &= \frac{1}{3} ((\phi_{11,4} + \phi_{12,4} + \phi_{13,4}) - (\phi_{1,5} + \phi_{2,5} + \phi_{3,5})) \end{aligned}$$

11.2.4 Linear System

$$\mathbf{A}p = \Delta$$

Table 11.5. *Computation of the zone shift values.*

$$\begin{aligned}
\Delta_{12} &= \frac{1}{3} ((0.300878 + 0.0581127 - 0.2) - (0.500878 + 0.258113 + 0.)) \\
\Delta_{23} &= \frac{1}{3} ((-0.410084 - 0.211325 + 0.) - (-0.210084 - 0.0113248 + 0.2)) \\
\Delta_{34} &= \frac{1}{3} ((0.535748 + 0.373018 + 0.2) - (-0.0642517 - 0.226982 - 0.4)) \\
\Delta_{45} &= \frac{1}{3} ((-0.674887 - 0.541655 - 0.4) - (0.325113 + 0.458345 + 0.6))
\end{aligned}$$

Table 11.6. *Input data*

	Shift	Value
1	Δ_{12}	-0.2
2	Δ_{23}	0.0
3	Δ_{34}	0.6
4	Δ_{45}	-1.

$$\begin{bmatrix} 1 & -1 & 0 & 0 & 0 \\ 0 & 1 & -1 & 0 & 0 \\ 0 & 0 & 1 & -1 & 0 \\ 0 & 0 & 0 & 1 & -1 \end{bmatrix} \begin{bmatrix} p_1 \\ p_2 \\ p_3 \\ p_4 \\ p_5 \end{bmatrix} = \begin{bmatrix} \Delta_{12} \\ \Delta_{23} \\ \Delta_{34} \\ \Delta_{45} \end{bmatrix}$$

$$p_{LS} = \frac{1}{5} \begin{bmatrix} 4 & 3 & 2 & 1 \\ -1 & 3 & 2 & 1 \\ -1 & -2 & 2 & 1 \\ -1 & -2 & -3 & 1 \\ -1 & -2 & -3 & -4 \end{bmatrix} \begin{bmatrix} \Delta_{12} \\ \Delta_{23} \\ \Delta_{34} \\ \Delta_{45} \end{bmatrix} + \alpha \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix}$$

$$\mathbf{A}^\dagger \mathbf{b} = \frac{1}{25} \begin{bmatrix} -6 \\ -1 \\ 4 \\ -11 \\ 14 \end{bmatrix}$$

These are the actual plot values used in figure 11.5.

$$\Phi_{corrected} = \Phi_{measured} - \mathbf{A}^\dagger \mathbf{b}$$

Table 11.7. *Problem statement for linear regression.*

trial function	$p_k - p_{k+1} = \Delta_{k,k+1}, k = 1: n$
merit function	$M(p) = \sum_{k=1}^n (\Delta_{k,k+1} - p_k + p_{k+1})^2$
number of zones	$m = 5$
number of overlaps	$n = 4$
rank defect	$m - n = 1$
measurements per zone	$\lambda = \{11, 13, 13, 13, 12\}$
measurements	$\phi_{k,j}, k = 1: m, j = 1: \lambda_m$
input data	$\Delta_{k,k+1}, k = 1: n$
results	$p_k, k = 1: m$
residual error	$r = \mathbf{A}^\dagger \mathbf{b} - \Delta$
linear system	$\begin{bmatrix} 1 & -1 & 0 & 0 & 0 \\ 0 & 1 & -1 & 0 & 0 \\ 0 & 0 & 1 & -1 & 0 \\ 0 & 0 & 0 & 1 & -1 \end{bmatrix} \begin{bmatrix} p_1 \\ p_2 \\ p_3 \\ p_4 \\ p_5 \end{bmatrix} = \begin{bmatrix} \Delta_{12} \\ \Delta_{23} \\ \Delta_{34} \\ \Delta_{45} \end{bmatrix}$
gauge condition	$\sum_{k=1}^m p_k = 0$

11.2.5 Least Squares Arbitration

There is a fundamental ambiguity arising from gradient measurements stemming from the basic fact that

$$\frac{d}{dx}\phi(x) = \frac{d}{dx}(\phi(x) + c).$$

We can recover the function shape, but not the offset. In other words, there is a translation invariance. This lone constant is the poster child for the rank one deficiency in the linear system of (11.2.4). Realizing this, the one dimensional problem could be solved without resort to least squares.

The system can be solved, for example, by moving from left to right and manually forcing the data to match. If the the overlap difference between zone 1 and zone 2 is Δ_{12} , add Δ_{12} to every value in zone 2. Now zones 1 and 2 are stitched together. Compute Δ_{23} , add this value to every point in zone 3. Zones 1, 2, and 3 are now stitched together. Continue as needed.

The least squares problem is obviated. How did this happen? The process of least squares is an exercise error arbitration which takes a peanut butter approach by trying to distribute the error evenly. In one dimension, there is no need for arbitration as there is no conflict in measurements.

In two dimensions, the problem changes. Consider the typical cell with a neighbor to the right and a neighbor above. The right-left overlap adjustment conflicts with the up-down overlap adjustment. The least squares process takes

Table 11.8. *Results for stitching with piston.*

fit parameters	$p_k, k = 1: m$	pistons
computed solution	$p = \frac{1}{25} \begin{bmatrix} -6 \\ -1 \\ 4 \\ -11 \\ 14 \end{bmatrix}$	$\mathbf{A}^\dagger b$
data vector	$\frac{1}{5} \begin{bmatrix} -1 \\ 0 \\ 1 \\ -2 \\ 3 \end{bmatrix} = \frac{1}{25} \begin{bmatrix} -6 \\ -1 \\ 4 \\ -11 \\ 14 \end{bmatrix} - \frac{1}{5} \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix}$	$p_{\mathcal{R}} + p_{\mathcal{N}}$
residual error	$r \cdot r = 0$	
problem statement	table 11.7	
measurements	table 11.3	
input data	table 11.6	
plots	figure 11.2 figure 11.4	raw data (bottom) corrected data

all off the overlap conflicts and provides a set of adjustments which minimizes the global error. To close, note that the least squares solution was used even though it is not necessary until dimension 2 or higher.

One last tidbit. Figure 11.7 shows the piston values that were input to distort the values. Least squares chooses a distinct set of corrections. Why was this set selected? A tantalizing clue is given by the null space vector in (11.2.4). Notice this vector is perpendicular to every column vector in \mathbf{A}^\dagger which implies that the sum of each column vector must be 0. Therefore, the gauge condition is that the solution vector will have sum 0:

$$p_1 + p_2 + p_3 + p_4 + p_5 = 0.$$

We may now eliminate a variable; choose the last one:

$$p_5 = -p_1 - p_2 - p_3 - p_4$$

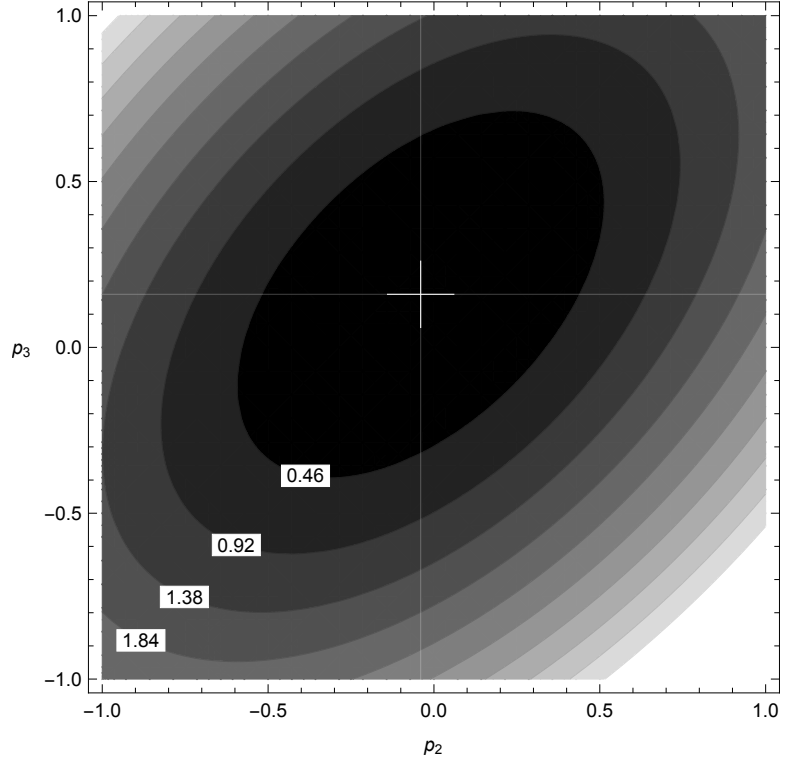


Figure 11.6. Looking at the merit function on the $p_2 - p_3$ axis.

