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FLUID METHODS FOR MODELING LARGE, HETEROGENEOUS NETWORKS

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13. ABSTRACT (Maximum 200 Words) Researchers from the University of Massachusetts developed fluid-based methodologies for characterizing the behavior of large IP networks handling large numbers of TCP and UDP flows. These methodologies provide for rapid and efficient rates of individual and aggregate flows. These fluid models were also applied to the problems of characterizing the spread of worms and viruses and to the cascade of failures within the BGP routing infrastructure. The resulting fluid models were used to develop novel active queue management mechanisms resulting in more stable TCP performance and novel rate controllers for the purpose of providing minimum rate guarantees to TCP flow aggregates. Last, methodologies and tools were developed to integrate fluid network simulation with packet-level simulation.				
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1 Introduction

The goals of this project were (1) to develop a theory for modeling large-scale IP networks based on fluid models and the resulting solution techniques, (2) to integrate these methodologies into existing tools, such as *ns* [1]; and to use the theory and tools to develop new control strategies for large-scale networks.

The main outcomes of the project were the following:

- **Fluid models for large networks carrying TCP and non-TCP flows:** Based on the use of stochastic differential equations, we developed simple models characterizing the behavior of TCP and non-TCP flows as they traverse through a network. These models characterize the control loop behavior of TCP flows and their interactions with routers. The time varying behavior of average quantities, such as queue length, packet loss probability, etc., is captured through a set of differential equations. Stationary behavior is characterized through a fixed point problem.
- **Fluid models for Internet worm and BGP failure spread:** We developed similar models that describe the spreading behavior of Internet worms, such as Code Red, and of cascading router failures.
- **Congestion controllers for routers:** We developed new active queue management policies, based on proportional integral control (PI) that provides more robust and stable TCP behavior. These controllers were extended to provide quality of service in the form of minimum throughput guarantees to TCP flows.
- **Fluid simulation tools:** We developed tools for solving the performance of large networks based on the fluid techniques. Moreover, tools that combine packet-level with fluid-level simulation were also developed. These were integrated into several existing simulation tools including *ns*.

We describe each of these outcomes in the remainder of this report. Prior to this, we motivate the need for our project.

2 Motivation

Networks continue to increase in size and complexity. The variety and interaction of the applications, the middleware and transport protocols, the routing protocols, and the router/switch resource management algorithms make the design, development, control and management of future networks an exceptionally difficult task. These difficulties can be eased by the development of analytic and numerical techniques (as opposed to detailed, brute force simulation) of such systems.

In the past, discrete-event simulation has been the performance-modeling tool of choice. This has been primarily due to the predominance of responsive flows (e.g., TCP) in network workloads, and the difficulties of analytically or numerically characterizing the performance of such flows. Although considerable progress has been made in modeling open-loop streaming applications

(e.g., video over UDP), these advances have had little impact on the modeling of responsive flows. While significant progress has been made in simulating large networks with responsive flows, such simulations are still very time consuming and require significant computing resources. This motivates the need for scalable models of networks that can be efficiently solved to obtain performance measures such as throughput, link loss rates, end-to-end packet delays, which in turn can be used to estimate application-level performance measures.

The remainder of this report is organized as follows: we describe the fluid modeling methodology for TCP/IP networks in Section 3 and its adaptation to the characterization of worm and failure spreading behavior in Section 4. In Section 5, we show how the differential equations describing the behavior of TCP/IP networks can be used to develop congestion controllers at routers to interact with TCP to provide different performance to different classes of sessions. Following this, we describe the tools developed during the project (Section 6). Section 7 offers some conclusions.

3 Fluid Models for TCP/IP networks

In this section, we summarize our work on modeling TCP/IP networks. We begin with a description of our network model. Let V be a collection of routers making up a network. Each router¹ $v \in V$ has a transmission capacity of C_v bits per second. In addition, router v can buffer up to B_v bits. Associated with v is an active queue management policy characterized by a probability discard function, $p_v(x_v)$, which takes as its argument x_v , the average queue length of v . One popular discard function, associated with the RED (random early discard) active queue management mechanism [2], is

$$p_v(x_v) = \begin{cases} 0 & 0 \leq x_v < t_v^{\min} \\ \frac{x_v - t_v^{\min}}{t_v^{\max} - t_v^{\min}} p_v^{\max}, & t_v^{\min} \leq x_v < t_v^{\max} \\ 1, & t_v^{\max} \leq x_v \leq B_v \end{cases} \quad (1)$$

where $t_v^{\min}$, $t_v^{\max}$, and $p_v^{\max}$ are configurable parameters. We make the reasonable assumption that p_v is non-decreasing in x_v .

3.1 Time varying models of TCP/IP networks

The behavior of a TCP/IP network is stochastic in nature. Hence, we find that it is best described using stochastic differential equations. We begin with a treatment of a single link and then show how it generalizes to a network setting.

3.1.1 Single link

Stochastic differential equations (SDEs) have been widely used in system modeling. Most of the effort has been directed to Wiener-process-driven SDEs. However, the primary sources of

¹ More precisely, each router interface has a transmission capacity, a buffer, etc. We will use the term router to refer to an outgoing router interface.

randomness in high-speed networks are discrete events such as losses in the TCP congestion control setting. Such a network can be modeled by stochastic differential equations driven by jump processes. For example, the additive increase of the TCP window size can be modeled by a differential equation, while the multiplicative decrease triggered by the losses can be modeled by a jump process. We use jump processes to model the packet losses inherent in the TCP-AQM dynamics. Since the marks delivered by proposed AQM algorithms are assumed to be randomly generated, we use Poisson processes with random rates that are functions of the (average) queue lengths at the routers. The jumps in these processes represent the losses in the TCP-AQM dynamics.

Consider a single congested router operating with a RED AQM scheme, as described earlier, supporting N TCP flows. The discard probability is $p(x(t))$ where $x(t)$ is the average queue length at the router at time $t > 0$. In particular, x is an *estimate* based on the instantaneous queue length $q(t)$ via an exponential weighing parameter α and sampling interval δ (i.e., the queue length $q(t)$ is sampled periodically every δ units of time). The rate, λ_i , at which losses are incurred by flow i at time t is $p(x(t)) T_i$, where T_i is the throughput of that flow, $W_i/R_i(q)$.

