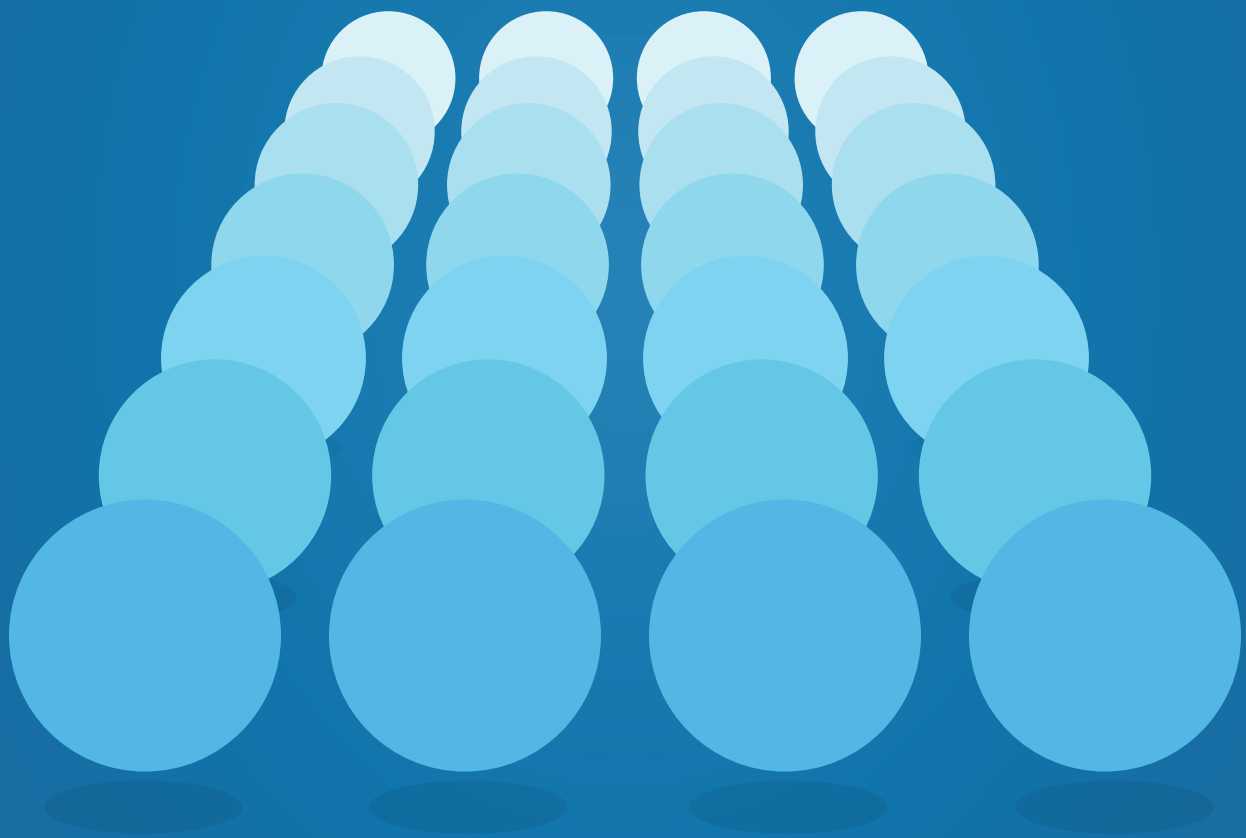


Everything You Need To Know About

Scalability



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VividCortex

Meet the Author

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Baron is a database expert who is well-known for his contributions to the MySQL, PostgreSQL, and Oracle communities. An engineer by training, Baron has spent his career studying how teams build reliable, high performance systems, and has helped build and optimize database systems for some of the largest Internet properties. Baron has applied his systems thinking skills to both computer systems and teams of people, and has written several books, including O'Reilly's best-selling High Performance MySQL. Prior to founding VividCortex, Baron was an early employee at Percona, where he managed teams including consulting, support, training, and software engineering. Baron has a degree in Computer Science from the University of Virginia.



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Introduction

What is Scalability?

Scalability is ambiguous for many people—a vague term often tossed about in conference presentations, for example. It's often used in ways confusingly similar to performance, efficiency, capacity, availability, and many other terms related to making things big and fast.

Wikipedia's definition of scalability, borrowed from a 2000 paper by André B. Bondi, is "Scalability is the capability of a system, network, or process to handle a growing amount of work, or its potential to be enlarged in order to accommodate that growth." This isn't wrong, but it's still a bit informal, and this ebook needs a more formal definition.

Dr. Neil J. Gunther provides one such definition: scalability is a *function*. I read Dr. Gunther's books and heard him speak for quite a while before that sunk in for me. Scalability can be defined as a mathematical function, the relationship between independent and dependent variables (input and output).

- Most important is getting the x-variable right

Bondi's definition provides a good clue: *work* is the driving factor of scalability. Useful ways to think about work include, to mention a few,

- units of work (requests)
- the rate of requests over time (arrival rate)
- the number of units of work in progress at a time (concurrency)
- the number of customers or users sending requests

Each of these can be sensible independent variables for the scalability function in different circumstances. For example, in benchmarks it's quite common to configure the number of threads the benchmark uses to send requests to a database. The benchmark usually sends requests as fast as possible, assuming zero think time, so the arrival rate is related to, but not strictly controlled by, the benchmark configuration (since it is determined by how quickly the database finishes each request).

One might say that load or concurrency is the input to the benchmark's scalability function, and the completion rate is the output.

In another scenario, you might vary the number of CPUs for the system under test (SUT) while holding constant the load per CPU, or if it's a clustered database, vary the cluster size and hold constant the load per node. In this case, the independent variable is the system resources.

In most cases I've analyzed, either concurrency or resources are usually sensible independent variables for a scalability function.¹ So for the purposes of this book, we'll consider scalability to be *a function of concurrency or capacity*. The dependent variable is usually the rate at which the system can process work, or *throughput*.

XXX chart with axes

For those who are like me and need extra emphasis, I'll repeat that this is a mathematical function, with concurrency or capacity on the X axis, and throughput on the Y axis.

Linear Scalability: The Holy Grail

In my experience, never was a marketecture slide deck created that mentions scalability without also including the word "linear." But like many

¹ I have seen other scenarios, such as [scaling by number of shards](#), but we don't need to dig into that in this book.

other things in scalability, that word is *horribly* abused.

XXX PRINCESS BRIDE MEME

Hand-waving claims of linear scaling usually coincide with vague definitions of scalability, and people who know a lot about scalability rarely say the word “linear.” Here are a few of the misdefinitions of linear scalability I’ve heard:

