Human Lifespan is Well-Modeled by Linear Regression

Sergio E. Garcia Tapia

March 28, 2025 (last updated April 2, 2025)

1 Introduction

It has been suggested that the maximum human lifespan is fixed at around 122 years, and is subject to natural constraints [Bla21]. In contrast, the average life expectancy of a newborn increase from 32 years in 1900 to 71 years in 2021; on average, humans are living longer [Dat+23].

As lifetime expectancy increases, it is natural to ask why we should care. One idea is that humans inherently strive for survival, so we naturally look for ways to live longer. Another answer is that a longer life brings more opportunities to nurturing social relationships such as with family and friends or pursuing career and education goals. To what extent, then, can people influence their longevity to enable them to fulfill these life goals? The Danish Twin Study established that only 20% of how long the average person lives is dictated by our genes, whereas the other 80% is dictated by our lifestyle [BS16]. While on a quest to uncover aspects of lifestyle that increase longevity, Dan Buettner from National Geographic discovered regions of high longevity known as blue zones [BS16]. Blue zones are regions that consist of a large number of centenarians (people aged over 100) whose longevity is attributed to healthy lifestyles.

There are 9 primary lifestyle habits that people in blue zones adhere to which contribute to slow-aging. These factors can be summarized as: being in environments that naturally encourage movement; having a meaningful life purpose; having effective routines to cope with stress; eat until 80% full and eat less later times of the day; eat beans and little pork; drink alcohol moderately; get a sense of belonging by being part of some community; prioritize loved ones, such as family and friends; surround yourself with groups of people that encourage healthy behaviors [BS16]. A study by Harvard's School

of Public Health further strengthens the evidence that a healthy lifestyle can increase longevity. The study considered the influence of 5 factors similar to the one we considered for blue zones, such as maintaining healthy eating patterns, not smoking, getting 3.5 hours of vigorous physical activity, drinking alcohol in moderation, and maintaining a normal weight. The study estimated that women at age 50 who did not adopt any of these 5 healthy habits were estimated to live until 79, whereas those who did were estimated to live to 93.1 years [Hic18].

In this paper, I explore the extent to which specific factors can lengthen a person's expected lifespan.

2 Understand Reality via Models

A phenomenon is a naturally occurring event whose manifestation can be observed. Phenomena range from natural disasters such as earthquakes and hurricanes to population growth and change in stock prices and even to gravity and magnetism. In this paper, the phenomenon of interest is human lifespan. Phenomena can be quite complicated, and their study often requires making simplifying general assumptions to simplify their analysis. The result of this simplification is called a model.

A model is an abstraction, or an approximation, or a simplification, of reality. The goal of a model is to remove elements that affect a phenomenon whose consideration may be infeasible or whose contribution may be relatively insignificant. The adage "an apple a day keeps the doctor away" may be thought of as a model for living a healthy life. Another one is "Sitting is the new smoking", which warns that a sedentary lifestyle is a risk factor for cardiovascular morbidity [BSC16]. A person who intends to lead a healthy life may heed these proverbs by exercising more and eating healthier foods, but may quickly realize a problem. How many apples should they eat? How much sitting is too much sitting? The problem is that models can be ambiguous; though the qualitative description is useful for everyday talk, our inability to quantify the effect of each component makes it difficult to apply. If we are to use models to guide the choices we make, we need something more rigorous.

A mathematical model is a specialization of a model that aims to establish a precise relationship between two or more quantities. To create a mathematical model, it is necessary to identify responses, also known as outputs and inputs, also known as factors, as well as ways to measure them. The Navier-Stokes equation, for example, are a mathematical model believed to aid in explaining and predicting breeze and turbulence. [Ins25]. Even without the availability of exact solutions, the model has been useful in navigation in boats in water and modern jet flight across the sky. It is interesting to note that in spite of the ingenuous principles that Navier-Stokes equation embody, they are still far from perfect. In his book, Silver shared that a mathematician once said "The best model for a cat is a cat" [Sil12]. The mathematician meant that no model is perfect, for missing any detail inevitably leads to some inaccuracy. It is also important to remember that a model, mathematical or not, is not reality. Thus, even though models are useful, we ought to be careful not to let them be the final word on the phenomenon that we are attempting to describe with them.