We have the following stochastic differential equation for the window size of flow $i = 1, \dots, N$:

$$dW_i = dt / R_i(q) - W_i / 2 dN_i.$$

Since the loss generated by the RED AQM policy has independent increments (i.e., $dN_i(t)$ is independent of W_i), the expected window-size process is described by

$$E[dW_i] = dt / R_i(q) - E[W_i] / 2 E[dN_i].$$

Since $E[dN_i] = p(x) E[W_i]/R_i(q)$, we have

$$\frac{dE[W_i]}{dt} = \frac{1}{R_i(q)} - \frac{E[W_i]^2}{2R_i(q)} p(x) \quad (2)$$

We write the exponentially weighed moving queue length average as

$$x((k+1)\delta) = (1-\alpha)x(k\delta) + \alpha q(k\delta) \quad (3)$$

This equation contains the AQM sampling parameter δ . Finally, we need an equation for the expected instantaneous queue length $q(t)$. For nodes that are busy most of the time the following equation is a good approximation:

$$\frac{dq(t)}{dt} = -\mathbf{I}(q(t) > 0)C + \sum_{i=1}^N E[W_i] / R_i(q). \quad (4)$$

This results in $N+2$ coupled equations and $N+2$ unknowns, $x, q, E[W_i(t)], i = 1, \dots, N$ which can be solved numerically using a numerical scheme such as the Runge-Kutta method. The solution provides an estimate of the average transient behavior of the system for a given sampling interval δ .

3.1.2 Extension to the network case

The extension to the network case is fairly straightforward. We represent the network by a binary matrix $\mathbf{A}$ where the rows represent the different flows and the columns represent the different routers (queues). We modify equation (2) to account for losses arriving from each router in the path. If $\mathbf{W}$ is defined as the vector of the window sizes of the flows, $\mathbf{x}$ the vector of estimated average queue lengths at each router, $\mathbf{q}$ the vector of instantaneous queue lengths at each router, and $\mathbf{P}(\mathbf{x})$ the vector of loss probabilities at each router, then we define a matrix $\mathbf{AP}$ where every column of $\mathbf{A}$ is multiplied by the corresponding element of $\mathbf{P}$. Thus, $\mathbf{AP}$ is an indicator of the loss probabilities seen by each flow at each router. If a flow does not traverse a particular router, the corresponding loss probability is set to 0. The combined loss seen by a particular flow i is then given by $1 - \prod (1 - \mathbf{AP}(\mathbf{x})_i)$. Thus, (2) is modified to

$$\frac{dE[W_i]}{dt} = \frac{1}{R_i(\mathbf{q})} - \frac{E[W_i]^2}{2R_i(\mathbf{q})} (1 - \prod (\mathbf{AP}(\mathbf{x})_i)) \quad (5)$$

The equations for the average queue size and instantaneous queue size at each router remains unchanged.

The model described above captures the important details of a network supporting TCP flows. There are some characteristics of such networks that are missing. For example, the model so far assumes that congestion feedback is instantaneous. This, of course, is not true, as there is at least a round trip propagation delay before feedback is returned. In addition, the above model does not capture certain flow constraints imposed by a link having a maximum capacity. These issues are dealt with in more detail in [3] and [4], as well as in Appendices A and B of this report.

Figure 1 illustrates the accuracy of the fluid model in comparison to discrete event simulation. The average window size of a TCP session is shown over time for a network of six routers and several thousand TCP sessions. The number of sessions is reduced by half at $t = 30$ and doubled at $t = 60$. We observe that the fluid model agrees with the simulation.

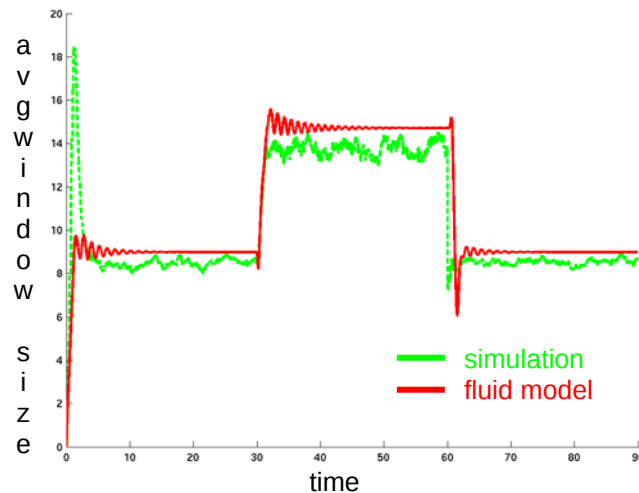


Figure 1: Comparison of discrete simulation and fluid model, average window size plotted as function of time

3.2 Steady State Models of TCP/IP Networks

We have described models that can be used to characterize the time varying behavior of large networks. In many cases an analyst is interested only in the steady state behavior of the network. In such cases, it is easier to characterize this behavior by solving a fixed-point problem. We briefly describe this approach in this section.

We consider a simple workload consisting of N infinite duration TCP flows labeled $i = 1 \dots, N$. Let $V_i = (j_{i,1}, j_{i,2}, \dots, j_{i,n(i)})$ be the ordered set of routers (i.e., route, path) taken by packets of flow i , where $j_{i,m} \in V$, $m = 1, \dots, n(i)$ and $n(i)$ is the number of links on the route. From the perspective of a link $v \in V$, it is useful to introduce S_v to be the set of TCP flows that traverse v . In addition, let $V_i(u) = (j_{i,1}, \dots, u)$ be the portion of the path from the data source to router $u \in V_i$ inclusive.

We (and others) have observed through measurements on the Internet and in numerous simulations that

- in the absence of a maximum rate constraint, each TCP flow traverses at least one congested router (here a congested router is one in which a flow suffers packet loss);
- each congested router is nearly fully utilized;
- each TCP flow exhibits a throughput that can be expressed as a function of the packet loss rate and average round trip time that it incurs. Denote this by $T(q, R)$ where q denotes the end-end loss rate and R the round trip time. We have derived accurate expressions for $T()$ in earlier work [5].

Consider the simple example where the N TCP flows share a single congested router with bandwidth C and buffer size B . Let us assume that this router includes the RED active queue management policy with a discard function such as that in equation (1). Let x denote the long term average queue length of this congested router. We denote the probability discard function as $p(x)$. We approximate the average round trip time of the i -th flow as

$$R_i(x) = A_i + x/C \quad (6)$$

where A_i denotes the sum of the propagation delay and transmission time at all routers on i 's path. The second term captures the contribution of congestion delay at the congested router to the round trip time. The assumption that the congested router is fully utilized (i.e., operates at full throughput) yields

$$\sum_{i=1}^N T_i(p(x), A_i + x/C)(1 - p(x)) = C, \quad (7)$$

which can be solved to obtain the average queue length of the congested router and, subsequently, the round trip time and the loss rate. This yields the steady state behavior of the system. We note that there exists a unique solution for x in the range $[0, B]$.

Before leaving this example, we comment on the generality of our approach. First, the N flows need not be TCP flows. It is sufficient that the i -th flow be characterized by a throughput function, $T()$, of the form postulated above. Most congestion control algorithms that have been proposed for next generation unicast and multicast transport protocols exhibit such throughput functions. These functions can be determined analytically, through simulation, or through measurement. Second, the approach applies equally as well if the active queue management policy marks packets rather than drops them. Last, although described in the context of a best effort service class, the basic approach is easily extended to the case of proposed Diff-Serv mechanisms. Consequently, the expression derived in [6] for the throughput of a flow operating in a Diff-Serv environment can similarly be used in our model. Finally, we note that constant bit rate, “non-responsive” flows can be accounted for in this model by simply subtracting their aggregate flow rate from the link capacity.

