
Integrating Domain Knowledge With Data: From Crisp To Probabilistic and Fuzzy Knowledge

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Abstract

It is well known that prior knowledge about the domain can improve (often drastically) the accuracy of the estimates of the physical quantities in comparison with the estimates which are solely based on the measurement results. In this paper, we show how a known method of integrating crisp domain knowledge with data can be (naturally) extended to the case when the domain knowledge is described in statistical or fuzzy terms.

1 INTEGRATING DOMAIN KNOWLEDGE WITH DATA: FORMULATION OF THE PROBLEM

A large part of information about the world comes from measurements. However, for many complex systems, some characteristics are very difficult to measure: e.g., for a jet engine, it is difficult to measure the temperature and pressure inside the jet chamber, where the temperatures are very high; for a human body, it is difficult to measure the characteristics of the internal organs, etc. In many such cases, experts have some knowledge about the domain. It is therefore desirable to use this knowledge to improve the measurement results.

It is known that prior knowledge about the domain can indeed improve (often drastically) the accuracy of the estimates of the physical quantities in comparison with the estimates which are solely based on the measurement results; see, e.g., (Reznik et al. 1985), (Kreinovich et al. 1986), (Kreinovich et al. 1990), (Luo et al. 1989), (Kreinovich Chang 1992), (Kreinovich Quintana 1992).

In this paper, we show how a known method of integrating data with crisp domain knowledge can be (nat-

urally) extended to the case when the domain knowledge is described in statistical or fuzzy terms.

2 INTEGRATING CRISP DOMAIN KNOWLEDGE WITH DATA

2.1 GENERAL IDEA AND ILLUSTRATING EXAMPLE

As we have already mentioned, if we know the *a priori* relation between the measured quantities x_1, \dots, x_n , then we can often use this relation to correct the measurement results $\tilde{x}_1, \dots, \tilde{x}_n$ and thus, to increase the accuracy of our measurements.

Let us show, on a simple example, that such an improvement is indeed possible. Let us assume that we are measuring two quantities x_1 and x_2 , that both quantities are measured by a measuring instrument with 0 average and standard deviation σ , and that the results of these measurements are \tilde{x}_1 and \tilde{x}_2 .

If we do not have any *a priori* information about the relationship between x_1 and x_2 , then the only thing that we know about, e.g., x_1 is that x_1 is approximately equal to \tilde{x}_1 with a standard deviation σ .

If we know that, in reality, these two quantities should have the same value, i.e., that $x_1 = x_2$, then, for each of these quantities, we have, in effect, not one but *two* independent measurements with the same accuracy σ , resulting in the values \tilde{x}_1 and \tilde{x}_2 . It is well known that in this situation, we can get a better estimate for x_1 if we take an arithmetic average of these two measurements: $\tilde{x} = (\tilde{x}_1 + \tilde{x}_2)/2$; this new estimate has a standard deviation $\sigma/\sqrt{2}$.

2.2 GENERAL DESCRIPTION

Most traditional statistical methods for processing measurement results are organized as follows: Crudely speaking, for each possible combination of actual val-

ues x_1, \dots, x_n , we compute the conditional probability $L(x_1, \dots, x_n)$ of the observed measurement results under the assumption that the actual values are x_1, \dots, x_n . Then, as an estimate for x_i , we choose the values for which this conditional probability is the largest possible.

The conditional probability $L(x_1, \dots, x_n)$ is called a *likelihood function*, and the corresponding method of finding the estimates for x_i is therefore called the *Maximum Likelihood Method*.

An *a priori* constraint on the values x_1, \dots, x_n means that not all combinations (x_1, \dots, x_n) are possible. Instead, we know the set $S \subset R^n$ of possible combinations.

In this case, it is reasonable, instead of looking for the values x_i which provide a *global* maximum for the likelihood function, to return the vector $x = (x_1, \dots, x_n) \in S$ for which the value $L(x)$ is the largest possible among all vectors $x \in S$.