Instead of (11.2.4), there is now

$$\begin{bmatrix} 1 & -1 & 0 & 0 \\ 0 & 1 & -1 & 0 \\ 0 & 0 & 1 & -1 \\ 1 & 1 & 1 & 2 \end{bmatrix} \begin{bmatrix} p_1 \\ p_2 \\ p_3 \\ p_4 \end{bmatrix} = \begin{bmatrix} \Delta_{12} \\ \Delta_{23} \\ \Delta_{34} \\ \Delta_{45} \end{bmatrix}.$$

The solution is the same:

$$\hat{p}_{gauge} = \begin{bmatrix} 4 & 3 & 2 & 1 \\ -1 & 3 & 2 & 1 \\ -1 & -2 & 2 & 1 \\ -1 & -2 & -3 & 1 \end{bmatrix} \begin{bmatrix} \Delta_{12} \\ \Delta_{23} \\ \Delta_{34} \\ \Delta_{45} \end{bmatrix}.$$

The 0 sum, or equivalently 0 mean, condition is a gauge condition which restores the column rank of the problem.

The piston values used to create the data set are decomposed into range and

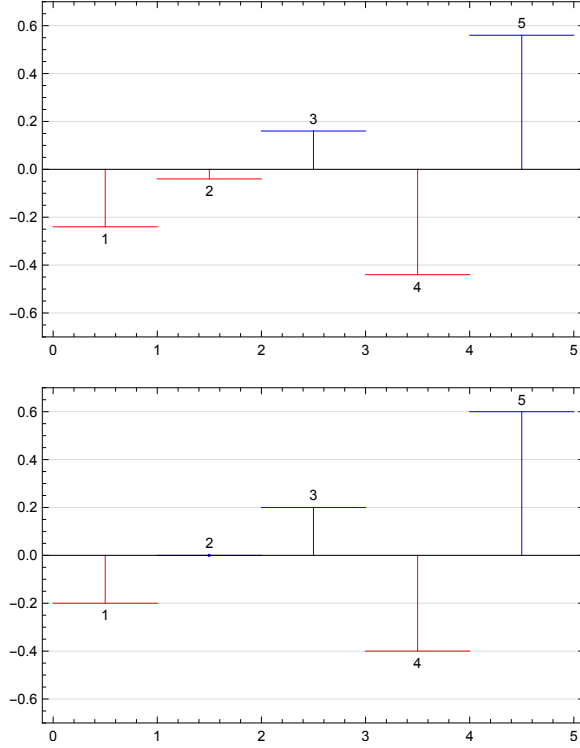


Figure 11.7. *On top, pistons output from the solution; on bottom, pistons input to create the data.*

null space terms.

$$\frac{1}{5} \begin{bmatrix} -1 \\ 0 \\ 1 \\ -2 \\ 3 \end{bmatrix} = \frac{1}{25} \begin{bmatrix} -6 \\ -1 \\ 4 \\ -11 \\ 14 \end{bmatrix} - \frac{1}{5} \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix}$$

11.3 Stitch $\nabla\phi$

The next challenge is to stitch data together using the gradient $\nabla\phi$ rather than the function value ϕ . The outputs now will be a set of piston adjustments called tilts which restore continuity of the gradient.

The problem arises in the field of wavefront sensing. Modern devices make exquisite measurements of tilts. The process of wavefront reconstruction takes these tilts and reconstructs the wavefront. Measure $\nabla\phi(x)$ and compute $\nabla\phi(x)$.

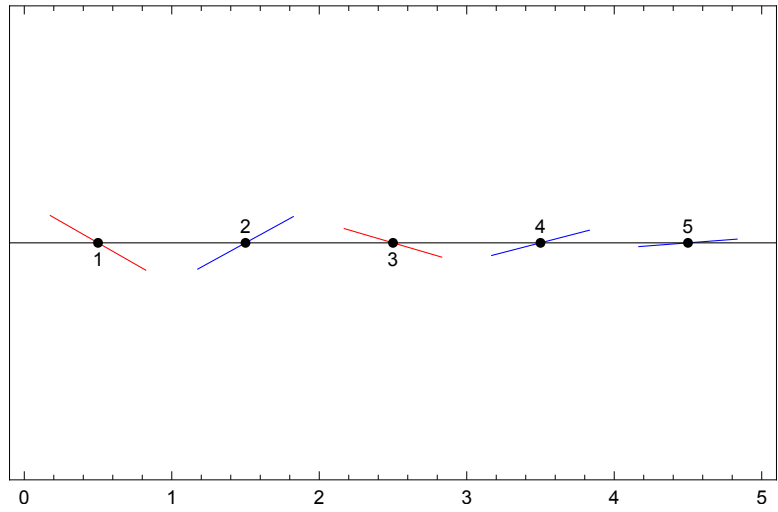


Figure 11.8. A set of tilt adjustments which restores continuity of the gradient across the domain.

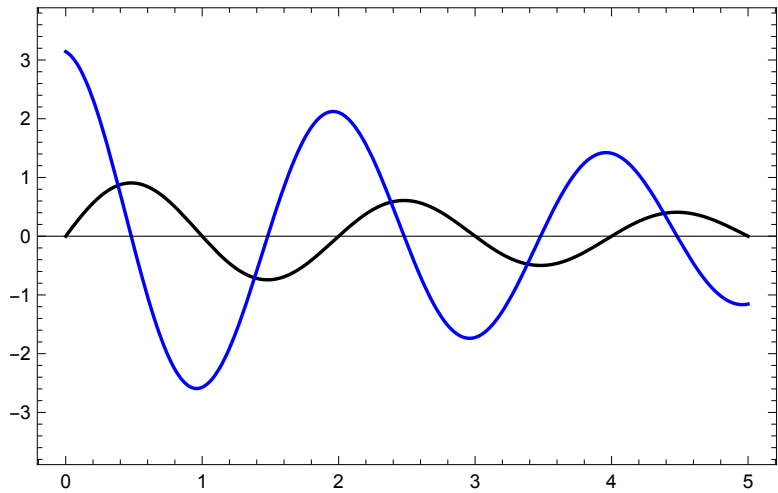


Figure 11.9. A function (black) and its gradient (blue).

Gradient of (11.3)

$$\nabla\phi(x) = \frac{1}{5} \exp\left(-\frac{x}{5}\right) (5\pi \cos(\pi z) - \sin(\pi z))$$

$$\tau = \frac{1}{100} \begin{bmatrix} -90 \\ 85 \\ -40 \\ 35 \\ 10 \end{bmatrix}$$

A scaled version of these values is plotted in figure 11.8.

Chapter 12

Gradient I

$$F = \nabla \phi$$

$$W^{1,2}(\Omega) = \{\phi \in L^2(\Omega) : \partial_x^1 \phi \in L^2(\Omega)\}$$

12.1 One Dimension

$$\Omega = \bigcup_k \omega_k$$

Interval

$$\omega = \{x \in \mathbb{R} : a < x < b\}$$

Average gradient

$$\langle \nabla \phi(x) \rangle_\omega = \phi(b) - \phi(a)$$

$$\begin{bmatrix} -1 & 1 & 0 & \dots & & \\ 0 & -1 & 1 & & & \\ \vdots & & \ddots & \ddots & & \vdots \\ & & & -1 & 1 & 0 \\ & & & 0 & -1 & 1 \end{bmatrix} \begin{bmatrix} \varphi_0 \\ \varphi_1 \\ \vdots \\ \varphi_{m-1} \\ \varphi_m \end{bmatrix} = \begin{bmatrix} x_1 \\ \vdots \\ x_m \end{bmatrix}$$

Part IV

Applications: Nonlinear Problems

Chapter 13

Linearization

13.1 Powers Laws and Exponentials

Exponential relationships, such as radioactive decay,

$$y = ae^{bx},$$

and power laws, like the learning curve,

$$y = ax^b \tag{13.1}$$

are inherently nonlinear. As such, they are outside of the scope of this work. But the problem is that the Internet has helped create and maintain the myth that such problems can be tamed and made linear. This is, empathically, not true.