The above single congested router scenario is easily generalized to an arbitrary network supporting a workload of infinite duration responsive flows. Let $\mathbf{x} = \{x_v\}_{v \in V}$. Generalizing (6), we approximate the average round trip time of session i by:

$$R_i(\mathbf{x}) = A_i + \sum_{v \in V_i} x_v / C_v \quad (8)$$

Similarly, let $q_i(\mathbf{x})$ denote the probability that a flow i packet is lost on its end-to-end path; this quantity is given by

$$q_i(\mathbf{x}) = 1 - \prod_{v \in V_i(v)} (1 - p_v(x_v)) \quad (9)$$

Note that when T_i is a function of the probability of packet loss and the average round trip time (as postulated earlier), the above relations allow us to express T_i as a function of the average queue lengths. Henceforth we write the flow throughput as $T_i(\mathbf{x})$.

Let $S \subseteq V$ denote the set of congested routers, i.e., those routers whose bandwidths are fully utilized. We have the following set of equations, one for each congested router,

$$\sum_{i \in S_v} T_i(\mathbf{x}) \prod_{u \in V_i(v)} (1 - p_u(x_u)) = C_v, \quad v \in S$$

In addition, for the routers that are not congested, we have

$$x_v = 0, \quad v \notin S$$

Thus we have $|S|$ nonlinear equations with $|S|$ unknowns. Solving these equations yields a set of $\{x_k^*\}$ that can be used in turn to obtain the throughputs of all flows, their average round trip times and loss rates. Generalizations of our approach to responsive flows, other than TCP and to classes of service other than best effort, are identical to those described above for the single bottleneck router. Additional details can be found in [6].

We present two examples to illustrate this approach, one with a single congested router, and the other with two congested routers.

A single bottleneck router. We have N flows traversing a common router. We are interested in the accuracy of the model as N grows. Hence, we take $C = 3N/40$ Mbs. The probability discard function is taken to be that corresponding to a RED mechanism with parameters $p_0^{max} = 0.1$, $t_0^{min} = N/2$, $t_0^{max} = 3N/2$, and $B_0 = 3N$. The constant term in the average round trip delay for flow i is taken to be $(6+2i)$ ms, $i=1,\dots,N$. Thus, the round trip times range from 8ms to $(2N+8)$ ms.

We use the following expression for the throughput of a TCP flow taken from [5]:

$$T(q, R) = M \frac{(1-q)/q + W(q)/2 + Q(q, W(q))}{R(W(q) + 1) + Q(q, W(q))F(q)T_0/(1-q)}$$

where M is the packet size measured in bits and

$$W(q) = 2/3 + 2\sqrt{(1-q)/(3q) + 1/9},$$

$$Q(q, w) = \min\{1, (1 - (1-q)^3)(1 + (1-q)^3(1 - (1-q)^{w-3})) / (1 - (1-q)^w)\},$$

$$F(q) = 1 + q + 2q^2 + 4q^3 + 8q^4 + 16q^5 + 32q^6.$$

This has been shown to be accurate for a wide range of parameters [5]. Although this expression was derived for the case of a TCP flow using TCP-Reno, it has been shown to accurately predict the performance of other versions of TCP such as TCP-SACK.

We solve this model for different numbers of flows, $N=20, 40, 80, 160$ and compare its predictions with those obtained by simulating the system with ns [1]. The predicted average queue size and loss rate are compared to those obtained through simulation in Table 1. We observe that the loss rate predictions are accurate, within 8% of those obtained from the simulation. The predictions for the average queue lengths are not as accurate; they fall within 20% of those obtained through simulation. Note that this will result in errors in the average round trip times of 5%-20%.

No. flows	Avg. queue length		Loss rate	
	Model	sim.	Model	sim.
20	18.0	20.9	0.040	0.037
40	32.9	38.4	0.032	0.031
80	58.3	67.9	0.023	0.024
160	102.5	113.3	0.014	0.015

Table 1. Predicted vs. Simulated Average Queue Length and Loss Rate.

Figures 2 and 3 illustrate the individual flow throughputs as predicted by the model and as obtained from the simulation for $N = 80, 160$. We observe that the difference between analysis and simulation is always less than 12% and usually less than 5% in the case of $N = 80$. We also note that the difference decreases as the system size increases - a very desirable property. In the

case of $N = 160$, the worst individual flow error is less than 11%, and the errors are typically less than 3%.

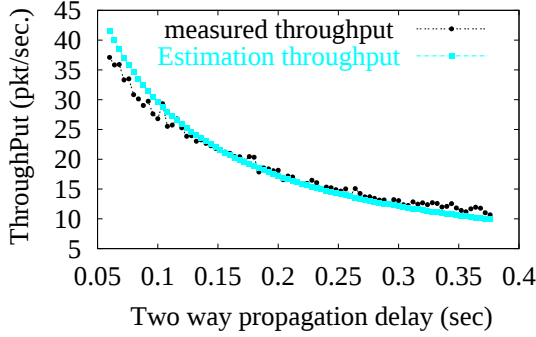


Figure 2: Flow throughputs, $N = 80$

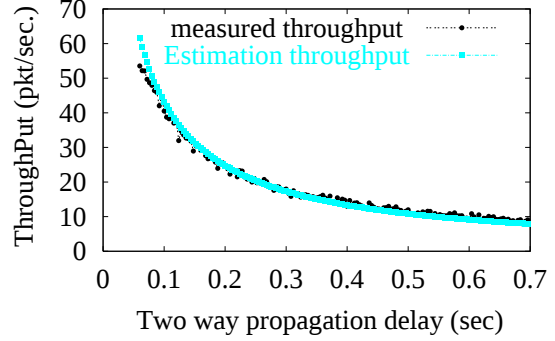


Figure 3: Flow throughputs, $N = 160$

Two congested links. The two-congested-router topology considered is shown in Figure 4. Routers 3 and 4 are configured so that they are not congested. There are three sets of flows in the simulation. The first set, consisting of 20 TCP flows, goes from router 3 to router 1. The second set of 40 TCP flows go from router 4 to router 2. The last set of 20 TCP flows go from router 0 to router 2. Router 0 is configured with a RED discard function with parameters $p_0^{max} = 0.1$, $t_0^{min} = 20$, $t_0^{max} = 60$, and a buffer capacity $B_0 = 120$. The capacity and propagation delay of router 0 is 3Mbps and 20ms, respectively. Router 1 is configured with a RED discard function with parameters $p_1^{max} = 0.1$, $t_1^{min} = 30$, $t_1^{max} = 90$, $B_1 = 180$. The capacity and propagation delay of router 1 are 4.5Mbps/sec and 20ms, respectively. The first two sets of TCP flows go through one congested router, while the third goes through both.