In the typical case when measurement errors $\Delta x_i = \tilde{x}_i - x_i$ are independent and normally distributed, with 0 average and standard deviations σ_i , the likelihood function $L(x)$ is the product of probability densities of several Gaussian distributions:

$$L(x) = \text{const} \cdot \exp \left(-\frac{(\tilde{x}_1 - x_1)^2}{2\sigma_1^2} \cdot \dots \cdot \frac{(\tilde{x}_n - x_n)^2}{2\sigma_n^2} \right).$$

This function $L(x)$ is complex, so maximizing it is difficult; it is known, however, that we can simplify the optimization problem if, instead of maximizing the likelihood function itself, we minimize its negative logarithm

$$-\log(L(x)) = \text{const} + \frac{(\tilde{x}_1 - x_1)^2}{2\sigma_1^2} + \dots + \frac{(\tilde{x}_n - x_n)^2}{2\sigma_n^2}.$$

In particular, if all the measurements are equally accurate, i.e., if $\sigma_1 = \dots = \sigma_n$, then minimization of $-\log(L(x))$ is equivalent to minimizing the distance between the points $x = (x_1, \dots, x_n)$ and $\tilde{x} = (\tilde{x}_1, \dots, \tilde{x}_n)$. Thus, when we know the measurement results \tilde{x} and the set S of all the vectors x which are consistent with the expert's knowledge, then the optimal estimate is the point $x \in S$ which is the closest to \tilde{x} .

The above toy example is in complete agreement with this geometric interpretation: namely, in this example, $S \subset R^2$ is the diagonal $S = \{(x_1, x_1)\}$, and for any point $\tilde{x} = (\tilde{x}_1, \tilde{x}_2)$, the closest diagonal point is indeed the point (x_1, x_1) with $x_1 = (\tilde{x}_1 + \tilde{x}_2) / 2$.

2.3 CASE STUDY: TELEMANIPULATORS

2.3.1 Telemanipulators Are Needed

In many real-life mechanical tasks, it is difficult or even impossible to use humans:

- Some environments are too dangerous for a human being: for example, when we manipulate objects in space, inside a radioactive part of a nuclear reactor, in a dangerous chemical environment, or even in a potentially dangerous environment such as handling viruses that cause deadly diseases.
- In other environments, there is no danger to the human operator, but there is a significant risk of contamination of the object: e.g., in handling microchips, lunar samples, etc.

In all these cases, reasonably simple mechanical tasks can be done by an automatic mechanical hand-arm. However, there is a limit on the complexity of the tasks that automatic devices can do. For more complicated tasks, for which we cannot use a completely automated system, we must use *telemanipulators*, i.e., devices in which a mechanical hand-arm copies the movements of a human operator.

2.3.2 Telemanipulators: Successes

The main goal of the telemanipulator is to reproduce the operator's movements as accurately as possible.

A human hand is a very flexible instrument. In mechanical terms, we can say that it has many degrees of freedom: we can move and rotate the hand itself, the arm, each finger, parts of each finger, etc. Thus, to reproduce its movements accurately, the manipulator also has to have many degrees of freedom.

At present, the best of widely available hand-arm manipulators, the Utah/MIT hand, has 22 degrees of freedom. It is still slightly less than a human hand, because, e.g., it only has 4 fingers and not 5. However, it can perform many important tasks that a human hand can do.

This hand was not designed for telemanipulation only. It has many other applications: e.g., it can even twist itself into the positions that would have been impossible for a human hand.

2.3.3 Telemanipulators: Problems

Both in the Hollywood movies and in the self-made movies that researchers show at robotic conferences, telemanipulation works perfectly well: a robotic hand

exactly reproduces the operator's movements. This is indeed happening in many application areas, but this reproduction accuracy is extremely difficult to achieve.

If we simply measure the pressure, etc., applied by the operator's arm, and send exactly proportional control signal to the electric motors that control different degrees of freedom of the robotic arm, we get a behavior that is often drastically different from what the operator did. For example, the operator's firm grip on the object may be distorted into the robotic arm dropping it, and vice versa, the operator's tender approach to a fragile object may result in a robotic arm's bumping into the actual object and damaging it.