So this chapter is about the proper way to handle power law and exponential relationships. Along the way, it will become clear why faux linearization does not work. In short, the answer is that the linearization *changes the problem*. That is, it produces a different problem with a different solution.

$$a_{LS} = \left\{ a \in \mathbb{R}^2: \|y_k - ax^b\|_2^2 \text{ is minimized} \right\}$$

13.2 Learning Curve

The trial function

$$y(x) = ax^b, \quad x \in \mathbb{R}^+.$$

If we restrict $x \in \mathbb{R}^+$, then $b \in \mathbb{R}$, avoiding the indeterminate form 0^0 . In practice $x \in \mathbb{N}$. The variable a is an initial value (at $x = 1$), and b is a rate parameter.

The merit function is

$$M(a, b) = \sum_{k=1}^m (y_k - ak^b)^2 \tag{13.2}$$

k	y_k	$y(k)$	r_k	k	y_k	$y(k)$	r_k
1	419.115	371.9	47.2149	27	119.695	128.685	-8.98984
2	251.263	297.505	-46.2419	28	131.931	127.187	4.74458
3	291.329	261.092	30.2371	29	126.702	125.758	0.944397
4	208.272	237.992	-29.7202	30	107.923	124.392	-16.4689
5	198.577	221.492	-22.9146	31	121.735	123.086	-1.35074
6	172.583	208.863	-36.2804	32	136.803	121.834	14.969
7	202.107	198.749	3.35859	33	97.0772	120.633	-23.5554
8	169.911	190.384	-20.4736	34	110.727	119.479	-8.75175
9	175.674	183.299	-7.62479	35	132.09	118.369	13.7215
10	191.393	177.185	14.2084	36	94.4017	117.3	-22.8979
11	152.023	171.83	-19.8067	37	111.213	116.269	-5.05649
12	183.7	167.082	16.6184	38	113.38	115.275	-1.89502
13	157.806	162.831	-5.02491	39	112.433	114.315	-1.88186
14	142.933	158.991	-16.0584	40	123.706	113.387	10.319
15	185.178	155.498	29.6797	41	125.563	112.489	13.0742
16	172.109	152.3	19.8089	42	101.05	111.619	-10.5695
17	174.763	149.356	25.4069	43	101.805	110.777	-8.9723
18	151.21	146.632	4.57821	44	107.492	109.96	-2.46808
19	132.162	144.101	-11.9391	45	89.7489	109.167	-19.4182
20	125.188	141.741	-16.5524	46	107.598	108.397	-0.799046
21	144.023	139.531	4.49148	47	117.91	107.649	10.261
22	117.048	137.457	-20.4084	48	94.22	106.922	-12.7018
23	125.011	135.503	-10.4919	49	108.116	106.214	1.90216
24	144.994	133.659	11.3349	50	108.17	105.526	2.64485
25	126.14	131.914	-5.77349	51	116.083	104.855	11.2277
26	146.929	130.258	16.6713	52	123.93	104.201	19.7289

To advance the discussion, consider the sample set of data for 52 points in table 13.2. What does the surface of the merit function in (13.2) look like in solution space? Very much like all the other merit functions so far. There is a unique minimum, a distinct feature of the landscape. The question becomes how to find it.

13.2.1 Problem Statement

The problem stated in table 13.1 embodies the major points discussed up to this point and displays a few details that will become clear as the discussion advances.

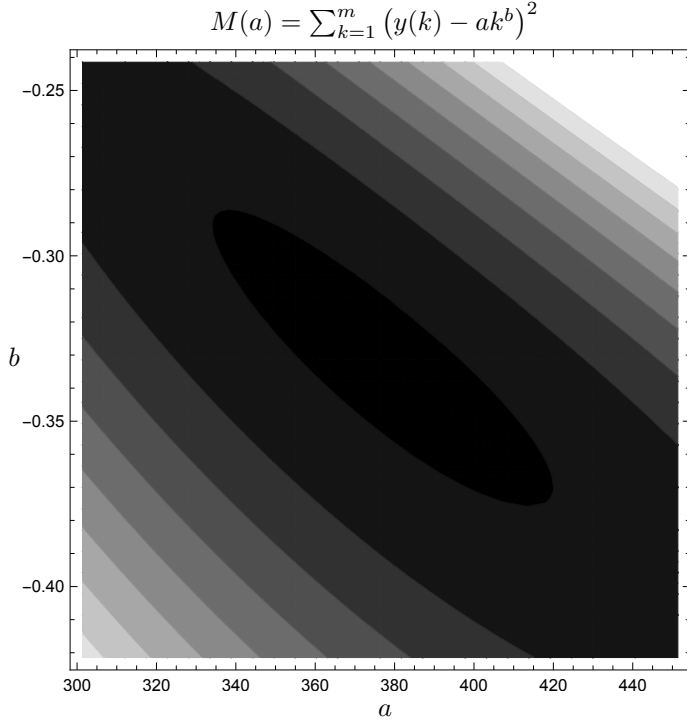


Figure 13.1. Merit function for the learning curve in solution space.

13.2.2 Solution

Without a linear system, attack with calculus as in §3.2.2. Find the roots of the gradient $\nabla M(a, b) = 0$:

$$\begin{aligned}\partial_a M &= -2 \sum_{k=1}^m (y_k - ak^b) k^b = 0, \\ \partial_b M &= -2 \sum_{k=1}^m (y_k - ak^b) ak^b \ln k = 0.\end{aligned}$$

Look at the equations in explicit form in table 13.2. There is no linear system which isolates the solution parameters in a fashion such as this:

$$\mathbf{A} \begin{bmatrix} a \\ b \end{bmatrix} = \zeta.$$

The condition $\partial_a M = 0$ allows the separation of the solution parameters.

$$a(b) = \frac{\sum y_k k^b}{\sum k^{2b}}$$

This permits removal of one variable, and reduces this merit function to a single

Table 13.1. *Problem statement for learning curve.*

trial function	$y(x) = ae^{b \ln x}$	$a \in \mathbb{R}^+, b \in \mathbb{R}$
residual error	$r_k = y(k) - ak^b$	
merit function	$M(a, b) = \sum_{k=1}^m (y(x) - ak^b)^2$	
unit numbers	$k, k = 1:m$	
measurements	$y_k, k = 1:m$	
results	a	initial value
	b	learning rate
# of measurements	$m = 52$	
# of parameters	$n = 2$	
system matrix	no system matrix – problem is nonlinear	
linear system	no linear system – problem is nonlinear	
unperturbed solution	$\begin{bmatrix} a \\ b \end{bmatrix} = \begin{bmatrix} 371.9 \\ -0.322 \end{bmatrix}$	80% rate
input data	table 13.2	

Table 13.2. *The simultaneous conditions defining $\nabla M(a, b) = 0$.*

$\partial_a M$		$\partial_b M$	
$a1^{2b}$	$= y_1 1^b$	$a^2 1^{2b}$	$= ay_1 1^b \ln 1$
$a2^{2b}$	$= y_2 2^b$	$a^2 2^{2b}$	$= ay_2 2^b \ln 2$
	\vdots		\vdots
$a52^{2b}$	$= y_{52} 52^b$	$a^{52} 52^{2b}$	$= ay_{52} 52^b \ln 52$

parameter, b :

$$\hat{M}(b) = \sum \left(y(x) - \frac{\sum y_k k^b}{\sum k^{2b}} k^b \right)^2$$

(13.3)

where summations from $k = 1$ to 52 are implied.

Figure 13.2 shows the plot of $\hat{M}(b)$ with the minimum marked. The nonlinear least squares problem in two dimensions is now reduced to a minimization problem in one dimension. The solutions are posted in table 13.1. The curve in figure 13.2 represents a slice through the surface of the merit function in figure 13.1. Such a slice is shown in figure 13.3.