3 A Model for the Lifespan Phenomenon

Let us consider how we would make a mathematical model for predicting the lifespan of an individual. A phenomenon has a response or output, mathematically denoted as y, whose values belong to a response space, a set denoted \mathcal{Y} . In human lifespan, y is the number of years that a person lives from birth to death, which can be 67.4 years or 100.12 years, for example. This quantity can be measured accurately from a person's birth and death certificates. Thus, the response space $\mathcal{Y} = \mathbb{R}$, the set of all real numbers. To devise an exact mathematical relationship involving lifespan y, we need to identify its causal drivers, z_1, z_2, \ldots, z_t , and an exact functional relationship t so that we may write

$$y = t(z_1, z_2, \dots, z_t).$$

The causal drivers are the true causal input information of the phenomenon's response. Here are some possible causal drivers influencing the lifespan of an individual:

 $z_1 =$ Number of fatal accidents experienced

 z_2 = Presence or absence of genes to combat deadly disease

 $z_3 = \text{Ability to gather healthy foods}$

 z_4 = Average daily rate of sustained mental stress

 z_5 = Level of physical aptitude

 $z_6 = \text{Grit}$ and desire for survival and longevity

One can argue that these inputs may not include all that goes into living a long life, or that some of these are not in fact causal drivers. In general, we do not have the omniscience to know exactly which inputs directly cause the response of a phenomenon to change. However, even if we did, we face the challenge of how to precisely measure these causal drivers. For example, it is impossible to know how many accidents a person will experience throughout their life. Similarly, the mental stress that a person experiences is linked to their life experiences and circumstances, which are always changing in unpredictable ways. Even if we did know all the causal drivers and were able to measure them, we would be hard-pressed to establish an exact functional relationship t relating them to the response y, because there are uncountably-many ways to relate these quantities.

4 Approximating Causal Drivers with Features

To deal with the infeasibility of obtaining true causal information, we introduce quantities known as features. A list of features x_1, x_2, \ldots, x_p also known as predictors or covariates, act as proxies to the unattainable true drivers which are the z_1, z_2, \ldots, z_t . By proxies we mean that these features stand for them when describing the phenomenon. The list of features x_1, x_2, \ldots, x_p is said to belong to a covariate space \mathcal{X} , which consists of all the possible lists involving different values of these features. We can express y through a functional relationship in terms of the features, but we must account for the fact that the features do not carry the same information density as the true causal drivers. That is, whereas we have the exact equation $y = t(z_1, z_2, \ldots, z_t)$, we only an approximate equation $y \approx f(x_1, x_2, \ldots, x_p)$ for some function $f: \mathcal{X} \to \mathcal{Y}$, or

$$y = t(z_1, z_2, \dots, z_t) = f(x_1, x_2, \dots, x_p) + \delta$$

where δ is a positive real number. The quantity δ is known as ignorance error, encapsulating the error incurred by expressing y in terms the features x_1, x_2, \ldots, x_p that are not truly causal. The following is list of features that may affect lifespan:

 $x_1 =$ Number of steps walked each day, an integer

 $x_2 =$ Average daily number of hours of sleep

 x_3 = Daily average number of 12-oz cups equivalent to a wine glass

 x_4 = Daily average number of cigars a person smokes

 x_5 = Weekly average number of hours of pleasurable social interactions

We learned earlier how centenarians in blue zones heed the habits embodied by some of these features (and more), so it is reasonable to believe that they affect lifespan. The advent of smart watches makes measuring x_1 and x_2 , which can be a good measure of a person's level of physical activity and aptitude and a way to help cope with mental stress, respectively. Admittedly x_3 may not be as accurate as possible because not even all wines have the same alcohol content. However, a person can take note of when they purchase a box or a single unit of bottles of cans of some drink, and maintain a diary with a headcount of when and how much they consumed. Similarly, cigars are sold in boxes of multiple units, so a person can track their purchases to know x_4 , how many cigars they have smoked. For x_5 , a person can refer to their extracurricular activities, such as outings for restaurants or dancing, video or board game sessions, or even time spent at home talking with friends and family; a calendar can help track this.

5 Stationarity

A reasonable question is whether the causal drivers, features, and functional relationships between them and the response y changes over time. A phenomenon is said to be stationary when such change does not occur. The significance is that if we create a model for a non-stationary model, it may not be applicable in the future as the relationships between the inputs and outputs change. When that happens, we say that the model drifts. In general, most phenomena are not stationary, and we might only be able to achieve stationarity if we have a closed system. For example, though lifespan is thought to have a maximum value, humans might make a scientific discovery to extend it, and in turn it may render our current model obsolete. Human lifespan is not a stationary phenomenon, so we need to acknowledge that in developing for it, we will need to update it so that its accuracy does not suffer heavily. For example, if a scientific breakthrough suggests that there is no maximal lifespan or that there is a way to extend it, we should consider adding features that take this development into account.