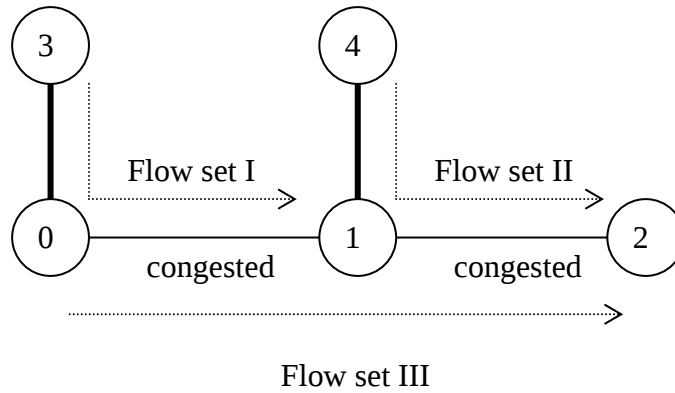


Figure 4: Two congested link example

Tables 2 and 3 show the average queue lengths and loss rates of each of the congested routers. Our preliminary investigations indicate that the differences between predicted performance and that obtained through simulation improve when we increase the number of flows while scaling up the RED parameters and the router bandwidth accordingly. Tables 4 and 5 show the average queue lengths and loss rates when all parameters are doubled, number of flows, capacities, t_0^{min} ,

t_l^{min} , t_l^{max} , and t_0^{max} . Once again, we see excellent agreement between analysis and simulation for loss and throughput measures, with good agreement for the average queue length.

Router	Avg. queue length		Loss rate	
	Model	sim.	Model	sim.
0	33.8	39.2	0.035	0.032
1	54.8	64.1	0.041	0.038

Table 2. Predicted vs. Simulated Average Queue Length and Loss Rate.

Flow set	Flow throughput	
	model	Sim.
1	28.6	27.4
2	23.9	23.2
3	8.9	9.7

Table 3. Predicted vs. Simulated Throughput.

Router	Avg. queue length		Loss rate	
	Model	sim.	Model	sim.
0	67.5	78.8	0.034	0.033
1	109.9	128.7	0.042	0.039

Table 4. Predicted vs. Simulated Average Queue Length and Loss Rate.

Flow set	Flow throughput	
	model	Sim.
1	27.2	28.7
2	23.2	23.7
3	9.9	8.8

Table 5. Predicted vs. Simulated Throughput.

4 Fluid Models for Worms and Failures

4.1 Worms

Computer “worms” are programs that self-propagate across a network exploiting security or policy flaws in widely-used services. In recent years, two major classes of worms, *scan-based worms* and *email worms*, have frequently attacked the Internet. Email worms propagate through emails and compromise computers when email users execute worm email attachments or simply view worm emails --- they require human interference to propagate and thus propagate relatively slowly. On the other hand, scan-based worms propagate by generating IP addresses to scan and directly compromise any vulnerable target computer --- they need no human activation and thus are able to propagate much faster than email worms. For example, Slammer in January 2003 infected more than 90% of vulnerable computers in the Internet within just 10 minutes. Recent well-known worms, Code Red, Code Red II, Slammer, Blaster, Sasser, are all scan-based worms. We concentrate on scan-based worms.

Attackers have tried many scanning strategies in previous worms. Code Red and Slammer uniformly scanned the entire IPv4 space. Code Red II used a local preference scan: it had a higher probability to scan an address within the same Class B or Class A network than a random address. Blaster sequentially scanned the Internet and chose its sequential-scan starting point from a local address with probability 0.4.

Based on classical epidemic spreading models, we have developed fluid models that can be used to study different scanning strategies such as those mentioned above. We present a simple model for a worm using a simple random scanning policy and then describe how it can be adapted to model other policies.

Consider a population of N vulnerable hosts, each of which has its own IP address. Let $I(t)$ denote the number of infectious hosts at time t . Then $[N - I(t)]$ is the number of susceptible hosts at time t . The epidemic model for a homogeneous system is:

$$dI(t)/dt = \beta I(t)[N - I(t)]$$

where β is the *pairwise rate of infection* in epidemiology studies from infectious hosts $I(t)$ to susceptible hosts $[N - I(t)]$. At $t = 0$, $I(0)$ hosts are infectious and the other $[N - I(0)]$ hosts are all susceptible.

Now β is given in terms of the scan rate of a computer, η , and the size of the address space, Ω , within which lie the vulnerable hosts. For example, in the case of Code Red, $N=360,000$, $\eta=358$ scans /min., and $\Omega=2^{32}$. Figure 5 shows how the number of infected hosts grows as a function of time. One observes that Code Red requires several hundred minutes to take off.

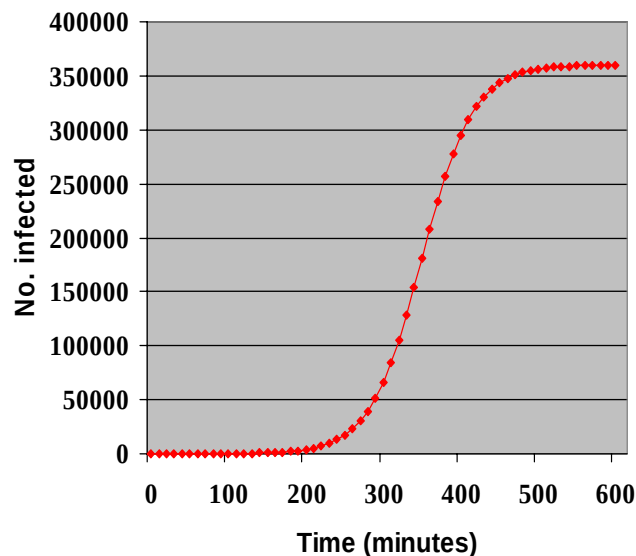


Figure 5: The classic Code Red worm

The model can be used to evaluate the performance of a number of different scanning strategies. These include:

- **Idealized worm:** All infected nodes are perfectly synchronized scanning only vulnerable hosts and only once.
- **Flash worm:** Infected worms know the identities of all vulnerable hosts and perform a random scan over them.

In both cases $\Omega = N$. The performances of such worms are illustrated in Figure 6 with and without a two second infection delay.