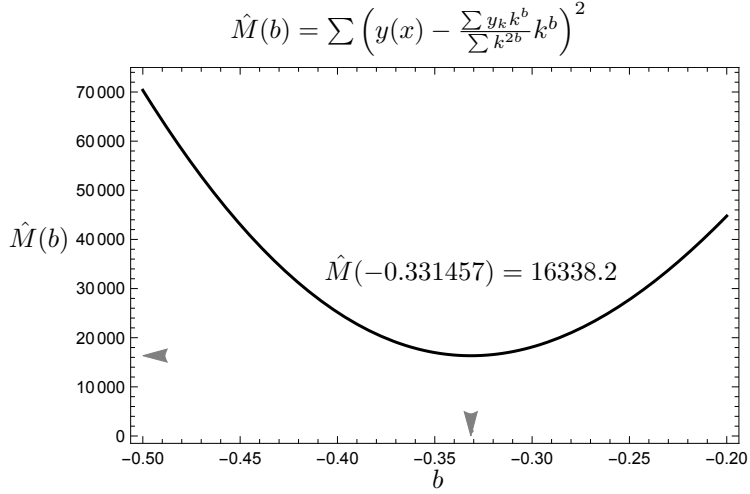


Figure 13.2. Merit function constrained to one parameter b .

13.2.3 Results

The minimizer is quoted to full double precision for debugging purposes:

$$b_{LS} = -0.33145698171782384,$$

the initial value is

$$a(b_{LS}) = 376.3700704535057,$$

and the minimum of the merit function is

$$M(a_{LS}, b_{LS}) = \hat{M}(b_{LS}) = 16338.235321721559.$$

The results and residuals are posted in figure 13.2.

The merit function in figure 13.8 has a new feature. In addition to the least squares solution and the error ellipses, the unperturbed solution is plotted with and “ \times ”. This is the ideal solution used to generate the raw data before the noise is added. The least squares solution will converge to this point as the magnitude of the error decreases.

13.3 What Not To Do

Folklore promotes the idea of transforming the problem to a more amenable form. And certainly transformation is a powerful and useful mathematical technique. But the popular transformation complicates the problem and provides an incorrect answer.

13.3.1 Logarithmic Transform

While the logarithmic transformation is exact,

$$y = ax^b \quad \Rightarrow \quad \ln y = \ln a + b \ln x,$$

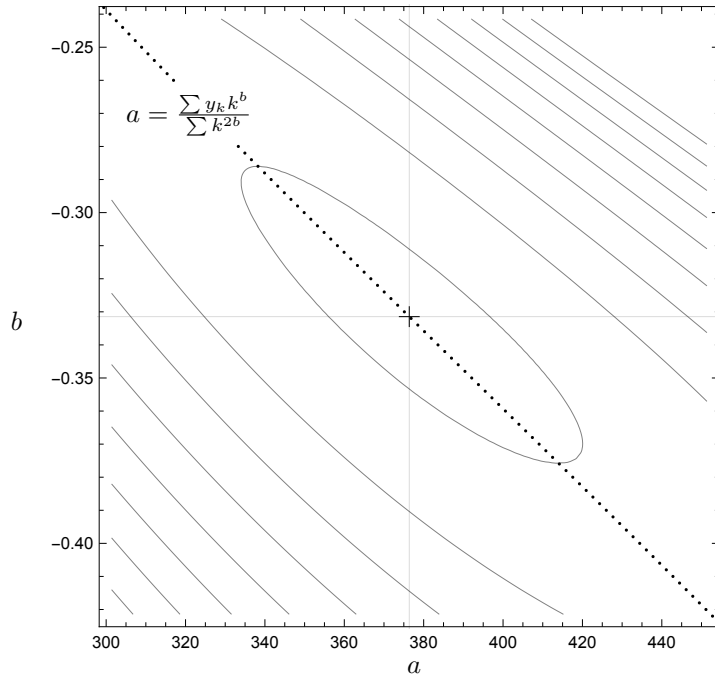


Figure 13.3. Merit function for the learning curve in solution space showing the constrained a parameter as a dotted line.

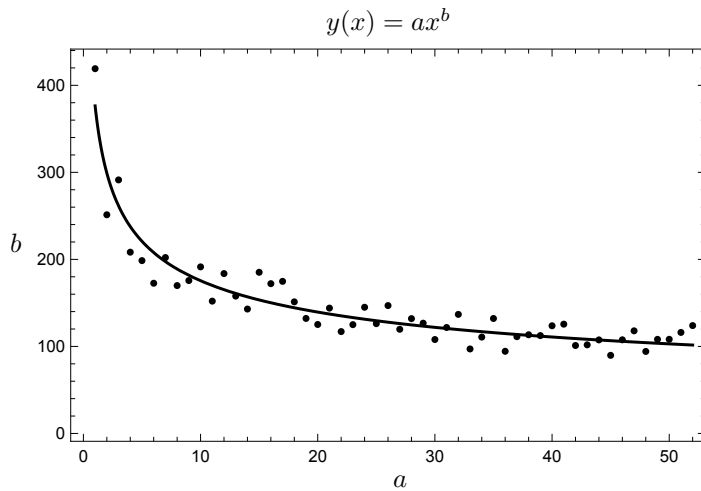


Figure 13.4. Solution for equations (13.2) and (13.3) using data in table 13.2.

the application is inappropriate because the problem is not exact. If there is no noise in the data set, the above transformation will preserve the solution. However, if the problem is exact, there is no need to employ least squares. One could take

Table 13.3. *Results for learning curve analysis.*

fit parameters	a	initial value
	b	rate
solution function	$y(x) = ax^b$	$[L]$
solution error		
computed solution	$\begin{bmatrix} a \\ b \end{bmatrix} = \begin{bmatrix} 376.37 \\ -0.331457 \end{bmatrix} \pm \begin{bmatrix} ? \\ ? \end{bmatrix}$	
unperturbed solution	$\begin{bmatrix} \tilde{a} \\ \tilde{b} \end{bmatrix} = \begin{bmatrix} 376 \\ -0.322 \end{bmatrix}$	
r^*r	16 338	
problem statement	table 13.1	
input data	table 13.2	
plots	figure 13.4	1. data and solution
	figure 13.5	2. residual errors
	figure 13.8	3. merit function

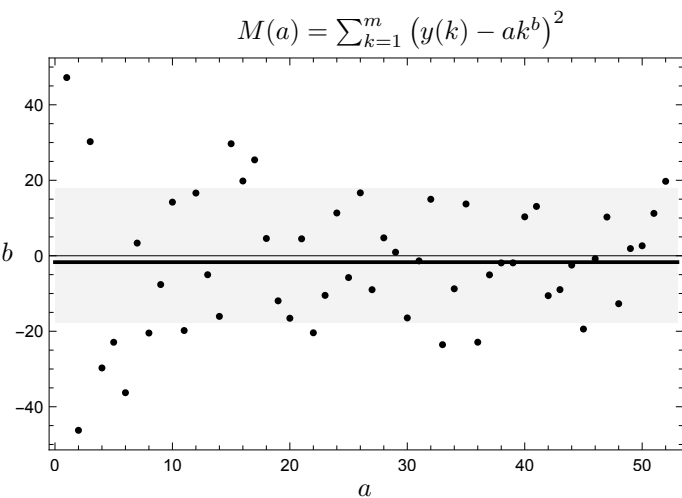


Figure 13.5. *The residual errors in figure 13.4.*

the ratio of two distinct data points,

$$\frac{y_j}{y_k} = \left(\frac{j}{k}\right)^b,$$

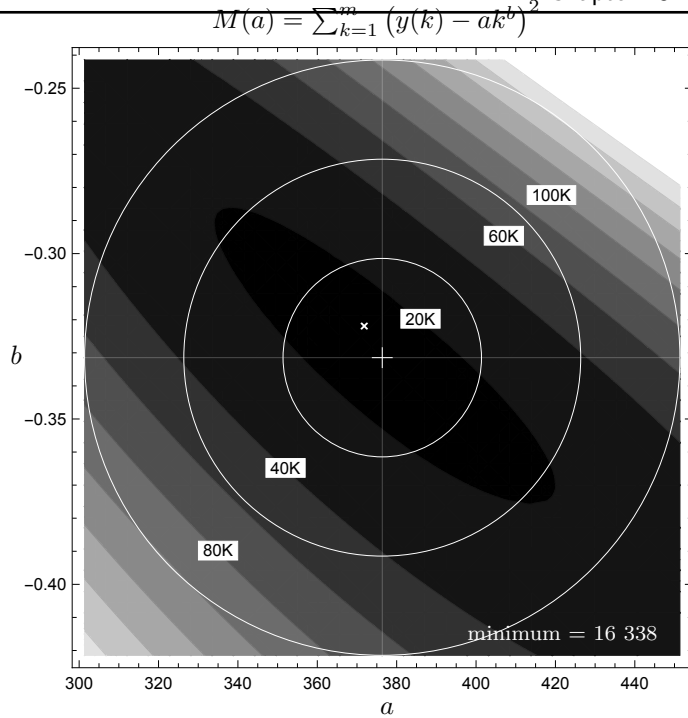


Figure 13.6. Minimization of the merit function for the learning curve.

and compute

$$b = \frac{\ln(y_j/y_k)}{\ln(j/k)}.$$

Again, there is no need for least squares methods.