6 Supervised Learning

So far, we can express the human lifespan response as

$$y = f(x_1, x_2, x_3, x_4, x_5) + \delta$$

where we rely on p=5 features. We must humble ourselves once more and acknowledge that it is practically impossible to know the exact functional form f. Just like for t, there are infinitely-many ways to relate the features to the response. Moreover, there may not be analytical or exact solution. An example of this was alluded to earlier when we mentioned the Navier-Stokes equation, for which no solution is known [Ins25]. The theory of differential equations is ambitious in its attempt to yield exact or analytical solutions, but we must accept that most problems do not admit such solutions, and the human lifespan phenomenon is no exception. Instead of giving up, we can adjust our expectations and settle for an empirical solution. An empirical solution is an approximate solution that can be obtained from historical data, and the framework for obtaining a solution in this way is called supervised learning.

6.1 The Historical Data

Historical data refers to a set \mathbb{D} of n data points that are measured by observing the phenomenon. We can organize the data points as ordered pairs and enumerate them

$$\mathbb{D} = \{(\boldsymbol{x}_1, y_1), (\boldsymbol{x}_2, y_2), \dots, (\boldsymbol{x}_n, y_n)\}$$

where y_k is the output for the kth observation, expressed as $y = f(x_k)$. For us, $x_k = \begin{bmatrix} x_{k,1} & x_{k,2} & x_{k,3} & x_{k,4} & x_{k,5} \end{bmatrix}^{\top}$, which identifies a person whose health habits are known, and who is known to have lived for y_k years from birth to death. In the framework of supervised learning, a large number n of data points is necessary to ensure the model can have a fair shot at capturing different patterns and thereby accurately make predictions. If our intent is to predict human lifespan, then our n must be reasonably large given that the world has about 8 billion people [Cen24]. An important consideration is how easily such data can be obtained.

In measuring the human lifespan phenomenon, collecting each data point literally takes a lifetime; an observation is not complete until someone has died. We could gather data on the features of live people through a survey and collect their features as a vector $x \in \mathcal{X}$, then when they have passed in the future we associate that vector with a lifespan $y \in \mathcal{Y}$. If we wanted to have data now, it might be expensive or impossible because of data protection laws. One crude way to gather some data about lifespans y would be to use Wikipedia, which often lists the birth and death dates of individuals (with accuracy subject to verification), but we may not find the information on the specific features x that we are interested for that individual. After all, not everyone tracks how many alcoholic drinks they intake every day, or exactly how many hours of sleep they get at night. Moreover, technology such as smart watches to track a person's daily footstep count was either not invented or widely unavailable years ago. Another consideration is that we may need to prefer more recent deaths to account for the fact that human lifespan is a non-stationary phenomenon.

6.2 Hypothesis Set

The next part of the supervised learning framework deals with the infeasibility of an analytical solution. Simply put, f is and will always be unknown, and it is likely much more complex than we can handle. Since we seek an approximate solution, or next task is to conjecture a functional form that we believe may well approximate f. The notation \mathcal{H} denotes a set of candidate functions, where $h \in \mathcal{H}$ means that $h : \mathcal{X} \to \mathcal{Y}$, meaning h is a function with the same domain and range as f. Our hope is to find the best function $h^* \in \mathcal{H}$ that approximates f. Such a function would be related to g, g, and g by

$$y = t(z_1, z_2, \dots, z_t) = f(x_1, x_2, \dots, x_n) + \delta = h^*(x_1, x_2, \dots, x_n) + \epsilon$$

Here, ϵ is known as noise, which is accrual of the ignorance error and a quantity known as misspecification error, expressed as $\epsilon - \delta = t - f$. To understand where it comes about, let us consider an example. Suppose we define the following set of candidate functions:

$$\mathcal{H} = \{ w_0 + w_1 x_1 + w_2 + w_3 x_3 + w_0 x_4 + w_5 x_5 \mid w_0, w_1, w_2, w_3, w_4, w_5 \in \mathbb{R} \}$$