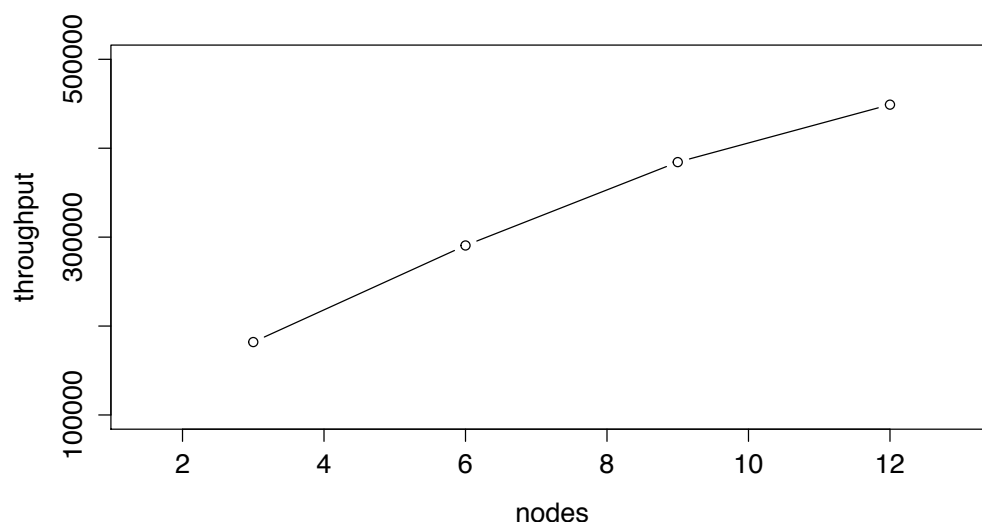
- A web architect at a conference said, “I designed our system to be shared-nothing so it would be linearly scalable.” He meant there was no single resource or system imposing a hard upper limit on how many servers could be added to the system. But he didn’t really know whether his system actually scaled linearly.
- A technical evangelist giving a presentation about a clustered database said, “adding a node to the cluster adds a predictable amount of capacity.” Predictable isn’t the same as linear.
- A sales presentation for another clustered database said the database “scales linearly, with a linearity factor of 97%,” meaning that each additional node increases the system’s capacity by 0.97 times the amount the previous node added. That’s a curve, not a line. (Later you’ll learn how to instantly determine the asymptotic upper bound on such a system’s total capacity.)

This may seem like a pointless rant, but it’s actually important if you want to be able to design and improve highly scalable systems.

Spotting bogus linearity claims is fun. Here are some ways to make systems appear linear:

- Show graphs without numbers, so you can’t do the math.
- Show graphs with nonlinear axes.
- Begin the axes, especially the Y axis, at a nonzero value.

Here is an example that employs some of these tricks.



Looks pretty linear, doesn't it? Yet if you do the math, it's nowhere near linear. It's an optical illusion, because the X axis begins around 1.45 instead of zero, and the Y axis starts at 100,000, so you can't tell that the chart isn't going to intersect the origin if you extend it downwards.

The real test of linearity is whether the transactions per second per node remains constant as the node count increases. The chart's original source mentioned that throughput increased from "182k transactions per second for 3 nodes to 449k for 12 nodes." The math is easy: the system achieves 60,700 transactions per second per node at 3 nodes, but only 37,400 at 12 nodes, which represents a 39% *drop in throughput* versus linear scalability. If it actually scaled linearly, it would achieve 728k transactions per second at 12 nodes.

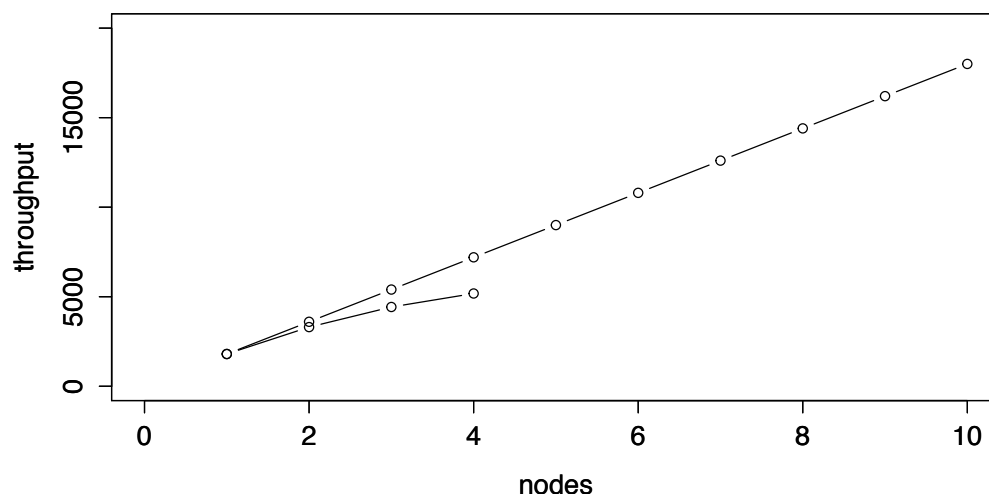
Linear means linear, folks! And seemingly small amounts of nonlinearity really matter, as you'll see later, because small sublinear effects grow very quickly at larger scale.¹

¹ In fact, they grow—wait for it—nonlinearly!

Why Systems Scale Sublinearly

Linear scalability is the ideal, yet despite the claims, systems that actually scale linearly are rare. It's very useful to understand the reasons for this, because a correct understanding of scalability, and the reasons and sources of sublinear scaling, is the key to building more scalable systems. That's why it's really important to be a linearity skeptic. It's not just being pedantic.

The best way to think about linearity is as a ratio of the system's performance at a size of 1, if possible. Neil Gunther calls this the *efficiency*. If a system produces 1800 transactions per second with 1 node, then ideally 4 nodes produce 7200 transactions per second. That would be 100% efficient. If the system loses a bit of efficiency with each node and 4 nodes produce, say, 5180 TPS, then the 4-node system is only 72% efficient:

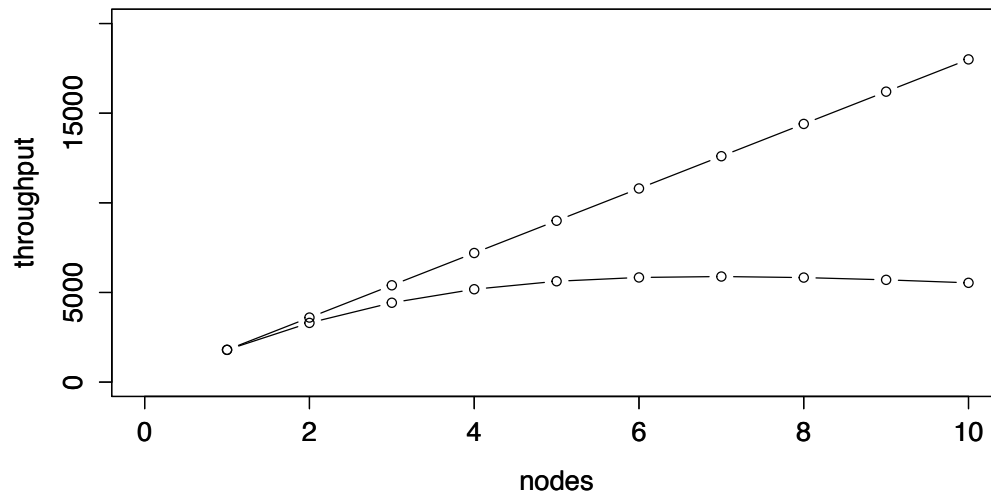


If you do this math, you'll often be surprised at how large the loss of efficiency is. Graphs can be deceptive, but the numbers are quite clear.¹

In the real world there's almost always some loss of efficiency, and if you

¹ Drawing a linear scaling line on the graph helps, too. Without that line, the eye tends to see the graph as more linear than it really is, and the loss of efficiency becomes less obvious.

can figure out why, you may be able to fix it. In fact, you've probably noticed that real systems tend not only to fall behind linear scalability a bit, but actually exhibit *retrograde* scalability at some point:



This is quite common in the real world—you scale things up and at some point your system starts going backwards and *losing* performance, instead of just gaining more and more slowly. In the MySQL 5.0 days, for example, it was common to see people getting better performance out of 4-core servers than 8-core servers.

Why does this happen? Why don't systems scale linearly, and why do they sometimes show retrograde scalability?

According to Dr. Gunther, there are two reasons: **serialization** and **crosstalk**. Serialization degrades scalability because parts of the work can't be parallelized, so speedup is limited. Crosstalk introduces a fast-growing coherency delay as workers (threads, CPUs, etc) block on shared mutable state and communication mechanisms. We'll explore these effects in the next section.