When the data contains noise, and the choice is to use least squares, the logarithmic transformation corrupts the data. Look at the transformation of trial function in the presence of noise, that is, with nonzero residual error:

$$y_k = ak^b + r_k \quad \Rightarrow \quad \ln y_k = \ln(ak^b + r_k).$$

The logarithmic transform is not linear and does not separate the parameters a and b in the presence of error.

Because this malady is so common, the solution invites belaboring. The function $\ln y = \ln a + b \ln x$ is linear in the new coordinates $\ln y$ and $\ln x$ but only when there is no noise – only when there is no need for least squares arbitration. The presence of noise creates an additive term and the logarithmic transformation creates a different problem which has a different solution.

13.3.2 Linear Transformation

A transformation T is linear if and only if

$$T(x + \alpha y) = T(x) + \alpha T(y).$$

The logarithmic transformation fails this primal test:

$$\ln(x + \alpha y) \neq \ln(x) + \alpha \ln(y)$$

The faux solution is

$$\begin{aligned} \ln a = 5.88094 & \quad \rightarrow \quad a = 358.146, \\ b = -0.314929. \end{aligned} \tag{13.4}$$

Note that the value of the merit function is higher than the true minimum:

$$M(358.146, -0.314929) = 17005 > 16338. \tag{13.5}$$

The eyeball is a poor instrument for distinguishing between the correct solution in table 13.6 and (13.4). However, the merit function sharply delineates the two.

The merit shows the solution (13.4) marked with \boxplus , distinctly away from the minimum of the merit function as seen in equation (13.5) and figure 13.8.

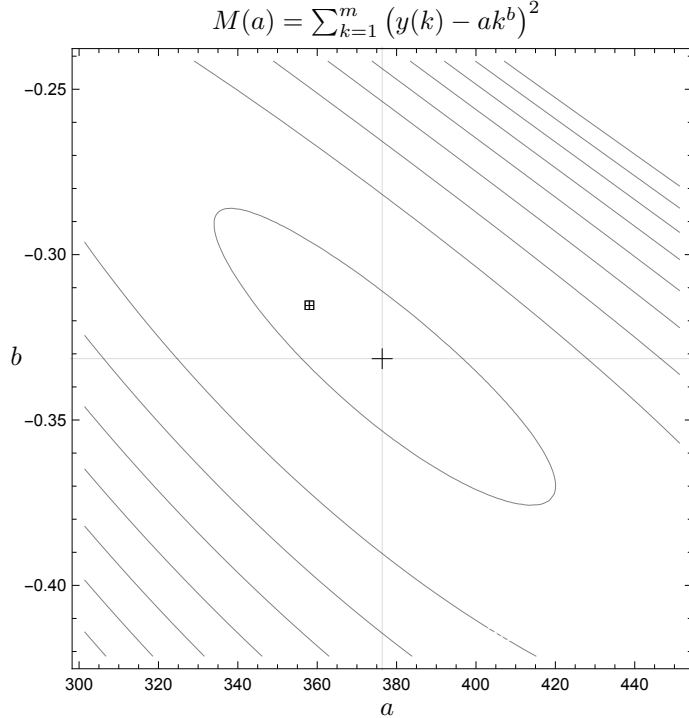


Figure 13.7. The merit function for the learning curve showing the minimum and the value .

13.3.3 Reflection Test Fails

13.4 Radioactive Decay

13.4.1 Theory

13.4.2 Problem Statement

Table 13.4. Problem statement for radioactive decay.

trial function	$c(t) = ae^{bt}$	counts
residual error	$r_k = c_k - ae^{bt_k}$	counts
merit function	$M(a) = \sum_{k=1}^m (c_k - ae^{bt_k})^2$	counts ²
measurements	$t_k, k = 1:m$	time, s
	$c_k, k = 1:m$	counts
results	$a \pm \epsilon_a$	initial value
	$b \pm \epsilon_b$	power
# of measurements	$m = 10$	rows in A
# of parameters	$n = 2$	columns in A
system matrix	$\mathbf{A} \in \mathbb{R}_2^{10 \times 2}$	
linear system	$\begin{bmatrix} 1 & x_1 \\ \vdots & \vdots \\ 1 & x_m \end{bmatrix} \begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \begin{bmatrix} T_1 \\ \vdots \\ T_m \end{bmatrix}$	
ideal solution	$\begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \begin{bmatrix} 0 \\ 10 \end{bmatrix}$	
input data	table 3.2	

$$C = C_0e^{-t/\tau}$$

13.4.3 Results

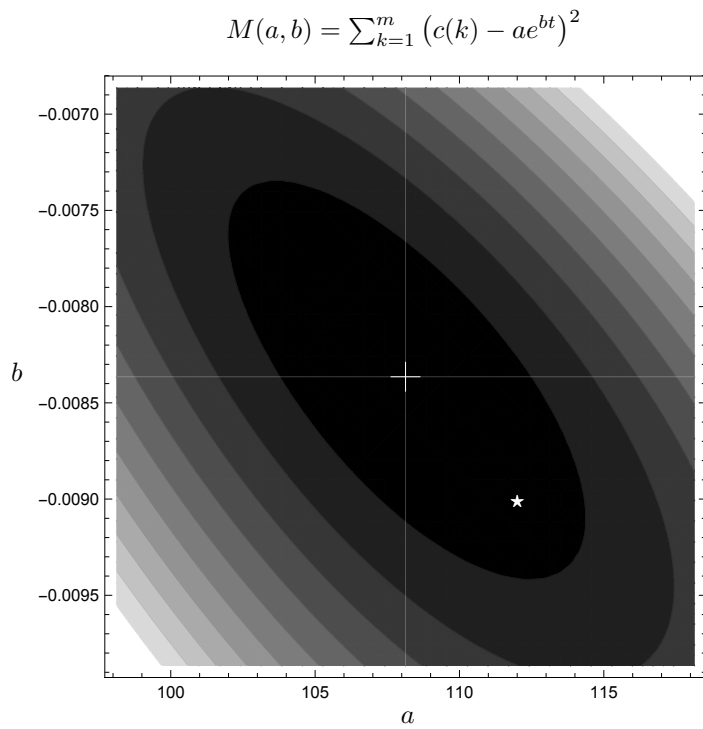
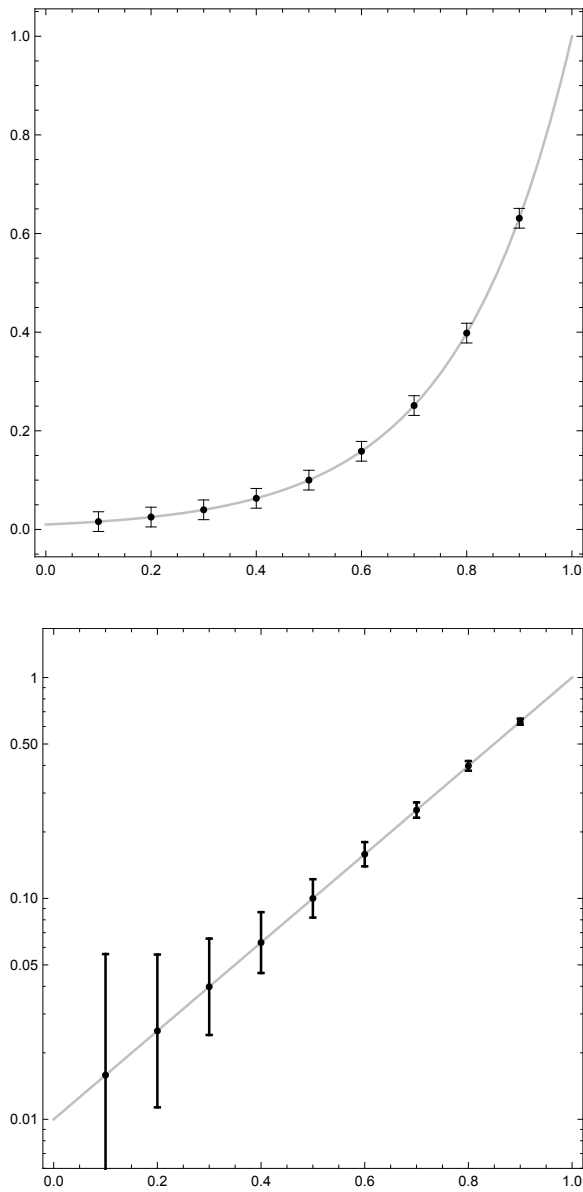