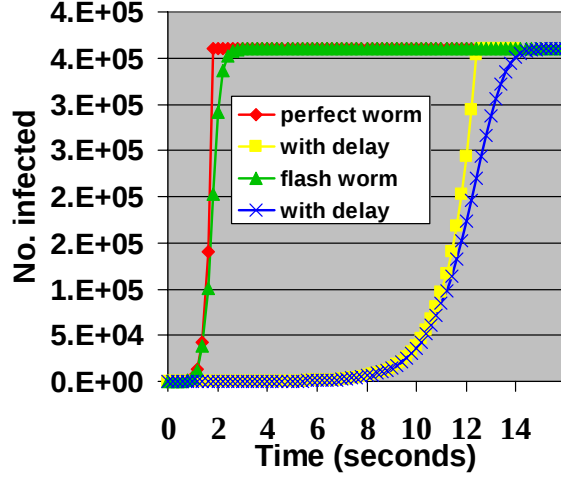


Figure 6: Idealized and flash worms with and without a two second delay

We have also adapted the model to study routing and local preference worms. Briefly, a routing worm is one that uses BGP routing tables to avoid portions of the IP address space that do not contain hosts. This has the effect of reducing the search space by 70%. A local preference worm scans addresses in the vicinity of the infected host with higher probability than distant addresses. Additional details of the models and its analysis are found in [8].

Figure 7 illustrates the performance of a routing worm and a local preference worm where the host scans addresses in its vicinity with probability 0.5.

From our analysis, we derive the following understanding of worm scanning strategies:

- Cooperation among infected hosts for scanning does not significantly increase a worm's spreading speed.
- A local preference scan increases a worm's propagation speed when vulnerable hosts are not uniformly distributed. The optimal local-scan probability increases as a function of the local-scan network size.
- When vulnerable hosts are uniformly distributed, divide-and-conquer scan, sequential scan, and uniform scan are equivalent in terms of a worm's propagation speed.

- A flash worm that uses uniform scan is an optimal spreading worm converged both from hit-list worm and from routing worm; a flash worm that makes sure no IP address is scanned more than once is the fastest spreading worm in terms of worm scanning strategy.
- For a sequential scan worm, such as Blaster, using local preference in selecting its starting point slows down the worm's propagation speed.
- For a selective attack worm, such as a routing worm [8], when the density of vulnerable hosts in the target domain is higher than the density in other domains, the worm propagates faster in the target domain if it only scans the target domain instead of all domains (and vice versa).

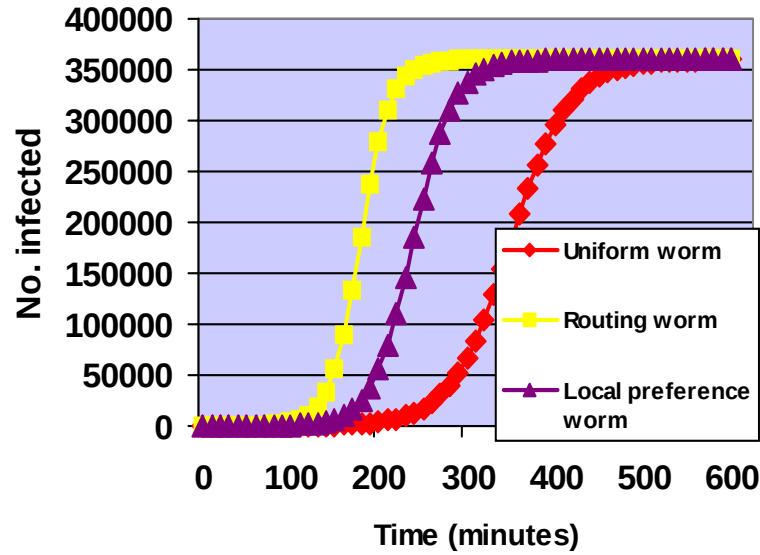


Figure 7: Comparison of routing and local preference worms with Code Red

In addition, based on our models, we have developed statistical detection algorithms for worms. Briefly, they look for an exponential increase in scan traffic that is characteristic of the spread of a worm. Additional details are found in [9], as well as in Appendix C of this report.

4.2 Cascading BGP failures

The Internet is a large collection of Autonomous Systems (AS) connected by the BGP routing infrastructure. The BGP maintains connectivity between AS's. It automatically reconfigures and computes routing tables when it detects a link failure. This computation starts locally around the failure point, and the information propagates through the Internet. The Internet routing table convergence has been observed to take tens of minutes. We developed models that reflect the behavior of BGP. Our analysis confirms that it is indeed possible to have cascading failures in the BGP routing infrastructure. Our results indicate the presence of phase transitions in these systems, and the presence and intensity of the phase transitions are strongly dependent on system parameters. We use the term *phase transitions* in the sense used by Erdős and Renyi in their work on random graphs that is “an abrupt change in a global system property”. The phase transition is to be interpreted as a sharp threshold rather than the definition used in statistical

mechanics. In our analysis, we observe that the propensity for phase transitions increases as clique size increases, and additionally also increases as the processing capacity of the routers decreases.

We consider a simple scenario with a finite set of N routers that are all connected to each other, i.e., form a clique. Let the number of down routers at any instant be $D(t)$. We define a “down” router to be one that does not have a functioning routing table, so a router that is in the process of rebooting and obtaining state dumps is also defined to be “down”. Now, we study the system of the number of down routers. The arrival and departure process to the system is defined as:

$$\alpha(t) := \text{number of arrivals in } [0; t)$$

$$\delta(t) := \text{number of departures in } [0; t)$$

Consider the process $\delta(t)$. The down routers come up with the help of the routers that are currently up ($N(t) - D(t)$). We define the service rate of an up router as k_s , where k_s is the average number of down routers a functioning router restores per unit time. Now, if $N(t) - D(t)$ servers are up and providing service, the service received by a single down node is its share of the total service capacity of the system. Thus, the share received by a single router is $(N(t) - D(t))/D(t)$. To account for the boundary condition $D = 0$, the denominator should actually be the term $(D(t) + k_a)$, where k_a represents the ambient load on the servers, representing, for instance, processing of normal route advertisements, and prevents $D \rightarrow -\infty$. However, for simplicity of exposition, we ignore the term as it does not affect the main observation we obtain later. Hence, the number of departures in an infinitesimal time dt , $d\delta(t)$, is defined by

$$d\delta(t) = D(t) (N(t) - D(t))/D(t) k_s dt = (N(t) - D(t)) k_s dt \quad (10)$$

Resetting BGP sessions leads to two kinds of messages, withdrawal announcements and subsequently full state updates when the BGP session is restored. We model the rate at which a functioning router goes down due to the load imposed by the resetting of a single BGP session as k . Typically, we expect $k_s \gg k$, as BGP resets are not uncommon and restoring a single session is unlikely to cause a peer router to go down. Now, the average arrivals in an infinitesimal time dt , $d\alpha(t)$, is given by the product of three quantities: the constant k , the number of routers (BGP sessions) that are currently down $D(t)$ (denoting the total load offered), and finally the number of routers currently up (that can go down), $(N(t) - D(t))$, i.e.,

$$d\alpha(t) = k D(t) (N(t) - D(t)) dt \quad (11)$$

Now $D(t)$ is $\alpha(t - \delta t)$, hence combining (9) and (10) yields after some manipulation the following relation

$$dD/dt = -k_i D^2 + (k_s + k_i N)D - k_s N$$

This is a Riccati equation, and without going into the actual solution of the equation, we immediately observe that the dynamical system described by this model exhibits a *phase transition*: If the initial state $D(0)$ of the system is above a certain threshold, then as $\lim_{t \rightarrow \infty} D(t) = N$, else $\lim_{t \rightarrow \infty} D(t) = 0$. In other words, if by some exogenous process (e.g. CODE RED) we

manage to bring a certain number of the routers down, thereby resetting the BGP sessions, then depending upon that number, the system either fully recovers or there is a *cascading failure*. A simulation of the system with different initial conditions is shown in Figure 8, where $k_s/k_i = 20$ and we plot two trajectories, one with $D(0) = 21$ and another with $D(0) = 19$. Additional details of this work can be found in [10], as well as in Appendix D of this report.