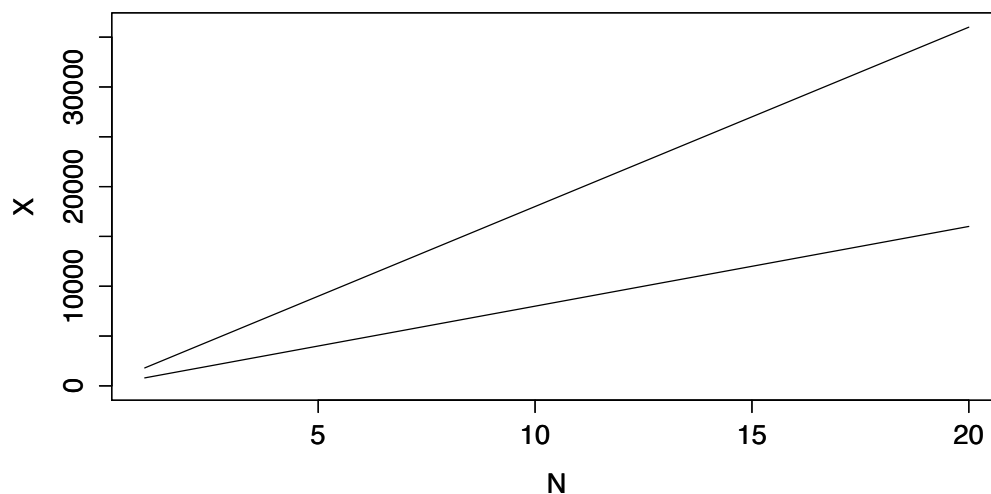
The Universal Scalability Law

Dr. Neil J. Gunther's Universal Scalability Law (USL) provides a formal definition of scalability,¹ and a conceptual framework for understanding, evaluating, comparing, and improving scalability. It does this by quantifying the effects of linear speedup, serialization delay, and coherency delay due to crosstalk.

Let's see how this works, piece by piece. An ideal system of size 1 achieves some amount λ of throughput X , in completed requests per second. Because the system is ideal, the throughput doubles at size $N=2$, and so on. This is perfect linear scaling:

$$X(N) = \frac{\lambda N}{1}$$

The λ parameter defines the slope of the line. I call it the *coefficient of performance*. It's how fast the system performs in the special case when there's no serialization or crosstalk penalty. Here are two ideal systems, with λ of 1800 and 800, respectively.



Note that every linearly scalable system is just as scalable as any other,

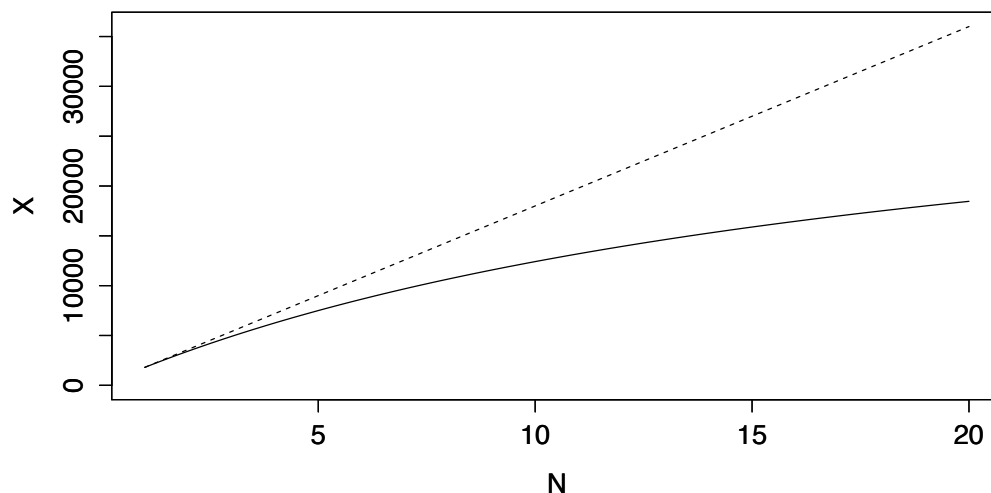
¹ Gunther originally called the USL "superserial," and you may encounter this terminology, especially in older books and papers.

regardless of the slope of the line. They have different performance but identical scalability characteristics: speedup is unlimited.

Serialization appears in most human and computer systems at some point, for example as a final stage of assembling the multiple outputs generated in parallel into a single final result. As parallelization increases, serialization becomes the limiting factor. This is codified in [Amdahl's Law](#), which states that the maximum speedup possible is the reciprocal of the serial fraction. This fits neatly into the bottom of the equation. Here σ is the serial fraction of the work, the *coefficient of serialization*:

$$X(N) = \frac{\lambda N}{1 + \sigma(N - 1)}$$

Serialization creates contention, and a system that is partially serialized will asymptotically approach a ceiling on speedup. If σ is .05, for example, the speedup approaches 20.



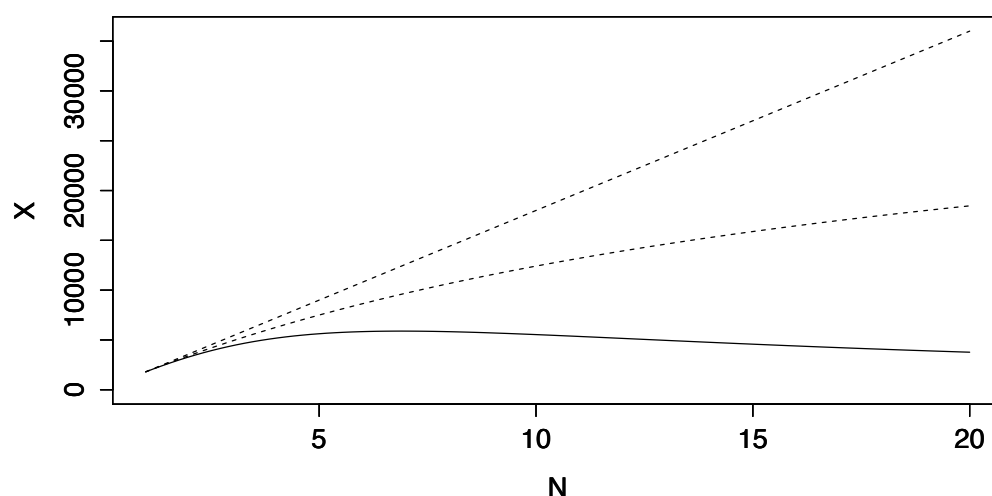
The last bit is the crosstalk penalty, also called the consistency or coherency penalty.¹ Crosstalk potentially happens between each pair of workers in the system (threads, CPUs, servers, etc). We represent the amount of crosstalk with another parameter, κ .

¹ I call it crosstalk because in my opinion it's the best description of the pairwise communication that must occur to make distributed data or other shared resources consistent or coherent.

You probably remember that the number of edges in a fully connected graph is $n(n - 1)$, which is $\mathcal{O}(n^2)$.¹ The κ coefficient quantifies the crosstalk in the denominator:

$$X(N) = \frac{\lambda N}{1 + \sigma(N - 1) + \kappa N(N - 1)}$$

The crosstalk penalty grows fast. Because it's quadratic, eventually it grows faster than the linear speedup of the ideal system we started with, no matter how small κ is. That's what makes retrograde scalability happen:



That's the Universal Scalability Law in all its glory. This plot has the same parameters as the ones I showed before, where a system of size 4 produced only 72% of its ideal output. That system has 5% serialization and 2% crosstalk, and now that I've plotted it out to size 20 you can see it's embarrassingly inefficient. In fact, we should have given up trying to scale this system after size 6 or so.

This shows visually how much harm a “small amount” of nonlinearity can do in the long run. Even very small amounts of these damaging coefficients will create this effect sooner or later (mostly sooner). This is why it's rare to find clustered systems that scale well beyond a couple

¹ If you're not familiar with it, [this blog post introduces Big-O notation](#).

dozen nodes or so. If you'd like to experiment with this interactively, I've made a graph of it at [Desmos](#).