There are three main reasons for the difference between the movements of the human and robotic hands:

- first, the sensors that measure the human hand's pressure are not 100% accurate;
- second, the motors and actuators are not perfect, and do not react precisely to the commands;
- third, the mechanical characteristics of the robotic hand itself are somewhat different from the mechanical characteristics of the operator's hand.

2.3.4 As Manipulators Get More Complicated, These Problems Get More And More Important

The above inaccuracy problems can be traced even on the example of simple manipulators that have a few degrees of freedom, but for more advanced manipulators, these problems become more and more acute. Indeed, for a manipulator, more advanced means that this manipulator has more degrees of freedom. Each degree of freedom bring its own inaccuracy, so if we have 22 degrees of freedom, then in principle, we get 22 sources of inaccuracy all leading to the huge inaccuracy of the resulting action.

Let us give a simple example.

- If we have a 3-finger manipulator, then for this manipulator to grip an object, it must place one finger below it, and two fingers above it. Due to inaccuracy, we may have a slightly distorted position, but we will still keep firmly 3 points on the object.
- For a 4-finger arm that is similar to the human arm, we need to place 3 fingers on top of the object. If, e.g., the upper surface is planar, we must have all 3 fingers on one line. Due to inaccuracy, one of these fingers may be higher than the others. As a result, this finger may not contact the

object at all, and hence, the grip will not be as firm as we desired.

So inaccuracy is harmful. In order to figure out how to decrease this inaccuracy, let us first analyze how we can describe it in precise terms.

2.3.5 How We Can Correct Manipulator Inaccuracy: The Idea of Virtual Tools

Although a human hand-arm has many degrees of freedom, we rarely use all of them in the same movement. Usually, the movement in different degrees of freedom is very much *coordinated*.

For example, if we have already firmly grasped an object, then we move the arm as a whole and, unless necessary, do not use the ability to move fingers and/or or fingertips separately.

There is a limited number of typical movements of this type, and a teleoperator can pretty well describe which of these typical movements he is applying at any given moment of time. When we get this information, we can use it to set up a similar coordination between the degrees of freedom of the robotic hand-arm. When the resulting constraints are in place, the originally flexible robotic hand acts as a new tool that is specifically designed for this type of movement. Since in reality, we are still using the same robotic hand, this is not a new *physical* tool, but a *virtual* tool.

We will see that using virtual tools can indeed be very helpful.

2.3.6 How to Describe Movement Type in Precise Terms

In precise terms, a fixed movement type means that we cannot have *arbitrary* values of x_1, \dots, x_n : these values must satisfy one or several restrictions (constraints).

For example, if we want the arm to move as a whole, then one of these constraints may take the form $x_1 = x_2$ (or $x_1 - x_2 = 0$), where x_1 and x_2 are pressures applied by two fingers. If we want to preserve the distance between the two fingertips, then we may require something like $(x_1 - x_2)^2 + (x_3 - x_4)^2 - \text{const} = 0$.

In general, we have one or several constraints of the type $F(x_1, \dots, x_n) = 0$. So, for the actual values $x_i = \tilde{x}_i - \Delta x_i$, in addition to the measurement results \tilde{x}_i , we have additional constraints of the type

$$F(x_1, \dots, x_n) = F(\tilde{x}_1 - \Delta x_1, \dots, \tilde{x}_n - \Delta x_n) = 0.$$

Since inaccuracies Δx_i are small, we can expand the function F in Taylor series and ignore terms that are quadratic or of higher order in Δx_i . As a result, each

original constraint $F_k(x_1, \dots, x_n) = 0$ becomes a *linear* constraint on possible values of Δx_i :

$$F_{k1} \cdot \Delta x_1 + \dots + F_{kn} \cdot \Delta x_n - F_k = 0, \quad (1)$$

where

$$F_{ki} = \frac{\partial F_k(x_1, \dots, x_n)}{\partial x_i} \Big|_{x_1=\tilde{x}_1, \dots, x_n=\tilde{x}_n}$$

and $F_k = F_k(\tilde{x}_1, \dots, \tilde{x}_n)$.