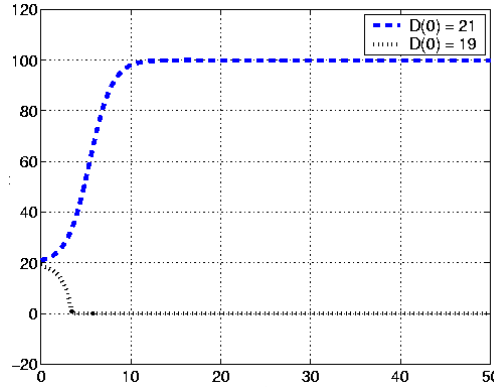


Figure 8: Illustration of how the initial number of failed routers can result in very different behavior

5 Congestion Controllers

In Section 3, we presented a model that describes the behavior of a set of TCP flows in a network of AQM-based routers. Their behavior is described by a set of differential equations, which is the language of control theory. For example, the behavior of a single TCP flow traversing a single bottleneck queue is illustrated in Figure 9.

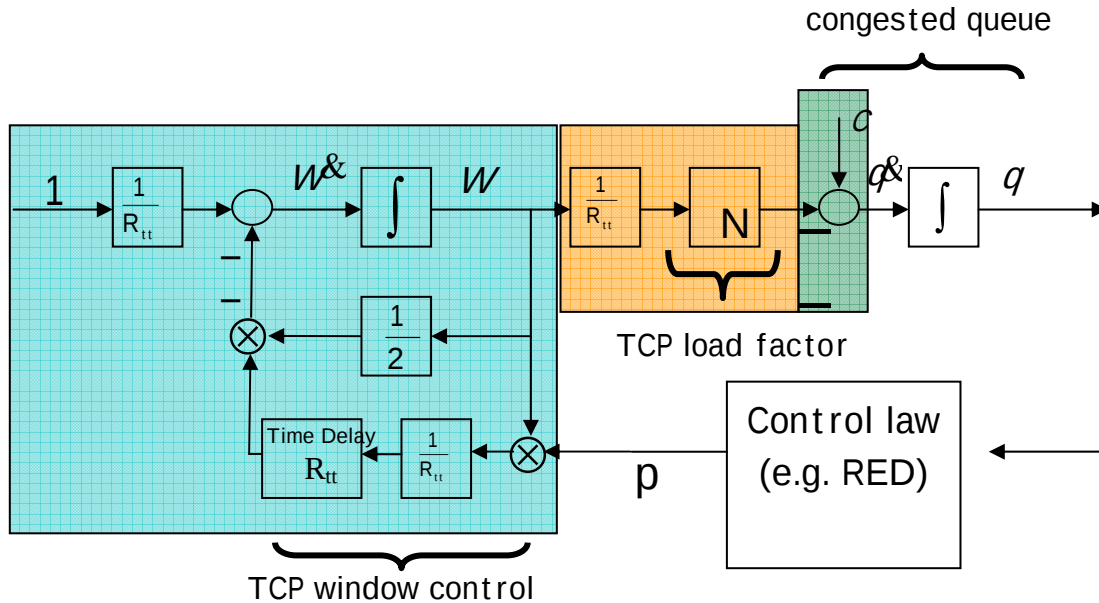


Figure 9: Control theoretic diagram of TCP

In the initial stages of the project, we performed a control theoretic analysis of the RED AQM protocol and presented guidelines for how to set its parameters. In the process of doing this, we uncovered a number of deficiencies of RED and proposed instead a new AQM algorithm based on proportional integral control. Two problems exhibited by RED are as follows:

- the queue averaging mechanism introduces additional delay in the TCP control loop which makes TCP unstable in many situations.
- RED couples the loss exhibited by a TCP session with the delay that its packets incur.

The new PI AQM controller introduces *phase lead*, which is ideal for systems with feedback delay. This leads to a speed of response faster than that of a stable RED mechanism. In addition, it decouples loss and delay; in particular, it is possible to specify a fixed queuing delay with a PI controller. Figure 10 compares the performance of a set of TCP sessions under RED and PI.

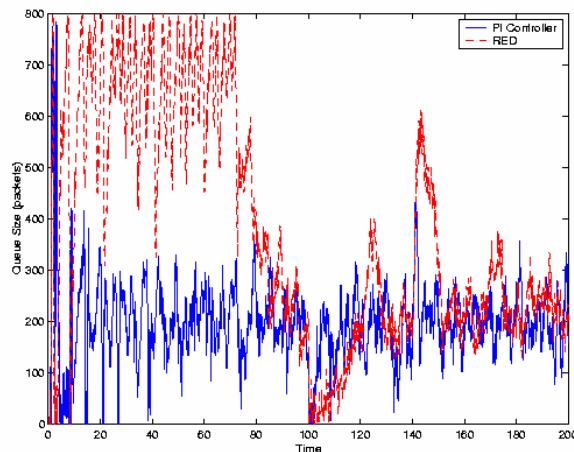


Figure 10: RED vs. PI control

Details of this work can be found in [11] which is found in the Appendix.

As part of the project, we also studied the problem of providing differential Quality of Service (QoS) to classes of TCP sessions. Specifically, we develop algorithms based on the IETF proposed DiffServ architecture [12], [13] to provide minimum throughput guarantees to different collections of TCP sessions. These algorithms rely on two components, multi-level PI AQM controllers residing at core routers that drop or mark packets differentially according to their colors and PI rate controllers residing at edge routers that monitor the throughput of these collections of TCP flows, marking packets as either red or green depending on whether or not the aggregate throughput exceeds its minimum rate. Details are found in [14], which is found in Appendix E.

Other results from this project include a study of the effect of uncontrollable flows on the behavior of TCP [15] and the development of algorithms to automatically tune AQM policies such as RED and PI [16].

6 Tools

As part of our project, we have implemented our fluid modeling methodology into the *ns* simulation software package. In particular, we developed a fluid network node that accepts packets as inputs, converts them into fluids, and then outputs packets after suffering appropriate losses and delays as predicted by the fluid model. This is illustrated in Figure 11. Additional details can be found in [17], as well as in Appendix F.