If you're curious, the USL is based on queueing theory. It's equivalent to synchronous repairman queueing. You can read more about that in Neil Gunther's books. If you're not familiar with queueing theory, I wrote a highly approachable introduction called [Everything You Need To Know About Queueing Theory](#).

Measuring Scalability

To recap, up until now we've figured out the right dimensions for a formal model of scalability that seems to behave as we know real systems behave, and examined Neil Gunther's USL, which fits that framework well and gives us an equation for scalability. (Are you excited yet?)

Now what do we do with it?

Great question! It turns out we can do a lot of extremely useful things with it. Unlike a lot of models of system behavior, this one is actually practical to apply in the real world. That's the real genius of it, in fact. Not only is the equation uncomplicated, but the variables it describes are easy to get from a lot of systems. If you've ever tried to model system behavior with queueing theory, e.g. the Erlang C formula, you're going to love how simple it is to get results with the USL. So many modeling techniques are foiled by the inability to get the measurements you need.

The thing I've used the USL for the most is measuring system scalability, by working backwards from observed system behavior and deriving the likely coefficients. To accomplish this, you need a set of measurements of the system's size as you see fit to define it (usually concurrency or node count) and throughput. Then you *fit* the USL to this dataset, using nonlinear least squares regression. This is a statistical technique that

finds the optimal coefficient values in order to calculate a best-fit line through the measurements. The result is values for λ , σ , and κ .

If you're reading about the USL in Neil Gunther's books, he takes a different approach. First, he doesn't use regression to determine λ , he assumes that is something you can measure in a controlled way at $N = 1$. (I've often found that's not true for me.) Secondly, there are a couple of different forms of the USL—one for hardware scaling and one for software scaling—which are the same equation, but with different parameters. I've found the distinction to be largely academic, although I will talk a bit more about this later.

Examples of systems I've analyzed with the USL include:

- Black-box analysis of networked software simply by observing and correlating packet arrivals and departures, looking at the IP addresses, port numbers, and timestamps. From this I computed the concurrency by averaging the amount of time the system was busy servicing requests over periods of time. The throughput was straightforward to get by counting packet departures.
- MySQL database servers. Some of its `SHOW STATUS` counters are essentially equivalent to throughput and concurrency.
- Linux block devices (disks) by looking at `/proc/diskstats`, from which you can get both instantaneous and average concurrency over time deltas, as well as throughput (number of I/Os completed).
- Lots and lots—and lots—of benchmark results.

I've built a variety of tools to help clean, resample, and analyze the data before arriving at satisfactory results. Most of them were commandline, though these days I use R more than anything else. This is an important topic: you will get dirty data, and that will make your results less useful. You need to visualize both in scatterplot form as well as in time-series form and ensure you're working with a relatively consistent set of data.

You can remove individual points or trim the time range you use, and you may need to experiment with averaging the data over time to get good results.

As for the R code, I'll give a little bit of a quickstart to show the soup-to-nuts approach. You'll save the data into a delimited file, with column headers `size` and `tput`. Then you'll load this into a variable in R and regress it against the USL.

Here's a complete sample, based on a [benchmark](#) that Vadim Tkachenko ran at Percona:

```
size tput
1 955.16
2 1878.91
3 2688.01
4 3548.68
5 4315.54
6 5130.43
7 5931.37
8 6531.08
9 7219.8
10 7867.61
11 8278.71
12 8646.7
13 9047.84
14 9426.55
15 9645.37
16 9897.24
17 10097.6
18 10240.5
19 10532.39
20 10798.52
21 11151.43
```

```

22 11518.63
23 11806
24 12089.37
25 12075.41
26 12177.29
27 12211.41
28 12158.93
29 12155.27
30 12118.04
31 12140.4
32 12074.39

```

Save that data into a file, say, `benchmark.txt`. Then load it and run the following commands:

```

benchmark <- read.csv("/path/to/benchmark.txt", sep="")
usl <- nls(tput ~ lambda*size/(1 + sigma * (size-1) + kappa * size *
  (size-1)), benchmark, start=c(sigma=0.1, kappa=0.01, lambda=1000))
summary(usl)
sigma <- coef(usl)['sigma']
kappa <- coeff(usl)['kappa']
lambda <- coef(usl)['lambda']
u=function(x){y=x*lambda/(1+sigma*(x-1)+kappa*x*(x-1))}
plot(u, 0, max(benchmark$size)*2, xlab="Size", ylab="Throughput", lty="dashed")
points(benchmark$size, benchmark$tput)

```

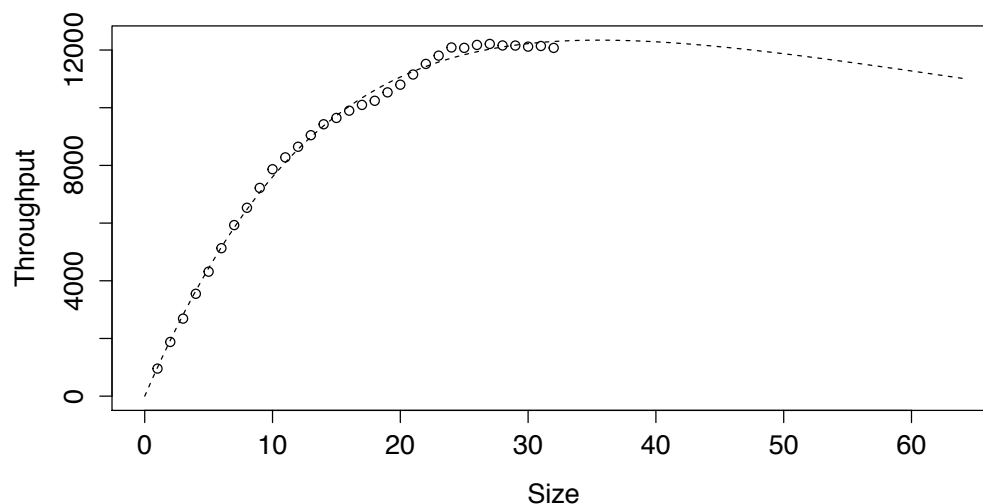
The results are as follows:

```

λ  995.6486
σ  0.02671591
κ  0.0007690945

```

Note the extremely small value for κ which nonetheless degrades scalability before N becomes very large. Here's the resulting plot:



If you're an R user, that's probably all you need to get going. You really should do more diligence, such as checking the R^2 value of the fit. But instead of doing all this work manually (which you can certainly do if you want), I suggest trying the [USL package from CRAN](#). It has many features built in, although it does have some limitations.

One final thing: if the κ coefficient has a nonzero value, the function has a maximum. You can find the size of the system at that maximum as follows:

$$N_{max} = \left\lfloor \sqrt{\frac{1 - \sigma}{\kappa}} \right\rfloor$$

Of course, to find the maximum predicted throughput, you just plug N_{max} into the USL equation itself. Doing so with the coefficients in this example predicts the system's throughput will increase until $N = 35$, which in this case means 35 threads, and the peak throughput will be 12341 queries per second. It also found λ , the throughput at $N = 1$, to be 995 QPS, which is close to the actual value of 955.

It's always interesting to use the USL on a subset of the performance data, such as the first third or so, to see how well it predicts the higher N values. This can be quite educational.

Note that you should have at least half a dozen or so data points in order

to get good results in most circumstances. In practice I usually try to capture at least a dozen for benchmarks, and more—often thousands—when analyzing systems that aren't in a controlled laboratory setting.

Relating Scalability and Performance

Throughput and latency are two common measures of performance. Most benchmarks measure overall system throughput, and claims of performance are almost always in throughput terms: “a million transactions per second,” and so on. Benchmarks usually define performance narrowly as how much work the system can do.