Thus, instead of the original vector \tilde{x} , we use the vector x which is the closest to \tilde{x} among the hypersurface S of all the vectors which satisfy the conditions (1), i.e., the vector x which is a projection of \tilde{x} on S . Computing this projection is a straightforward geometric task.

2.3.7 A Simple Example Showing That Constraints Can Decrease Uncertainty

Suppose that we have a movement in which two fingers have to move in the exact same way, i.e., in which $x_1 = x_2$. Let us also suppose that the measurement accuracy is $\sigma_1 = \sigma_2 = 0.2$. Then, if we take the constraint into consideration, we can decrease the standard deviation $\sqrt{2}$ times.

2.3.8 Implementation

The papers (Nelson et al. 1995), (Fuentes et al. 1996), (Jagersand et al. 1997), (Kreinovich et al. 1998) describe the actual implementation of several virtual tools for grasping and manipulation with the Utah/MIT hand. This implementation does lead to an improved telemanipulation performance.

3 INTEGRATING STATISTICAL DOMAIN KNOWLEDGE WITH DATA

Often, instead of crisp constraints of the type $f(x_1, \dots, x_n) = 0$, in which we know the exact value (e.g., 0) of some function $y = f(x_1, \dots, x_n)$ of the measured quantities, we have *statistical* knowledge, which is described, e.g., in terms of a known probability distribution for the quantity $y = f(x_1, \dots, x_n)$.

In this case, we can still use the Maximum Likelihood Method as for crisp constraints; the only difference is that in addition to the measurement errors, we also need to take into consideration the probability of the corresponding value y . Namely, we have $L(x) = \rho_m(\tilde{x} - x) \cdot \rho_c(f(x_1, \dots, x_n))$, where $\rho_m(\Delta x)$ is the probability density function which describes measurement errors, and $\rho_c(y)$ is the probability density which describes the statistical domain knowledge.

If both distributions are Gaussian, then, instead of maximizing a complex function $L(x)$, we can minimize a much simpler (quadratic) function $-\log(L(x))$, and get a simple system of linear equations from which we can easily find the desired estimate of x .

4 TOWARDS INTEGRATING FUZZY DOMAIN KNOWLEDGE WITH DATA

4.1 THE NEED FOR FUZZY DOMAIN KNOWLEDGE

Often, experts can only formulate their knowledge in terms of words from natural language. For example, an expert can say that “the temperature inside the chamber is about 10,000 degrees” (or “huge”), where the word “about” does not have a precise meaning.

A natural way to formalize such a knowledge is to use *fuzzy logic*, in which each such term like “small” is described by a *membership function* $\mu_{\text{small}}(x)$ which tells, for every value x , the degree $\mu_{\text{small}}(x) \in [0, 1]$ to which this value is small. After this formalization, we get a *fuzzy knowledge domain*.

4.2 HOW CAN WE COMBINE THIS KNOWLEDGE WITH DATA?

A natural question is: *How can we combine fuzzy domain knowledge with data?*

There exist many heuristic methods of combining such data, but since we want to *guarantee* that this combination indeed improves the measurement results (and, ideally, improves them in the best possible way), we would like to have *guaranteed* methods of combining fuzzy domain knowledge with data.

4.3 OUR MAIN IDEA: LET US TREAT AN EXPERT AS YET ANOTHER MEASURING INSTRUMENT

From the viewpoint of getting new information about the world, we can consider an expert as one more measuring instrument, an instrument which brings us additional information about the actual (unknown) values of different physical quantities. From this viewpoint, the only difference between an expert and a more traditional measuring instrument is that a measuring instrument returns a *number* while an expert returns a *word* (“small”, “large”, etc.). Within this approach, we can determine the statistical characteristics of an expert along the same lines as we determine the statistical characteristics of the measuring instrument.

Namely, to determine the statistical characteristic of a measuring instrument, we use it to measure several quantities for which the precise value has already been measured by a (more) precise (standard) measuring instrument. Then, for each possible value x of the measured quantity, we group together all the measurements in which the actual value was x (or close to x). For all measurement results \tilde{x} from the group, we compute the measurement error $\Delta x = \tilde{x} - x$, and then we use the standard statistical analysis to determine the (conditional) probability distribution of this error $P(\Delta x|x)$.