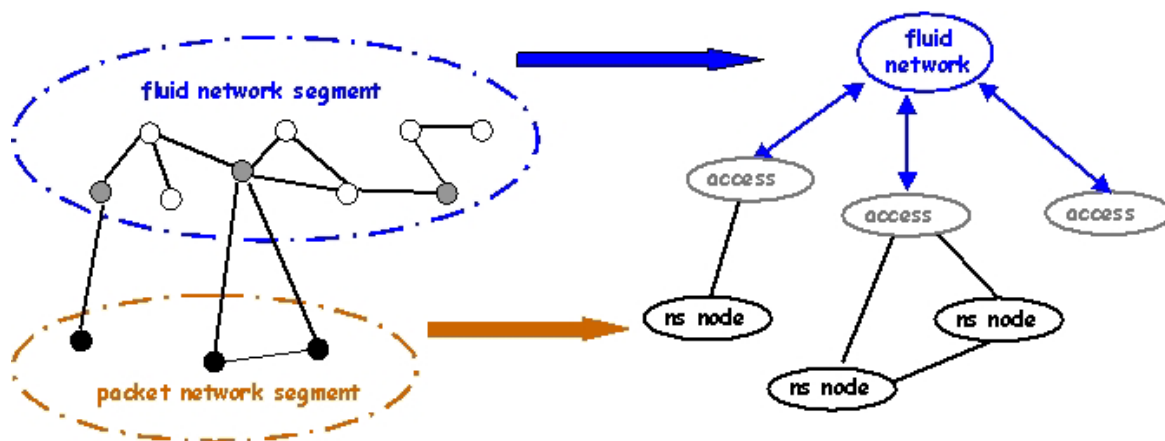


Figure 11: Fluid simulation module within *ns*

7 Summary

Our project significantly advanced the state-of-the-art in scalable modeling and simulation. We were the first to introduce the use of fluid models for the characterization of TCP flows within a large network. This is not standard practice. We were the first to begin to examine TCP behavior, coupled with AQM, using control theoretic principles. This is now standard practice and has resulted in significant advances in the development of better control mechanisms. Finally, we were the first to use fluid models to characterize worm spreading and cascading failures within the Internet. Again, this has become standard practice.

Finally, references [18] – [43] include additional results related to what has been discussed so far in the report. They are not included in the Appendix.

8 Acknowledgments

We would like to acknowledge a number of individuals who made contributions to the work reported here. These individuals are Tian Bu, Yossi Chait, Yu Gu, Jim Kurose, Francesco LoPresti, Yujing Wu, Honggang Zhang, Cliff Zou. We also would like to acknowledge Greg Hadynski for his help in preparing this final report.

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Appendix A: Fluid-based Analysis of a Network of AQM Routers Supporting TCP Flows with an Application to RED

A1. INTRODUCTION

Active queue management techniques were proposed in [8], [3] to both alleviate some congestion control problems for IP networks as well as provide some notion of quality of service. Modeling and analysis of such networks is important to understand their dynamics. While traditional discrete event simulations work well in general, even the most efficiently coded simulators suffer from the problem of scaling. We exploit fluid modeling to present a general methodology for the analysis of a network of routers that support active queue management with TCP flows. We use Poisson Counter Driven Stochastic Differential Equations to model the behavior of TCP traffic and derive a set of differential equations that describe the AQM policy and the router queuing process. Next, we develop a numerical scheme for obtaining the transient average behavior of a number of metrics including queue length, round trip time and TCP flow throughput from a set of coupled ordinary differential equations that result from our analysis. Given an AQM policy, we are able to get (expected) transient behavior of networks from our solution. We are able to handle large flows without a significant increase in computational complexity.

We illustrate our approach by analyzing RED [8] using the techniques developed in Section 3.1. Our solution technique yields predictions that match well with those obtained through simulation. Our modeling and solution methodology lead to a straightforward discovery of a critical problem with the RED averaging mechanism, one not addressed elsewhere. Our scheme has similar aims, to obtain transient behavior by numerical solution of a system of equations, to a heuristic approach proposed in [9]. It is however not clear how and why the heuristics work in cases reported.

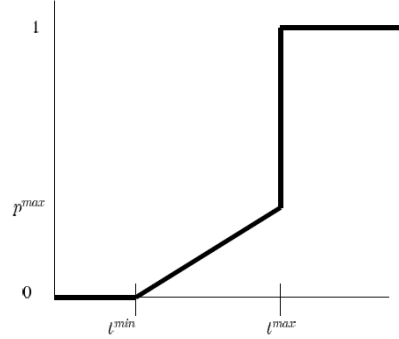
A2. APPLICATION TO THE RED ACTIVE QUEUE MANAGEMENT POLICY

We now present an application of the system, taking RED as the AQM policy, which has been shown to outperform Drop-Tail queues under certain scenarios. RED is a powerful mechanism to control traffic, potentially solving problems like flow synchronization, correlation of drop events while providing consistently high link utilizations. However, numerous problems have been cited with RED [10]. RED works well in certain scenarios, whereas it does very poorly, even worse than Drop-Tail, in other cases [6]. There is no clear understanding on how to tune various RED parameters that work well in all scenarios. Consequently, there is considerable nervousness in the community regarding deployment of RED, and numerous variations of RED have been proposed [5], [12], [4]. Some of the schemes have self-tuning parameters, while others maintain per-flow state. It is clear that there is a great need for thoroughly understanding the behavior of RED. We believe our techniques can help in that effort.

We first consider a system in which there is a single congested router with a transmission capacity of C . Associated with this router is an active queue management (AQM) policy that is characterized by a packet discard function $p(x)$ that takes as its argument an *estimate* of the average queue length at the router. The queue length of the router is denoted by $q(t)$, $t > 0$. The

classical example of an AQM policy is RED [8] for which $p(x)$ takes the form in the Figure below where t^{min} , t^{max} , and p^{max} are configurable parameters.

Figure 12: RED drop function



We will compare the results obtained from our model to those obtained from simulating an equivalent system using the well-known ns simulator. Our differential equation solver was implemented as a MATLAB program (a simple implementation with 42 total lines of code) which takes the matrix A , where the rows represent the different flows and the columns represent the different routers (queues) in a network, and a link capacity vector C as input. We did not incorporate slow-start in our program. All routers are assumed to have the same RED parameters. The propagation (non-queueing) delay for each class of flows is kept at 200 ms. The buffers are sized so that all losses are RED-related, i.e. no drop-tail losses occur on the network. The t^{min} is 150 packets and the t^{max} is 200 packets. δ is a parameter in our solver which is not specified in RED. Instead RED updates the queue size estimate on the arrival of each packet. We can account for it in two different ways

- ns updates the queue at every packet arrival. Thus, we choose δ_v to be $1/b_v$, where b_v is the instantaneous arrival rate at a queue measured in packets. δ_v thus becomes a function of time.
- We choose a fixed value of S_v . If the queue is stable, then the steady state arrival rate = service rate. Then, for each (bottlenecked) queue, S_v is simply $1/C_v$ where C_v is the capacity of link v in terms of (average) sized packets per second.