On the other hand, users care mostly about the performance of individual requests. A user's opinion about your website's performance is based entirely on how quickly pages load and render. From this viewpoint, as Cary Millsap says, *performance is response time*.

Which view is right? Both. System performance is measured in throughput, and request performance is measured in latency.

We've seen that the USL can model and forecast how system size affects throughput. Can it also model latency? Yes, it can, because of a relationship called Little's Law:

$$N = XR$$

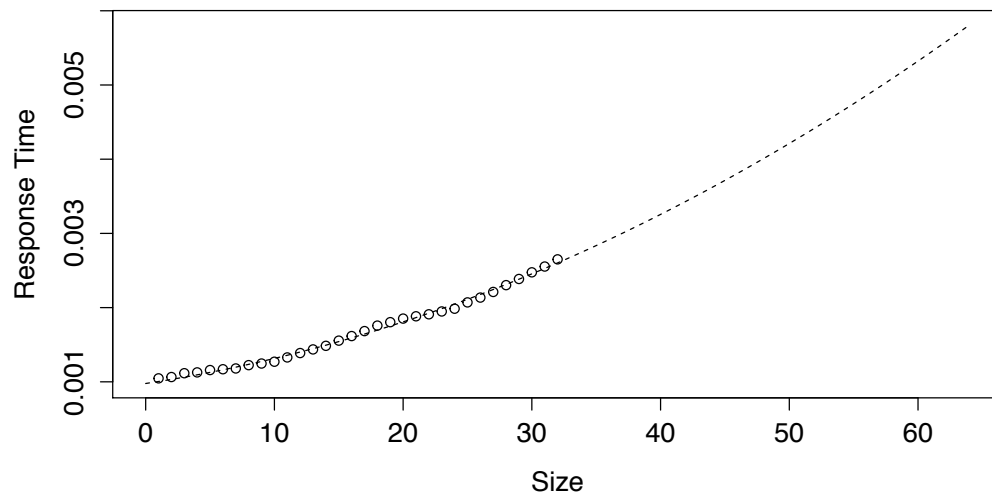
Little's Law says that the mean number of requests resident in a system is equal to the throughput times the mean response time. This relationship is valid for stable systems, in which all requests eventually complete.

It's straightforward to use Little's Law together with the USL and solve for

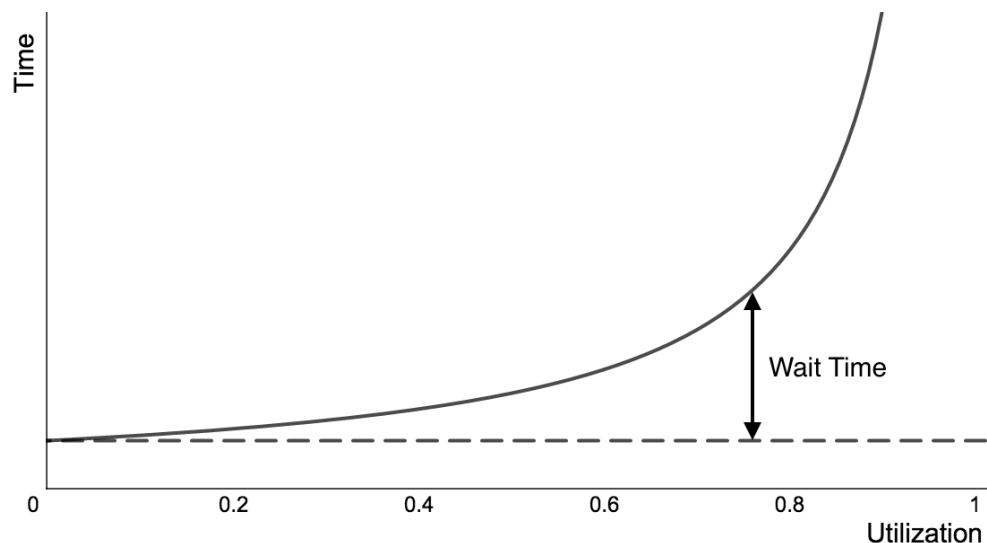
response time as a function of concurrency. The result is a quadratic function:

$$R(N) = \frac{1 + \sigma(N - 1) + \kappa N(N - 1)}{\lambda}$$

So if the USL is correct, response time is related to the square of concurrency. Here's the same data we saw previously, now inverted and used to predict response time.



Be careful not to confuse this chart with another famous "hockey stick" chart, that of response time versus utilization, which is familiar from queueing theory:



The difference is that utilization ranges only from 0 to 1, whereas concurrency has no fixed upper limit.

The inverted USL predicts mean response time. If you wanted to, you could use queueing theory to assess this in a more nuanced way, such as the probability that any given request needed to wait more than a set amount of time.

Capacity Planning with the USL

“How much load can this system sustain?” is a common question in capacity planning. The practical purpose is usually something like the following:

- How soon will the system begin to perform badly as load increases?
- How many servers will I need for the holidays?
- Is this system close to a breaking point?
- Is the server overprovisioned? By how much?

Capacity planning is often a difficult problem because it's hard to tell what a system's capacity really is. The USL can be helpful in understanding this, to a point.

There are conventional ways to determine system capacity, but they're often difficult, expensive, and don't give results you can really believe in. For example, you can set up load tests, but it takes a lot of work and time, and the results are really suspect because the workload is always artificial in some way.

You can also run benchmarks, but most benchmarks are pretty useless for predicting a system's usable capacity. In addition to being an artificial workload, they push a system to its maximum throughput and beyond.

As I mentioned, it's rare for benchmarks to be run by people who understand the importance of latency. But when I do see benchmarks that measure latency percentiles, the systems almost always perform very badly at their peak throughput.¹

Another way I've tried to predict system capacity in the past is with queueing theory, using the Erlang C formula to predict response time at a given utilization. Unfortunately, this requires that you know request service times, which are often impossible to obtain. You can measure total response time, but that includes waiting time in the queue, so it's not the same thing as the service time. The utilization is also often deceptive, because the real utilization of the resources you're trying to measure can be difficult to measure correctly too. Most people I know consider the Erlang approach to be difficult to apply.

If load tests, benchmarks, and queueing theory are difficult to use, can the USL help? To a point, yes, it can. The USL's point of maximum predicts the system's maximum throughput, so it's a way to assess a system's capacity. Because the USL is a *model*, it can help you use observed data to predict how a system will perform under load beyond what you can observe. As a result you can get a much better idea of how close you are to the system's maximum capacity.

The USL has a few nice properties:

- It's a "black box" technique, which uses data that's usually easy to get.
- Gathering data and using regression to analyze it is also easy.
- The USL is a relatively simple model, so people like me can understand the math.
- The USL is highly intuitive in comparison to most other approaches.

¹ This is a problem with the way benchmarks are usually designed, in my opinion. I'd really prefer for the benchmark system to be intelligent enough to back off and reduce pressure on the system under test if it violates a service level objective (SLO). Ideally, this is defined as a quantile, such as 99th percentile latency less than 10ms. A smart benchmark would throttle load, eventually finding a longterm stable arrival rate at which the SUT can consistently perform well.

One of the major caveats is that maximum throughput isn't maximum *usable* capacity. Again, when the system is at its maximum throughput, response time is probably terrible, and will be extremely inconsistent. That's why it's more important to focus on the system's maximum throughput within the constraints of a service level objective.