Figure 13.8. *The merit function for the learning curve showing the minimum and the value .*

Table 13.5. *Logarithmic scaling distorts errors.*



<i>k</i>	<i>t</i>	<i>c</i>
1	0	106
2	15	80
3	30	98
4	45	75
5	60	74
6	75	73
7	90	49
8	105	38
9	120	37
10	135	22

Table 13.6. Results for radioactive decay.

fit parameters	a_0	intercept, °C
	a_1	slope, °C / cm
solution function	$T(x) = a_0 + a_1x$	°C
solution error	$\epsilon_T^2(x) = \epsilon_0^2 + x^2\epsilon_1^2 + a_1^2\epsilon_x^2$	°C
computed solution	$\begin{bmatrix} a_0 \\ a_1 \end{bmatrix} = \begin{bmatrix} 4.8 \\ 9.41 \end{bmatrix} \pm \begin{bmatrix} 4.9 \\ 0.87 \end{bmatrix}$	
ideal solution	$\begin{bmatrix} \tilde{a}_0 \\ \tilde{a}_1 \end{bmatrix} = \begin{bmatrix} 0 \\ 10 \end{bmatrix}$	
r^*r	316.6	
curvature matrix $(\mathbf{A}^*\mathbf{A})^{-1}$	$\frac{1}{180} \begin{bmatrix} 95 & -15 \\ -15 & 3 \end{bmatrix}$	
problem statement	table 13.4	
input data	table 3.2	
plots	figure 3.2	1. data and solution
	figure 3.3	2. residual errors
	figure ??	3. merit function

Chapter 14

Population Growth

In this section we take a nonlinear model for population growth and separate the linear and nonlinear terms.

14.1 Model

$$y(\tau) = a_1 + a_2\tau + a_3e^{d\tau} \quad (14.1)$$

$$\mathbf{A}(d + \gamma) \neq \mathbf{A}(d) + \mathbf{A}(\gamma)$$

$$\begin{array}{ccc} \mathbf{A}(d) & a & = y \\ \left[\begin{array}{ccc} 1 & \tau_1 & e^{d\tau_1} \\ 1 & \tau_2 & e^{d\tau_2} \\ 1 & \tau_3 & e^{d\tau_3} \\ 1 & \tau_4 & e^{d\tau_4} \\ 1 & \tau_5 & e^{d\tau_5} \\ 1 & \tau_6 & e^{d\tau_6} \\ 1 & \tau_7 & e^{d\tau_7} \\ 1 & \tau_8 & e^{d\tau_8} \end{array} \right] & \left[\begin{array}{c} a_0 \\ a_1 \\ a_3 \end{array} \right] & = \left[\begin{array}{c} y_0 \\ y_1 \\ y_3 \\ y_4 \\ y_5 \\ y_6 \\ y_7 \\ y_8 \end{array} \right] \end{array}$$

$$\min_{\substack{a \in \mathbb{R}^3 \\ d \in \mathbb{R}}} \left\| \mathbf{A}(d) \left[\begin{array}{c} a_1 \\ a_2 \\ a_3 \end{array} \right] - y \right\|_2^2 \quad (14.2)$$

Table 14.1. *Problem statement for population model with linear and exponential growth.*

trial function	$y(\tau) = a_0 + a_1\tau + a_2e^{d\tau}$	$a \in \mathbb{R}^3$ $d \in \mathbb{R}$
merit function	$M(a, d) = \sum_{k=1}^m (y_k - a_1 + a_2\tau + a_3e^{d\tau_k})^2$	
# measurements	$m = 8$	
# parameters	$n = 4$	
rank defect	$\rho = n$	overdetermined
input data	$(\tau_k, y_k), k = 1: 8$	table 14.2
results	a_0 a_1 a_2 d	constant linear exponential power term
residual error	$r = \textcolor{blue}{A}^\dagger b - \Delta$	
linear system	$\textbf{A}(d) \begin{bmatrix} a_1 \\ a_2 \\ a_3 \end{bmatrix} = y$	

14.2 Problem Statement

$$r^2 = \left\| \textbf{A}(d) \begin{bmatrix} a_1 \\ a_2 \\ a_3 \end{bmatrix} - y \right\|_2^2$$

(14.3)

14.3 Data

14.4 Example

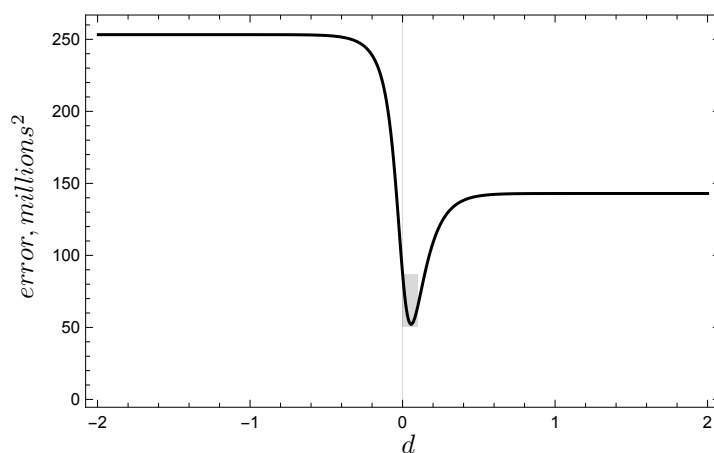
year = 1900 + 10($\tau - 1$)

14.5 Polynomials

There is the model we choose and the model which nature chooses. Are they the same?

Table 14.2. *Data v. prediction.*

year	census	fit	r	rel. error
1900	76.00	77.51	1.51	2.0%
1910	91.97	90.98	-0.99	-1.1%
1920	105.71	104.87	-0.84	-0.8%
1930	122.78	119.48	-3.29	-2.7%
1940	131.67	135.36	3.69	2.8%
1950	150.70	153.46	2.76	1.8%
1960	179.32	175.45	-3.87	-2.2%
1970	203.24	204.26	1.029	0.5%

**Figure 14.1.** *Residual error for (14.3) with a_1 , a_2 , and a_3 at optimal values.*

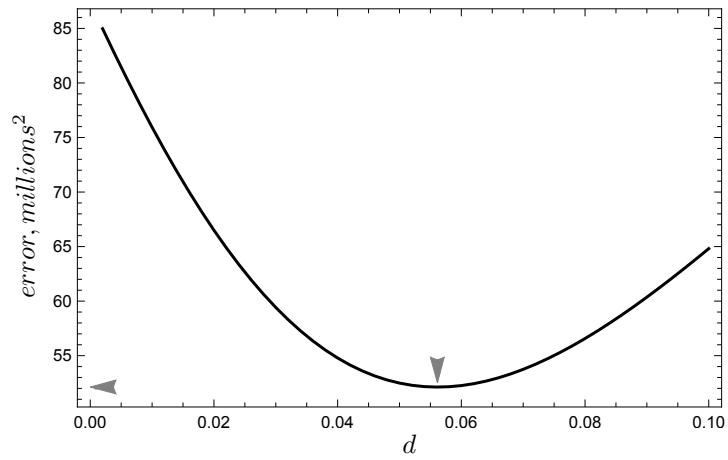


Figure 14.2. Residual error for a_1 and a_2 fixed at optimal values.