Set \mathcal{H}_1 is the set of hyperplanes, which is a generalization of the idea of lines in higher dimensions. Such a choice is attractive because of the ease of interpretation of the results, and is appropriate if the data appears to have a "linear" pattern. On the other hand, suppose our colleague proposes the following candidate set:

$$\mathcal{H}_2 = \{ w_0 + w_1 x_1 x_2 x_5 + w_2 x_3 x_4 \mid w_0, w_1, w_2 \in \mathbb{R} \}$$

Our colleague reasons that people who smoke are also most likely to drink, so the multiplicative term x_3x_4 , which is known as a first-order interaction of features, accounts for this. Similarly, they argue walking, sleeping, and interacting with friends compound on each other some way, leading to other lifespan gains. Whether or not that is accurate, this example demonstrates that there are many ways to design a candidate set. Not all candidate sets may be equally effective in approximating f. For example if the data has a parabolic pattern, fitting a line may not be a good idea. The misspecification error $\epsilon - \delta$ quantifies this "mismatch" between the candidate set and the data set $\mathbb D$ that presumably came from f. I think our colleague may be right in considering a hypothesis set with interactions. For example, studies shows that smokers tend to sleep less than non-smokers [Jae+12]. Thus, interactions between the features may improve our model. However, they can also make our model more complex to reason about, so going forward we will use \mathcal{H} as our set of candidate functions.

6.3 Algorithm

Once we have measured data \mathbb{D} in the real world and chosen a candidate set of functions \mathcal{H} , we need a way to obtain a function h from \mathcal{H} . The set \mathcal{H} is infinite and we need a way to search the space efficiently to obtain an approximation to f. To do this, employ an algorithm, denoted \mathcal{A} , which is a finite sequence of steps devised to solve a problem. We can view an algorithm as a function $\mathcal{A} = \mathcal{A}(\mathbb{D}, \mathcal{H})$ that accepts a data set and a set of candidate functions and outputs an approximation $h \in \mathcal{H}$ to f. Earlier we said that h^* is the best approximations to f in \mathcal{H} , so naively we might hope for an algorithm h^* . Once again, though, we must temper our expectations; as we have emphasized, \mathcal{H} is huge, uncountably-infinite as a matter of fact. Another way to think about it is that many functions do not attain all possible outputs in their co-domain, and the algorithm \mathcal{A} that we use is no different. We will say that the best possible function that our algorithm can produce is $g \in \mathcal{H}$, related to g, g, g, and g by

$$y = t(z_1, z_2, \dots, z_t) = f(x_1, x_2, \dots, x_p) + \delta = h^*(x_1, x_2, \dots, x_p) + \epsilon = g(x_1, x_2, \dots, x_p) + \epsilon$$

The symbol e denotes a new type of error called a residual, and introducing g also introduces a new type of error called estimation error, given by the difference $h^*-g=e-\epsilon$. In a future section, we will discuss errors in more detail.

Many ingenuous algorithms have been developed by very smart people, each subject to certain constraints and weighing different parts of the supervised learning problem differently. For example, if the response space in our phenomenon was binary, such as it would be if we were modeling the presence or absence of a disease, then we might encode that as the set $\mathcal{Y} = \{0,1\}$. One algorithm, known as perceptron, is applicable if the data set \mathbb{D} satisfies a condition known as linear separability, which most data sets do not satisfy. Another algorithm, known as logistic regression, produces a candidate function g whose output is not in \mathcal{Y} , but rather it outputs a probability that a person has the disease. In the case of human lifespan with a numeric response space $\mathcal{Y} = \mathbb{R}$, some plausible choices of algorithms are K-nearest neighbors (KNN) and ordinary least squares regression (OLS). In this paper, I will choose OLS.

7 Predictions

Before considering OLS, let us re-center ourselves by remembering why we are using supervised learning in the first place. We want to understand our phenomenon, or at least, we want to make predictions about the response of the phenomenon given a unit and their features. Suppose that Luis is 24-year old individual who works part-time with an interest in planning for retirement and in pursuing a college education. He determines that knowing how long he will be alive for will help in his planning. We offer Luis help, but we tell him that we need some information about him. After some initial hesitation to share personal information, we gather the following vector:

$$\boldsymbol{x}_{\mathrm{Luis}} = \begin{bmatrix} 12042 & 7.3 & 0.1 & 0 & 15.7 \end{bmatrix}^{\mathrm{T}}$$