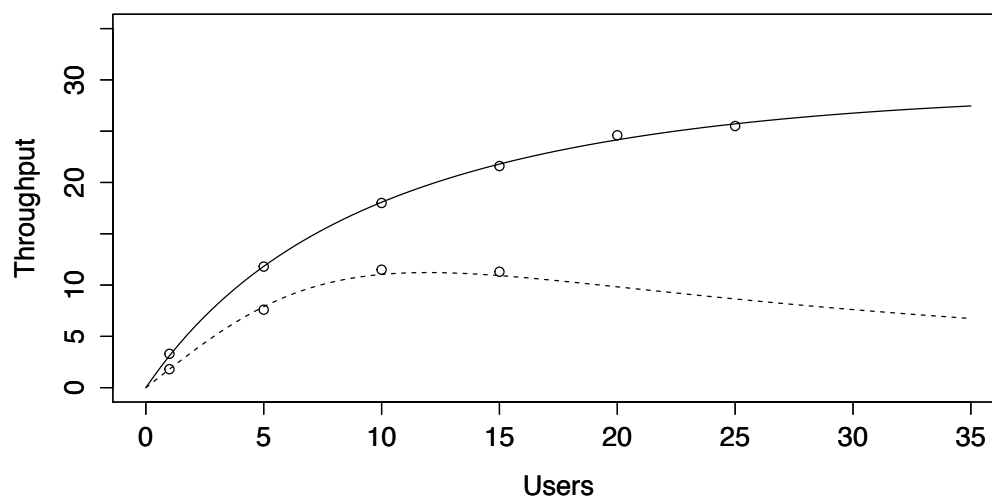
Another important thing to realize is that a lot of systems perform worse than the USL predicts they will, because their degradation in scalability at larger sizes is more severe than predicted. As a result, I suggest viewing the USL's prediction as optimistic: "I won't count on being able to scale this system as high as the USL predicts I can."

You can also combine the USL with selected techniques from queueing theory, such as the Square Root Staffing Rule, to understand how much capacity is needed and what quality of service it will provide. See the aforementioned [queueing theory ebook](#) for more on this topic.

Using the USL to Improve Scalability

One of the best uses of the USL is to evaluate and understand *why* a system doesn't scale as well as it might. Armed with this, you might know where to look for bottlenecks, so you can alleviate them and improve the system's scalability. With practice, you'll also develop a mindset of scalability, building intuition about which design decisions might cause serious sublinearity.

An example will help illustrate. A few years ago PayPal published [benchmark results](#) of a Java application they rewrote in NodeJS. I analyzed their benchmark results and [wrote](#) about it on the VividCortex blog. Here are the plots and the key scalability parameters:



	σ	κ
Java	0.000011	0.006323
NodeJS	0.080319	0.000222

These systems scale very differently,¹ and for very different reasons. In a nutshell, the Java benchmark shows much higher crosstalk penalty, whereas the NodeJS benchmark exhibits more serialization. Examining the architectures of the two systems reveals why: the Java app is multi-threaded and NodeJS is single-threaded with an event loop, and the PayPal blog post even mentions that they used “a single core for the NodeJS application compared to five cores in Java.”

This is a great real-life example of key scalability tenets:

- Avoid serialization and make things as parallelizable as possible.
- Avoid crosstalk and blocking or queueing.

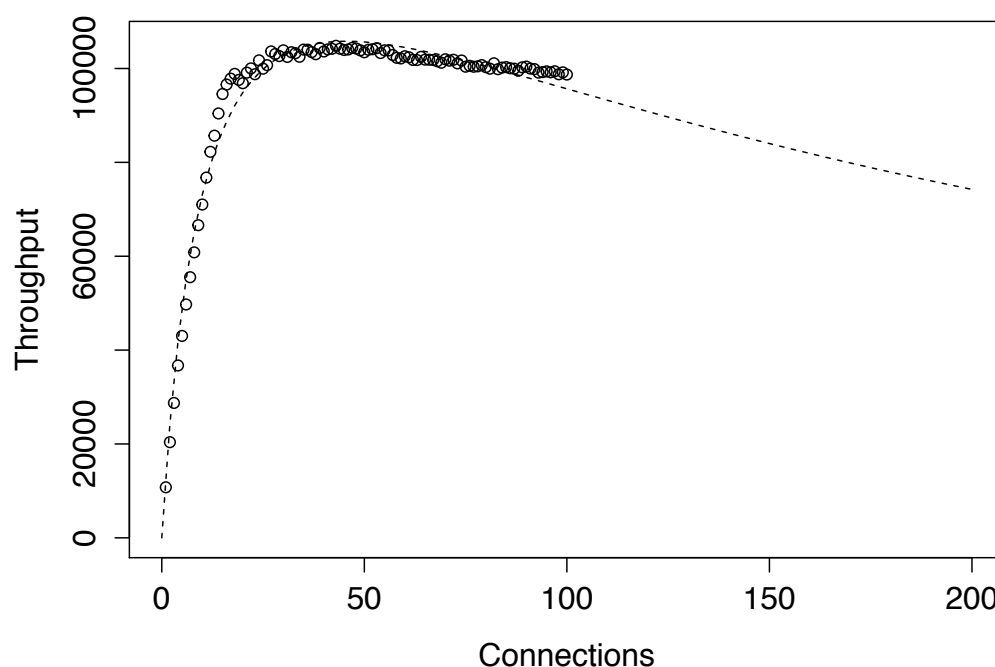
Thinking Critically About The USL

¹ Both of them scale pretty poorly, in fact.

For many people, the USL is a huge shift in mindset. I know it was for me. But is it the be-all and end-all? Of course not.

Not only is the USL unsuitable for solving every problem related to scalability or capacity planning, sometimes it doesn't work all that well for the problems it can solve. The examples I've shown thus far are remarkable in that they're extremely clean data. Real-world systems often have noisy data that doesn't "look like" the USL much at all. It may look more like something your cat threw up on the screen. Even when there is a nice-looking shape to the data, regression can produce unphysical results, or refuse to produce results.

And then there are the systems for which you have nice clean-room measurements, highly reproducible, but they just don't fit the USL very well. I could show many examples. I've seen systems that appear to scale nicely, following the USL, but suddenly flatline instead of continuing a graceful curve, or abruptly degrade way faster than predicted—or the reverse, appearing to degrade more *slowly* than predicted once we exit Amdahl territory and enter the region where the higher-order terms prevail. Here's an example:



Maybe this is queue saturation or resource saturation, or maybe it's something else, but the USL is unaware of it.

What, then, is the USL good for? The answer is lots.

First of all, a model is better than no model. In the absence of a model explaining the workings of system scalability, there isn't even a point of reference against which you can compare your expectations and results. There's no basis for a conversation. There's no frame of reference to say, "I think this system should scale better than it does," or "This system's behavior makes no sense." Whether the USL is usable or not, it still provides a framework. Without it, why not just draw lines at will? You could get a set of French curves and draw curves as directed by your muse. No one could say you're wrong.



In an objective sense, the USL is incomplete or wrong, perhaps both. I don't think it matters very much, but if it were complete and correct, it would describe more systems and scenarios than it does in my experience. As Richard Feynman said in a 1964 lecture at Cornell University, "If it disagrees with experiment, it's wrong. That simple statement is the key to science. It doesn't make a difference how beautiful your guess is, it doesn't make a difference how smart you are, who made the guess or what his name is—if it disagrees with experiment, it's wrong. That's all there is to it."