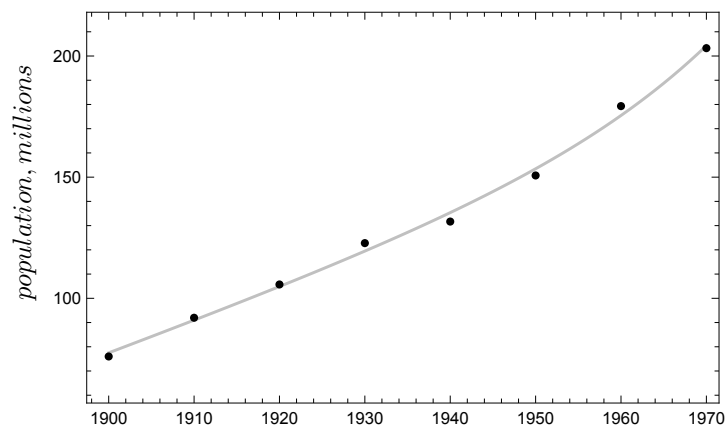


Figure 14.3. Solution plotted against data.

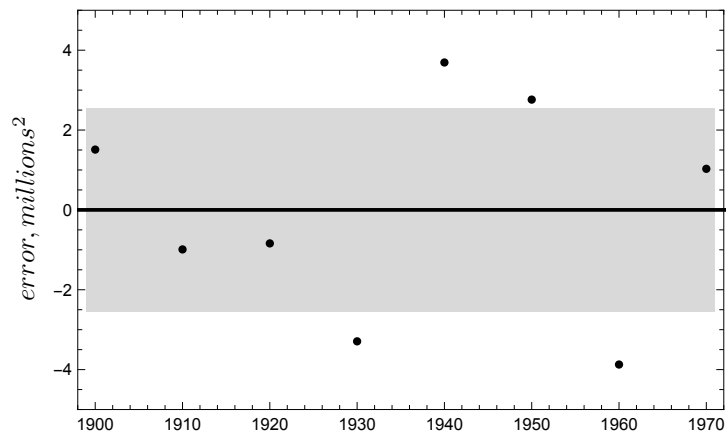


Figure 14.4. Scatterplot of residual errors.

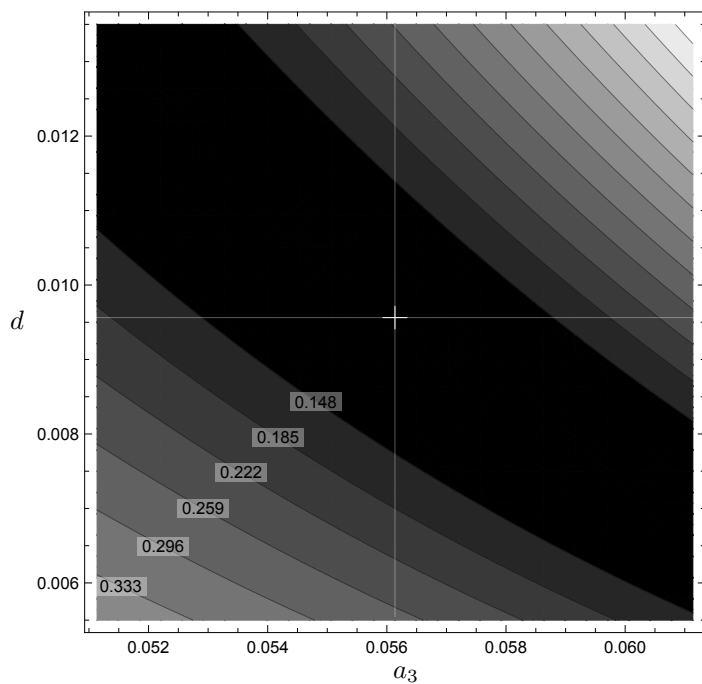


Figure 14.5. The merit function with a_1 and a_2 fixed at best values showing least squares solution (center cross).

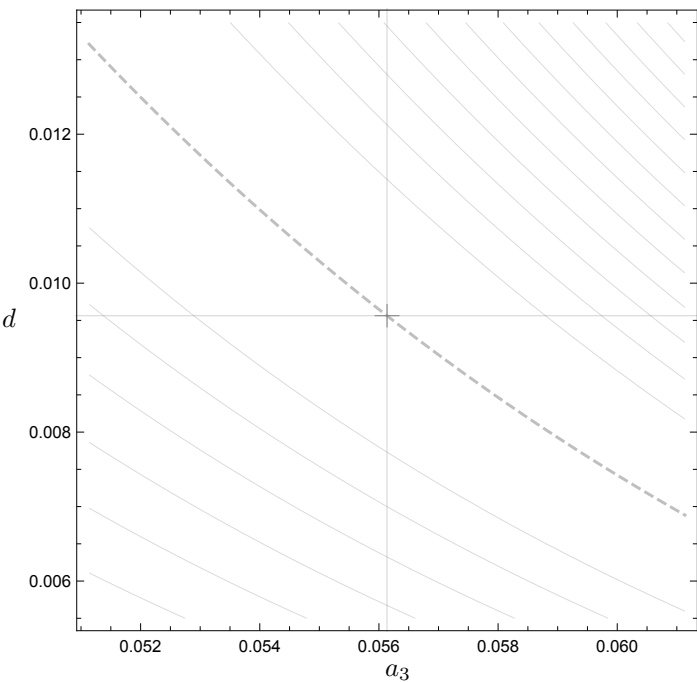


Figure 14.6. The merit function with a_1 and a_2 fixed at best values showing least squares solution (center cross) and the (dashed line).

Table 14.3. Results for census analysis

fit parameters	$c = \begin{bmatrix} 0.010 \\ 0.0170 \\ 0.0096 \end{bmatrix} \pm \begin{bmatrix} 0.031 \\ 0.0014 \\ 0.0020 \end{bmatrix}$
	$d = 0.056136 \pm ??$
r^*r	55.12
$\sum r_k$	-6.2×10^{-14}
σ_r	2.55
a	$\begin{bmatrix} 0.5397 & -0.0188 & 0.0165 \\ -0.0188 & 0.0011 & -0.0014 \\ 0.0165 & -0.0014 & 0.0022 \end{bmatrix}$
plots	data vs fit: figure 14.3 residuals: figure 14.4 merit function in $\mathcal{R}(\mathbf{A}^*)$: figure 14.6