That is, Luis walks on average 12,042 steps each day, sleep on average 7.3 hours each night, drinks 0.1 glasses of wine each day on average, does not smoke at all, and spends about 15.7 hours a week on average socialization with friends and family. Having applied our algorithm and obtained a function $g = \mathcal{A}(\mathbb{D}, \mathcal{H})$, we can use g to make a prediction. For example, suppose that our algorithm produced the vector

$$\boldsymbol{b} = \begin{bmatrix} 10 & \frac{1}{1000} & \frac{20}{8} & 5 & -10 & 3 \end{bmatrix}$$

Then our function g is given by

$$\hat{y} = g(\mathbf{x}) = 10 + \frac{1}{1000}x_1 + \frac{20}{8}x_25x_3 - 10x_4 + 3x_5$$

The quantity \hat{y} is the predicted number of years for a person whose feature as given by the feature vector \boldsymbol{x} . For Luis, we can compute this number explicitly using our made-up model g:

$$g(\mathbf{x}_{\text{Luis}}) = 87.892$$

That is, our model predicts that Luis will live for almost 88 years. Luis seems a bit disheartened, sharing that he wanted to live longer, but he is also curious as to whether our prediction is accurate. We tell Luis that our model is still experimental, so he should not let it have the final word. We humbly also tell him that there may be better models out there. How can we ensure that our model is as good as can be? We need an error metric.

8 Types of Error

Before discussing error metrics, we need to understand discuss how errors are introduced in our modeling framework. The first type of error, ignorance erro(r, arises from using features x_1, x_2, \ldots, x_p as proxies for the true drivers (casual inputs) z_1, z_2, \ldots, z_t and a function f domain contains the possible values of these features instead of the true phenomenon function f:

$$y = t(z_1, z_2, \dots, z_t) = f(x_1, x_2, \dots, x_n) + \delta$$

The error due to ignorance accounts for the fact that the features do not contain the same information density as the true drivers. One way to deal ignorance is to introduce new relevant features to describe the phenomenon, thus increasing the number of features p per observation. The next approximation we made was to introduce h^* from some candidate set \mathcal{H} :

$$f(x_1, x_2, \dots, x_p) + \delta = h^*(x_1, x_2, \dots, x_p) + \epsilon$$

The quantity ϵ is known as the noise, and the difference $\epsilon - \delta = f - h^*$ is the misspecification error. We touched on this earlier, and it is a result of choosing a candidate set \mathcal{H} that may not correctly capture the functional form of f. For example, if the points in \mathbb{D} , presumably produced by f (subject to some error due to δ) seem to follow a parabolic pattern, then a set of hyperplanes may accrue some misspecification error. One way to

reduce misspecification is to choose a larger set of candidate functions. The last type of error is known as the residual, which shows up when we obtain a modeling function g from our algorithm:

$$h^*(x_1, x_2, \dots, x_p) + \epsilon = g(x_1, x_2, \dots, x_p) + \epsilon$$

The residual contains the ignorance and misspecification error, but it adds a new type of error known as estimation error, given by $e - \epsilon = h^* - g$. The estimation error results from not being able to obtain the best possible function h^* out of \mathcal{H} . One way to reduce estimation error is to employ a better algorithm \mathcal{A} . Another way is to ensure that the features that we use are indeed relevant. For example, suppose that a colleague recommends that we use a new feature x_6 that represents the number of digits of π that you know. You find it hard to believe that this affects lifespan at all. Suppose you are omniscient and know that it indeed has nothing to do with lifespan. If your model accounts for it and produces a non-negative weight associated with x_6 , then we accrue estimation error. Estimation error can also increase if we decide to transform our features, perhaps by introducing a new feature that is the square of an existing feature. For example, if we had a single feature x_1 and believed the data set had a quadratic pattern relative to this feature, the we would create a transformed feature x_1^2 to allow for parabolic solutions out of our candidate set. This may lead to a decrease in misspecification error, but possibly an increase in estimation error.

We have the most control over the residual error e. Next, we discuss error metrics, and regularly consider residual error in our calculations.

9 Error Metrics

10 The Null Model

Before we consider OLS, it is worth thinking about what, if anything, we can achieve without one. In other words, suppose we go on Wikipedia and learned the lifespan y_1, y_2, \ldots, y_n of n different people. However, their Wikipedia page did not have information about the number of steps they walked each day, the number of hours they slept each night, or any of the features x_1, x_2, x_3, x_4, x_5 that we are using as proxies to the true drivers. In the absence of feature, there is still one model we can employ: the null model.