Similarly, we take several situations in which the desired quantity has different values and whose value has been measured by an accurate measuring instrument. For each of these situations, we can then ask an expert to provide us with a word which best describes her perception of the corresponding value x . Then, for each possible value x of the measured quantity, we group together all the measurements in which the actual value was x (or close to x). Then, for each of the words w which an expert used, we can compute the (conditional) probability $P(w|x)$ that an expert used the word w under the condition that the actual value was x .

As soon as we have this statistics, we can process the expert's estimates just like we would have processed one more measurement result. In this case, mathematical statistics guarantees that on average, combining expert estimates with the measurement results do not make things worse.

Comments. This idea is not as heretic as it may seem (reducing fuzzy to probabilities?). On one hand, our method of determining a *probability* function $P(w|x)$ corresponding to a word w is actually very similar to the main methods of determining the *membership* functions that describe a word w ; see, e.g., (Dubois et al. 1980), (Norwich et al. 1984), (Turksen 1991). So, we can view the function $P(w|x)$ as a membership function. On the other hand, the very fact that fuzzy set theory can be, in principle, reformulated in more statistical terms (namely, in terms of random sets) is known for more than two decades; see, e.g., (Nguyen 1979).

4.4 PEDAGOGICAL EXAMPLE

Let us consider a simple example of how this combination can be done. Let us assume that an expert estimated the value as “small”, and that the corresponding probability function for “small” is $P(\text{small}|x) = \mu(x)$. Let us also assume that we have made a single measurement by a measuring instrument whose measurement error Δx is normally distributed with 0 av-

erage and known standard deviation σ , and that the measurement result is \tilde{x} . The corresponding probability density will be denoted by $\rho(\tilde{x} - x)$.

We want to use Maximum Likelihood Method and find the value x for which the value of the likelihood function $L(x)$ is the largest possible. For each x , the probability of getting a value \tilde{x} is equal to $\rho(\tilde{x} - x)$, the probability of an expert's estimate's “small” is equal to $\mu(x)$. It is natural to assume that the measuring instrument and an expert are statistically independent. Therefore, the probability $L(x)$ of getting, for a given x , both the measurement result \tilde{x} and the expert estimate “small” is equal to the product $L(x) = \rho(\tilde{x} - x) \cdot \mu(x)$.

To find the maximum of this function $L(x)$, it is usually convenient to seek a minimum of its negative logarithm

$$-\log(L(x)) = \frac{(\tilde{x} - x)^2}{2\sigma^2} - \ln(\mu(x)) \rightarrow \min.$$

Differentiating this expression w.r.t. x and equating the derivative to 0, we conclude that

$$x = \tilde{x} + \sigma^2 \cdot \frac{\mu'(x)}{\mu(x)},$$

where $\mu'(x)$ denotes the derivative of $\mu(x)$. Since $x \approx \tilde{x}$, we have $\mu(x) \approx \mu(\tilde{x})$; thus, can neglect the terms which are of second order w.r.t. Δx , and get an explicit analytical expression

$$x \approx \tilde{x} + \sigma^2 \cdot \frac{\mu'(\tilde{x})}{\mu(\tilde{x})}.$$

The difference between x and \tilde{x} is a correction term which is motivated by the fuzzy domain knowledge and which makes the resulting estimate more accurate (provided, of course, that the experts were correct).

4.5 THIS IDEA HAS BEEN SUCCESSFULLY USED IN PRACTICAL APPLICATIONS

Our foundations are new, but the resulting formulas (which we obtained heuristically) have actually been successfully used in practical applications; see (Kreinovich et al. 1986) and references therein.

The domain to which we applied it was testing jet engines. In this domain, expert estimates are extremely important, because an important part of this testing is trying to figure out what is going on in the high-temperature regions, and the temperatures are so high there that we cannot place any sensors. So the only available information about these regions is the experts estimates.

One of the authors (L.R.) used this fuzzy representation of uncertainty in designing software for the automated jet engines testing system IVK-12. This system was successfully used to test jet engines for aircraft and spaceships.

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