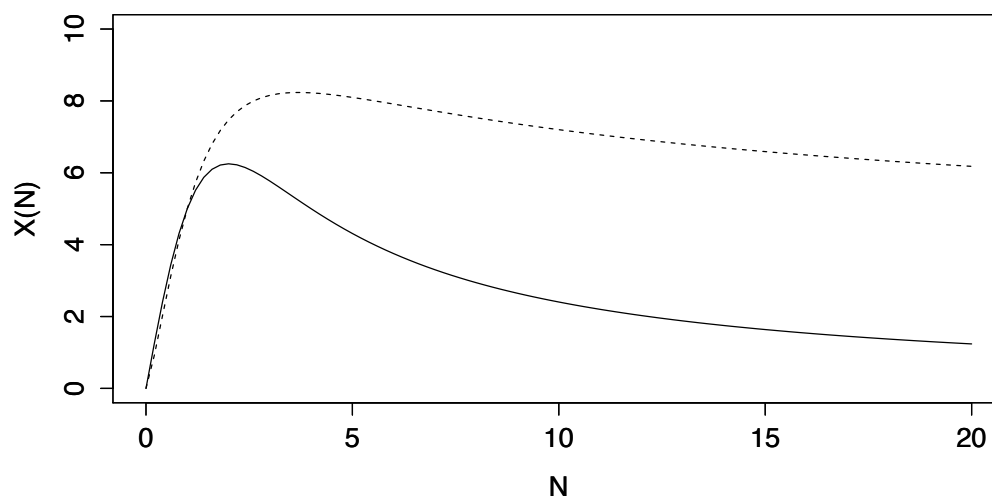
This is not to say the USL isn't useful. It's incredibly useful. But we must

not put it onto a pedestal and worship it.

I'd also like to note that one could quite easily conjecture and analyze other USL-like models. For example, given that many computer algorithms can be shown by analysis to be $\mathcal{O}(n \log n)$ instead of $\mathcal{O}(n^2)$, perhaps the quadratic κ term could be replaced by a logarithmic term, like so?

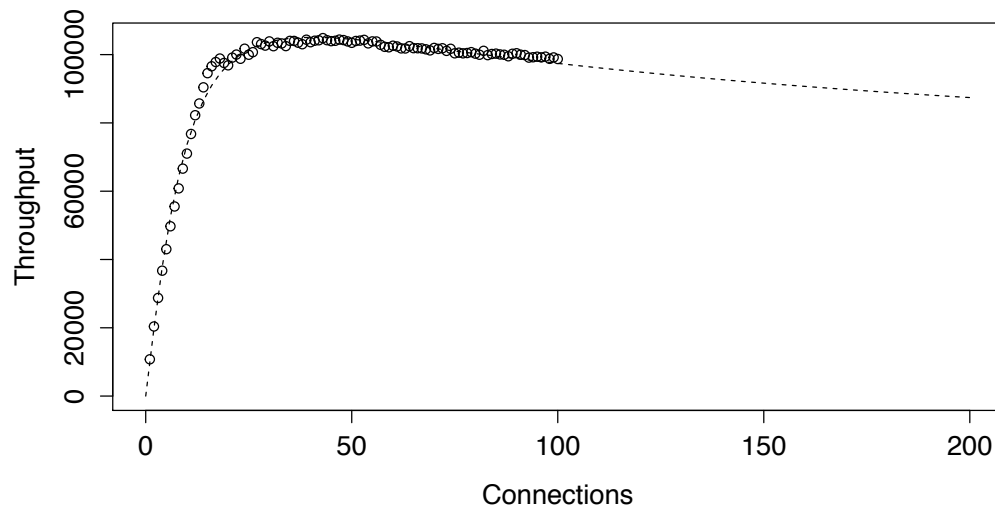
$$X(N) = \frac{\lambda N}{1 + \sigma(N - 1) + \kappa \log(N)(N - 1)}$$

To give some visual intuition of how this differs from the accepted form of the USL, here they are together on a single plot. The standard USL is the solid line and the logarithmic variant is the dashed line.



You might think that you could choose parameters to the standard USL to make it follow the dashed line better, but it doesn't work. The functions are of different order and won't behave the same.

Would this be a better model for the data I showed previously? Visually, it does appear to model the observed data somewhat more closely:

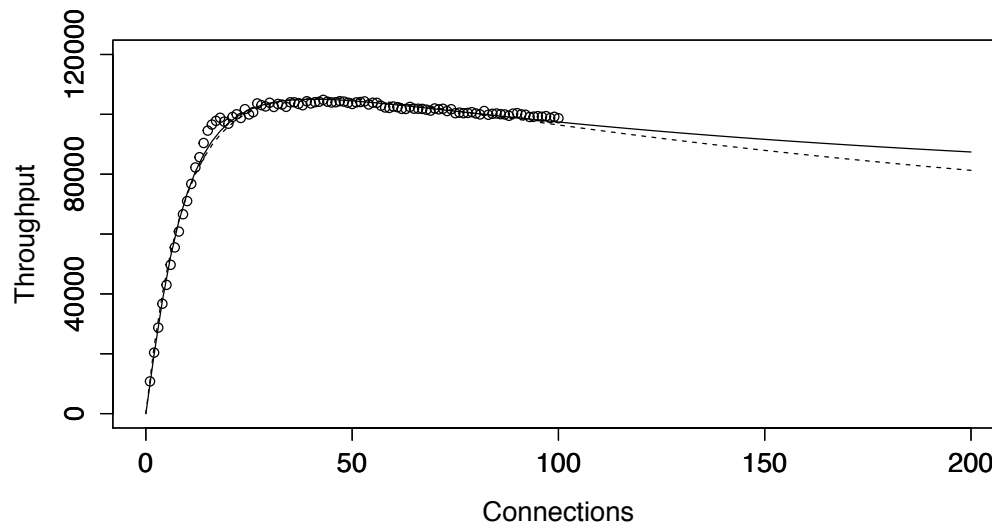


Jayanta Choudhury has suggested changes to the USL to model cases such as this, where resources apparently become saturated. His proposed model, called the Asymptotically Improved Super-Serial Law,¹ is slightly more complex:

$$X(N) = \frac{\lambda N}{1 + \sigma(N - 1) + \sigma \kappa N^{\beta}(N - 1)}$$

The β parameter ranges from 0 to 1, inclusive. Unfortunately it's more difficult to estimate parameters for this model, making it a bit harder to use than the USL. The following plot shows the logarithmic extension I proposed above in a solid line, plotted with the AISSL in a dashed line, on the same dataset. As can be seen visually, the logarithmic variation fits the data better.

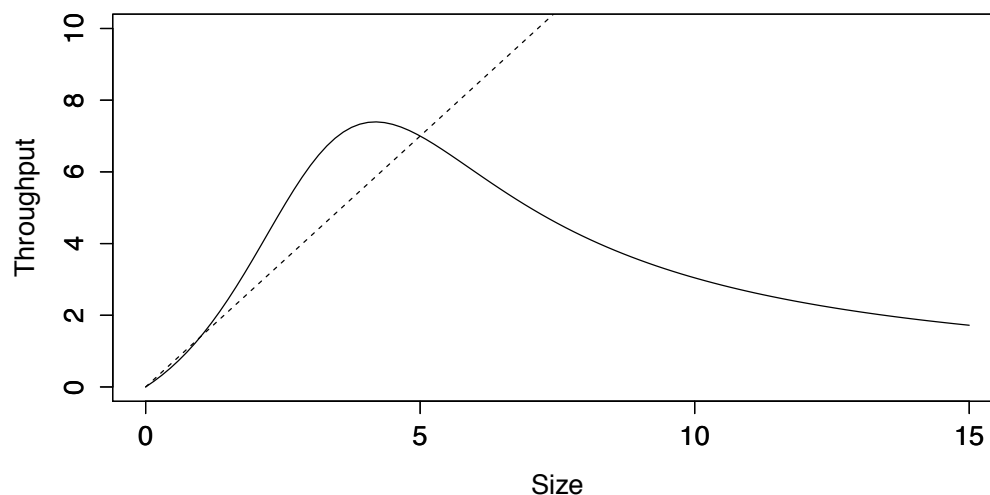
¹ See *Parameter Estimation of Asymptotically Improved Super-serial Scalability Law* by Dr. Jayanta Choudhury, of TeamQuest Corporation



As I'm sure you can imagine, you could play games like this all day long. Dreaming up an idea is much easier than proving that it's correct or showing how it might arise analytically from the underlying mechanisms as we understand them to operate in the system.