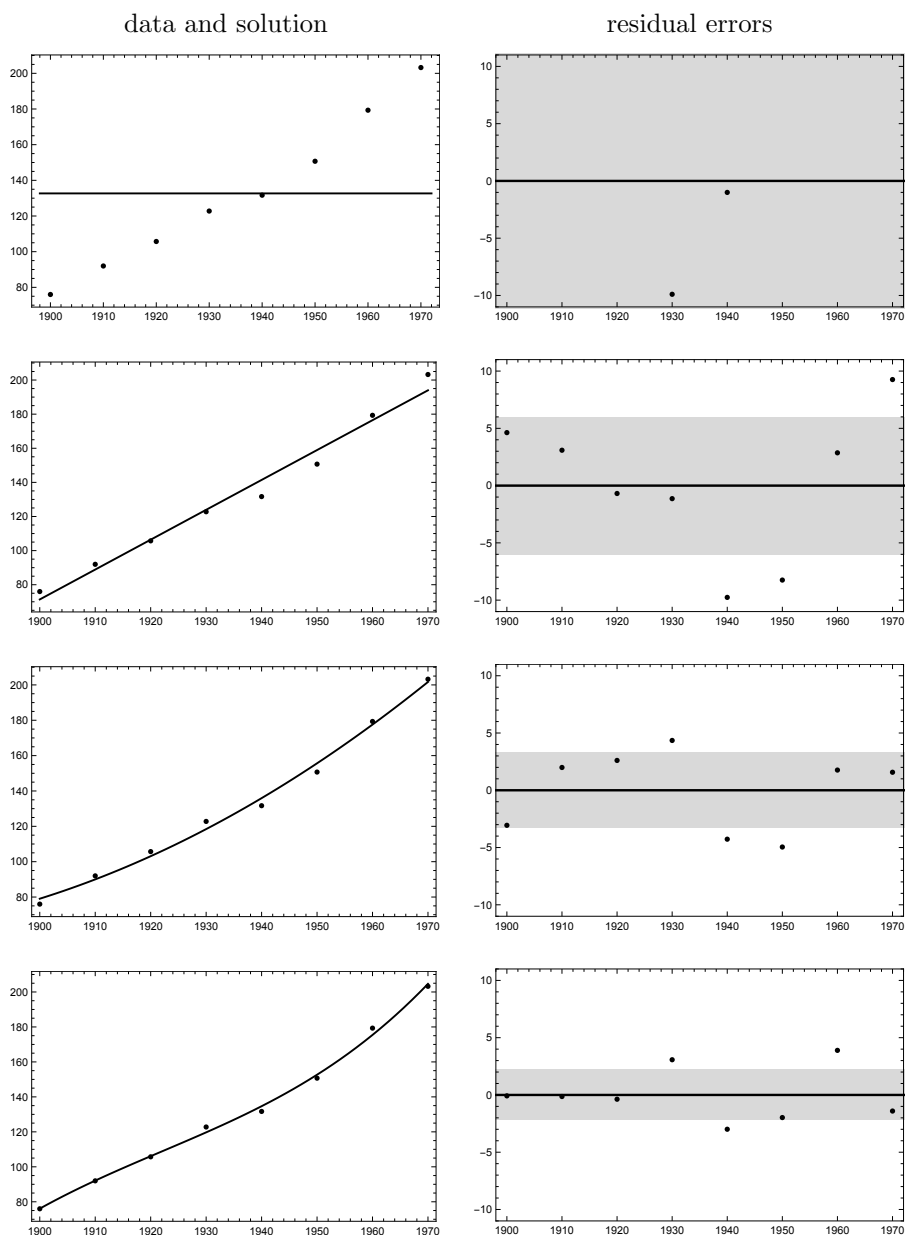
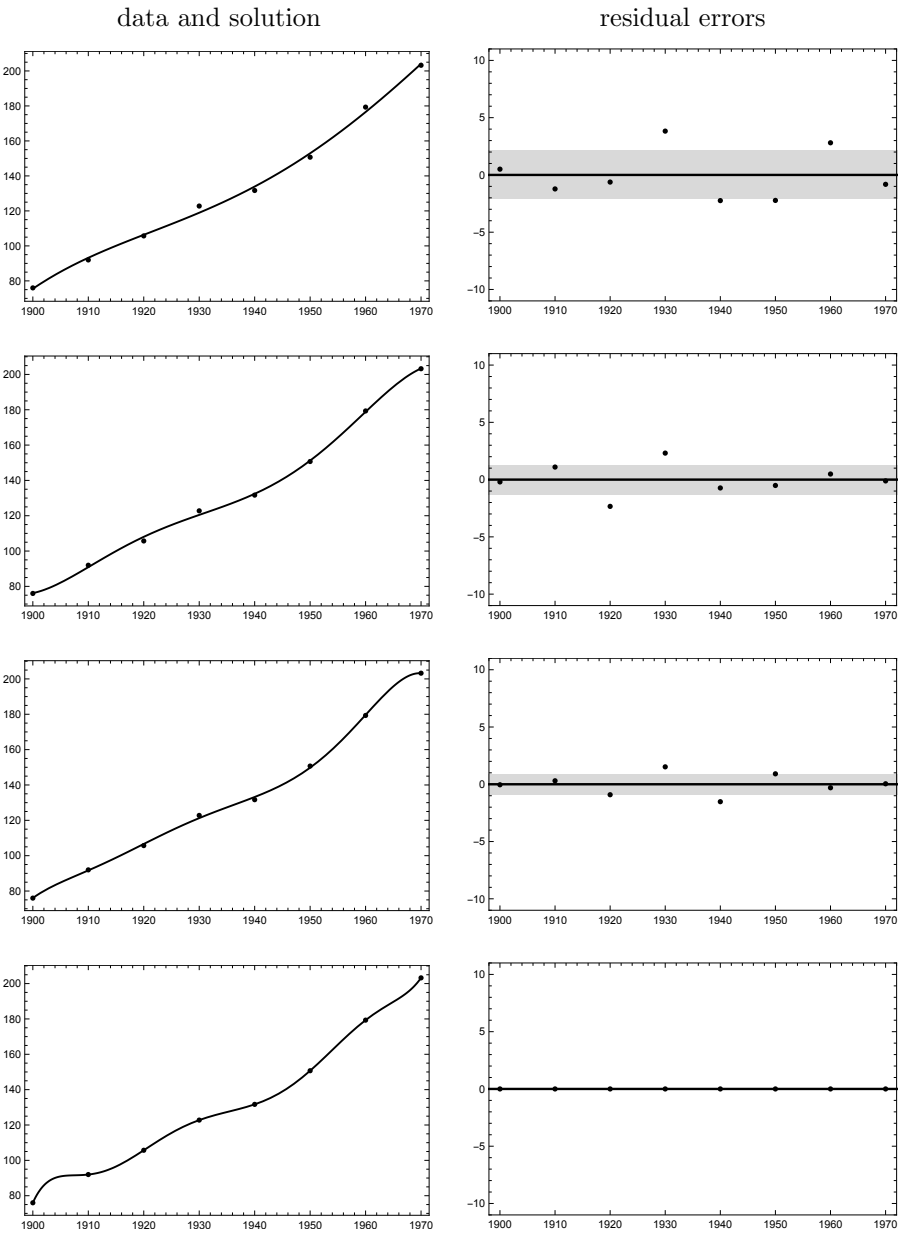
Table 14.4. *Fitting the census data with low order polynomials.*

Table 14.5. *Fitting the census data with higher order polynomials.*



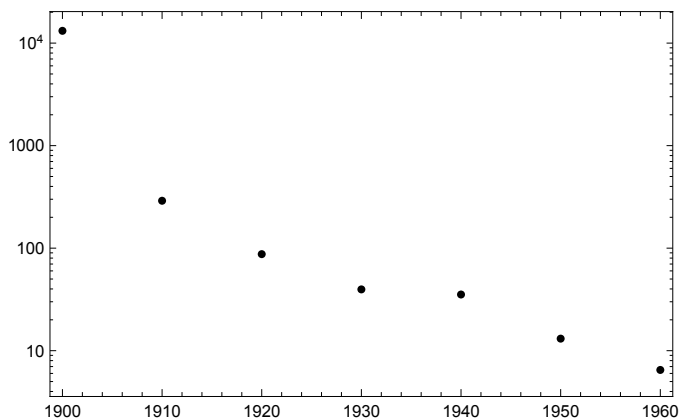


Figure 14.7. Total error $r*r$ by order of fit.

Table 14.6. Projections, by order of fit, for population in 2010.

order	population (millions)
0	133
1	264
2	320
3	420
4	300
5	-928
6	-4,540
7	22,183

Part V

Appendices

Appendix A

Exercises

A.1 Linear systems

§3.2.2 Equation (3.3) describes an extremum. How do we know this is a minimum, not a maximum? Because least squares solutions form a convex set. Proves this.

A.1.1 6.1.1

$$\begin{aligned} a_1 &= \Delta^{-1} \left((\mathbf{1}^T \mathbf{1}) (x^T T) - (\mathbf{1}^T x) (\mathbf{1}^T T) \right) \\ a_1 &= \Delta^{-1} \left((\mathbf{1}^T \mathbf{1}) (x^T T) - (\mathbf{1}^T x) (\mathbf{1}^T T) \right), \\ &\rightarrow \Delta^{-1} \left((\mathbf{1}^T \mathbf{1}) (x^T T + x_* (\mathbf{1}^T T)) - (\mathbf{1}^T x + x_* (\mathbf{1}^T \mathbf{1})) (\mathbf{1}^T T) \right), \\ &= a_1 + \Delta^{-1} \left(x_* (\mathbf{1}^T \mathbf{1}) (\mathbf{1}^T T) - x_* (\mathbf{1}^T \mathbf{1}) (\mathbf{1}^T T) \right), \\ &= a_1 \end{aligned}$$

§6.1.1 Show that under the affine transformation $T \rightarrow \alpha T + \beta$, where $\alpha, \beta \in \mathbb{R}$, the slope transforms as

$$a_1 \rightarrow \alpha a_1.$$

By inspection, argue that a_0 is invariant.

Part VI

Backmatter

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