Denoted g_0 , the null model also comes from the hypothesis set \mathcal{H} , so its form depends on the functional forms allowed by \mathcal{H} . Thus, we know at the very least that g_0 gets its input from \mathcal{X} and outputs a value in \mathcal{Y} . Since $\mathcal{Y} = \mathbb{R}$ is numeric, the null model corresponds to computing the arithmetic mean:

$$g_0(\boldsymbol{x}) = \frac{1}{n} \sum_{i=1}^n y_i$$

That is, to make a prediction about the lifespan of an individual whose features are x, we actually ignore the features, and predict that their lifespan is simply the average of the lifespan of the n individuals we have seen. In some way, this is justified if we truly know nothing about each person other than how long they lived. The null model is useful because it can act as a reference for performance. It stands to reason that if we have accurate features about the health habits of an individual, if those features are useful at all as predictors for lifespan, and if our algorithm is any good, the our computed model g ought to do better than g_0 , which literally knows nothing about the person. In a future session, we will discuss errors and how to determine whether our model performs better than g_0 .

11 Ordinary Least Squares Regression

To understand ordinary least squares, suppose that the number of features in our model is p = 1. For concreteness, say that we are only tracking x_1 , the number of steps a person walks each day, hereby referred to as x for this example. Then our data set \mathbb{D} is a list of points in the two-dimensional Cartesian coordinate plane:

$$\mathbb{D} = \{(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)\}\$$

Note that x_1 in this example is the number of steps the first person walks each day, not the generic name for the walk feature, and x_2 is the number of steps the second person sleeps, not the generic name for the daily sleep feature. We can visualize \mathbb{D} using a scatterplot, as in Figure 11. In the past, you may have learned about the idea of the line of best fit. A general line has the equation $\hat{y} = w_0 + w_1 x$ for some parameters $w_0, w_1 \in \mathbb{R}$. For the value of the feature x_k corresponding to the number of steps the kth person takes daily, the line would predict $\hat{y}_k = w_0 + w_1 x_k$. The actual response for the kth person is y_k , so the error in the prediction with this line is $e_k = y_k - \hat{y}_k$. Since we have n points, that implies n residual errors e_1, e_2, \ldots, e_n . The idea of ordinary least squares is to pick a pair of parameters $b_0, b_1 \in \mathbb{R}$ defining a line $\hat{y} = g(x) = b_0 + b_1 x$ so that the sum of the squares of the errors, denoted SSE and given by

$$SSE := \sum_{k=1}^{n} e_k^2 = \sum_{k=1}^{n} (y_k - \hat{y}_k)^2$$

is as small as possible. Implicit is our discussion is the assumption that out algorithm \mathcal{A} is searching through a space \mathcal{H} of all possible lines, which is indeed the case since earlier defined \mathcal{H} to be the set of hyperplanes. The algorithm, then can be expressed as

$$\boldsymbol{b} = \operatorname*{argmin}_{\boldsymbol{w} \in \mathbb{R}^{p+1}} \{ SSE \}$$

where $\boldsymbol{b} = \begin{bmatrix} b_0 & b_1 & b_2 & \cdots & b_p \end{bmatrix}^{\top}$ is a vector of all the least square parameters. In our example, p = 1, so we would have $\boldsymbol{b} = \begin{bmatrix} b_0 & b_1 \end{bmatrix}^{\top}$. Note the procedure for the algorithm can be read as "find the parameters that minimized the SSE".

Admittedly, ordinary least squares may not be a suitable algorithm to use if our data set \mathbb{D} does not suggest a linear trend. However, what is attractive is that under mild conditions it admits an exact solution $g = \mathcal{A}(\mathbb{D}, \mathcal{H})$. Indeed, throughout our discussion, we have had to give up hope on finding an exact solution for h^* , f, and t, so why would g be any different? It turns out that many minimization problems like the one for SSE do not have exact solutions and required complicated optimization algorithms to give an approximate solution. Thus, when it comes to being able to state a precise solution, OLS is the exception, not the rule.

That is not to say that g, the function determine by the OLS procedure, will be effective for making predictions. We know, for example, that there are some optimal weights $\beta_0, \beta_1 \in \mathbb{R}$ defining a linear function $h^*(x) = \beta_0 + \beta_1 x$ that our algorithm will not produce. Since our g is different from h^* , we accrue estimation error.

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