Superlinear Scaling

I've spent some time talking about the possible causes of sublinear scaling. Is superlinear scaling possible? As I've worked with the USL over the years, I've found a number of cases where systems apparently do scale superlinearly. It manifests as a negative κ coefficient and a USL curve that has a more complex shape:



At first I dismissed this result as unphysical (how can there be less than zero contention?), but after repeatedly seeing this happen and having many conversations with Neil Gunther about it, I started to wonder. So did Neil, and eventually he was able to reproduce and explain the effect on a [large-scale Hadoop TeraSort benchmark](#).

The TeraSort case is quite a specific one. In general, I would explain superlinear scalability as a disproportionate scaling of some resource relative to the load placed upon it, creating an economy of scale. For example, adding more nodes to a clustered database system adds more memory; if the dataset size is not scaled proportionately, then more of the data fits into memory on each node, and access times improve relative to disk reads. Any resource that is more efficient when shared than when used singly may cause this effect.

It's worth noting that this initial boost, depending upon its cause, may be countered by a correspondingly disproportionate “payback” later when performance falls quickly below linearity again.

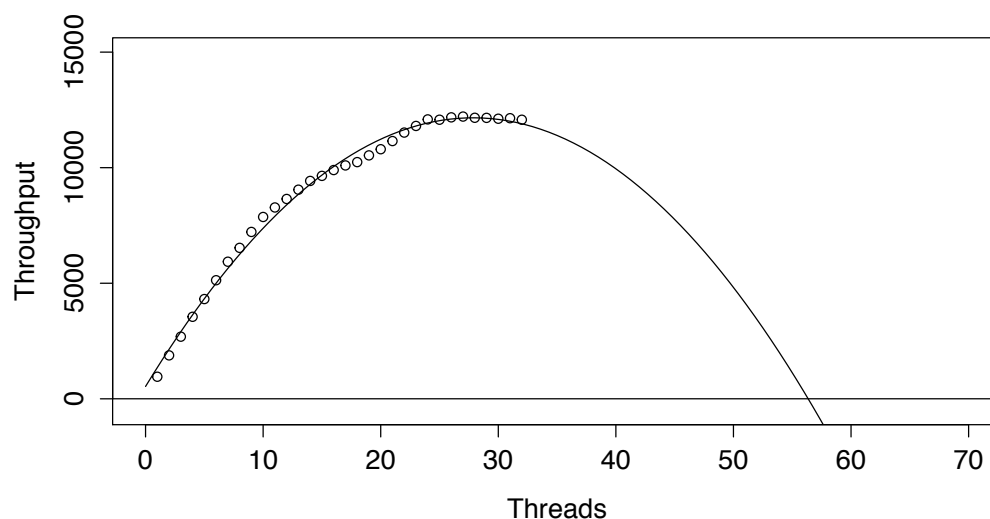
Another special case to be aware of is that some clustered systems behave differently at sizes 1 and 2 than they do at 3 and above. For example, at size 1 there is no crosstalk or contention—it isn't a distributed system. At size 2 special-cases may be in play. At size 3 and above,

usually generic algorithms and techniques suited for any size n are in use. Some clustered systems have to be benchmarked or measured at larger sizes in order to ignore these effects.

Other Scalability Models

In addition to the USL and the variants I've already discussed, there are many other potential models of scalability you could consider. In my opinion some are good, some are useful, some are not. I also believe that there is much more work to be done on this topic.

Some alternative theories, however, are just garbage. Chief amongst them is “quadratic scalability.” Observing that systems under increasing amounts of load will first increase in throughput, then level off and begin to decline, some people get the bright idea that they should “fit a curve to it.” The curve is always a quadratic polynomial and the fitted curve ends up being a parabola opening downwards. Let's see how this looks on the benchmark data from Percona once again:



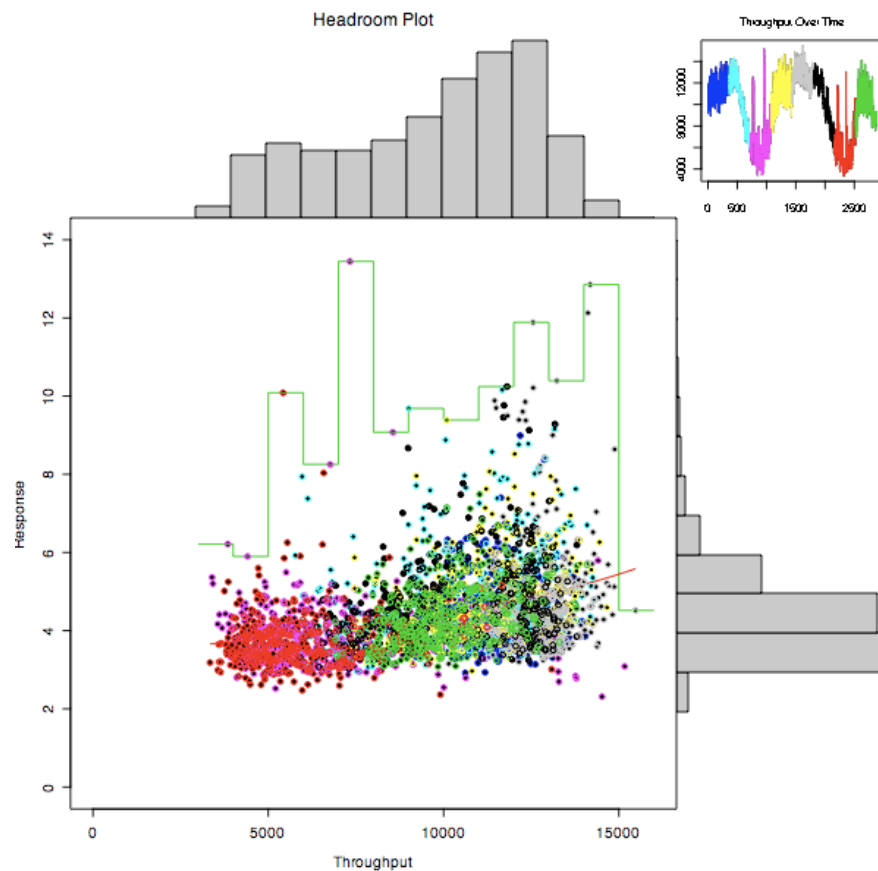
Look, Ma, it's a great fit! I don't even need to compute the R^2 value to know that. There's just one problem: it predicts that at some point we'll achieve negative throughput.

You should ignore this model because by definition it doesn't work. Rather than use it, I'd suggest that you get out that French curve set again and get in touch with your inner artist.

A few other models that I've seen expressed in graphical form, but not mathematically and lacking analysis, are:

- New Relic's scalability chart, which plots latency as a function of throughput and renders a smoothed line through the points. The line has no predetermined form and doesn't express any particular model.
- AppDynamics's scalability analysis feature does the same thing, but fits a parabola through the lines instead of a polynomial of arbitrarily high degree.
- Cockcroft headroom plots, by Adrian Cockcroft. These are also latency-versus-throughput charts, but add some histograms and other useful visual cues around the edges.

These three express essentially the same belief about the relationship between variables; the differences are relatively minor. Here's an example of a [Cockcroft Headroom Plot](#):



At some point I intend to analyze what it means to treat latency as a function of throughput, in the context of queueing theory or the USL. My initial reaction is that throughput isn't a proper independent variable, but I confess to being conflicted on this point. As long as you stay within the bounds of a stable system I think it might work. Please let me know if you do this work (e.g. rearranging the USL using Little's Law to derive response time as a function of throughput).

Further Reading

* GCaP * Look at the Percona white paper all of the books in my picture



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