Note that both techniques are approximate. In our implementation we use $1/C_v$ as an estimate of δ_v .

A2.1. Experiment topology

We use the simple topology shown in the Figure below. It consists of two RED queues Q1 and Q2. Both links have three sets of flows going through them. S2 goes through both the queues, whereas S1, S3 and S4, S5 go through only Q1 and Q2 respectively. The only bottleneck links are the queues Q1 and Q2. We'll show results from 5 different experiments performed using ns and our DE solver.

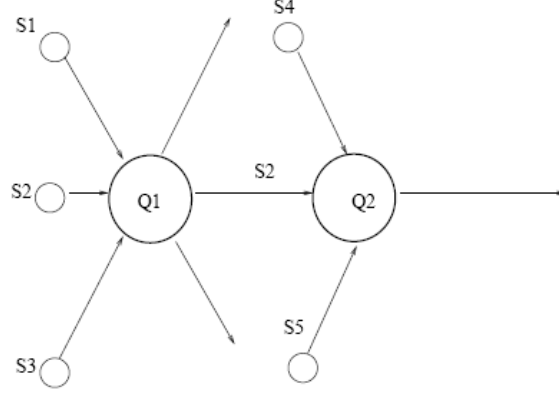


Figure 13: Simple network topology

We tabulate the parameter choices in the various experiments in the following table for reference.

Parameter sets for various experiments					
Exp. No.	α	Q1 cap.	Q2 cap.	Pkt size	p_{max}
1	0.0001	5 Mb/s	5 Mb/s	500 Bytes	0.1
2	0.0001	5 Mb/s	2.5 Mb/s	500 Bytes	0.1
3	0.0001	15 Mb/s	15 Mb/s	500 Bytes	0.1
4	0.0001	15 Mb/s	15 Mb/s	1500 Bytes	0.1
5	0.0001	15 Mb/s	15 Mb/s	500 Bytes	1

A2.2. Experiment 1

We first consider a symmetric case, where both RED queues have similar bandwidth capacity of 5Mb/s. α is kept at 0.0001, $p^{max} = 0.1$. Each class of flows consists of 40 individual flows which start at $t = 0$ (200 flows in all). At time $t = 75$, three fourths of the flows in each class drop out (so there are only 10 flows in each class). At time $t = 150$, those flows restart. We plot the Queue estimate and instantaneous queue length for Queue 1 and 2 in plots (a) and (b) of the following Figure respectively along with our model predictions of the expected values for the same. Note that we are plotting the results of *one* ns simulation along with our solution which gives the *expected* results. As we can see, the differential equation solution tracks the simulations pretty well, tracking both the average queue estimate and the instantaneous queue length well. Our method adapts to the changing nature of the load as well. Note that our solution differs initially from the simulation in all cases, because we didn't implement the effect of slow start.

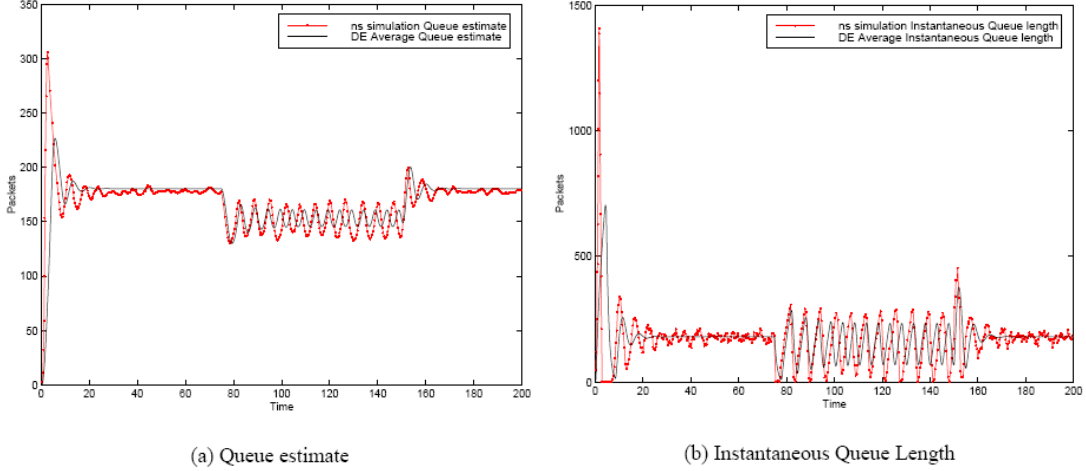


Figure 14: Symmetric case, plots for Queue 1, Experiment 1

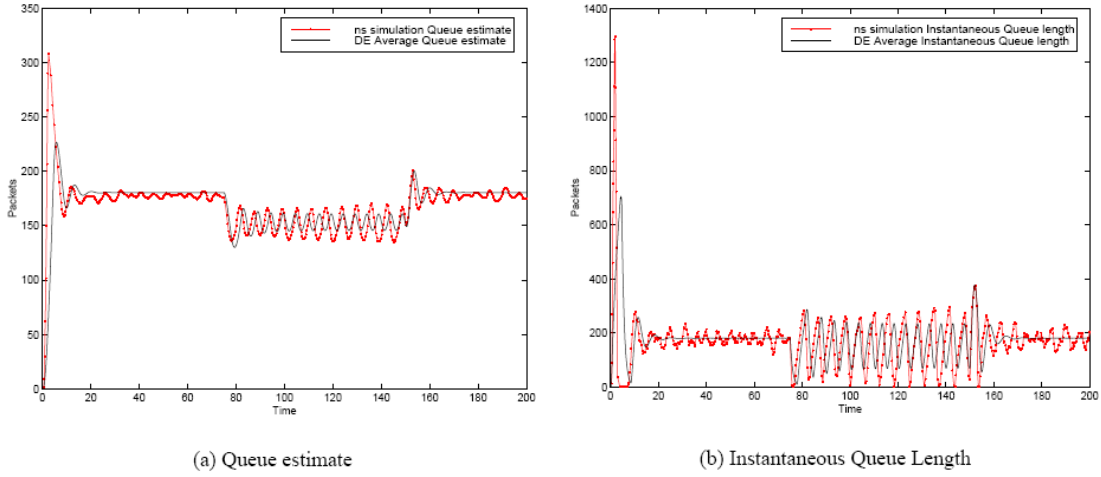


Figure 15: Symmetric case, plots for Queue 2, Experiment 1

A2.3. Experiment 2

Now we repeat the experiment in an asymmetric setting. We reduce the link capacity of the second queue to be 2.5Mb/s from 5Mb/s. Again we show the average queue estimate and instantaneous queue length along with our DE estimates for both the queues in the following Figures and our results match well with ns simulations for both the queues. Notice that the average queue estimate stays higher for Queue 2 which is correctly reflected in our DE solution.

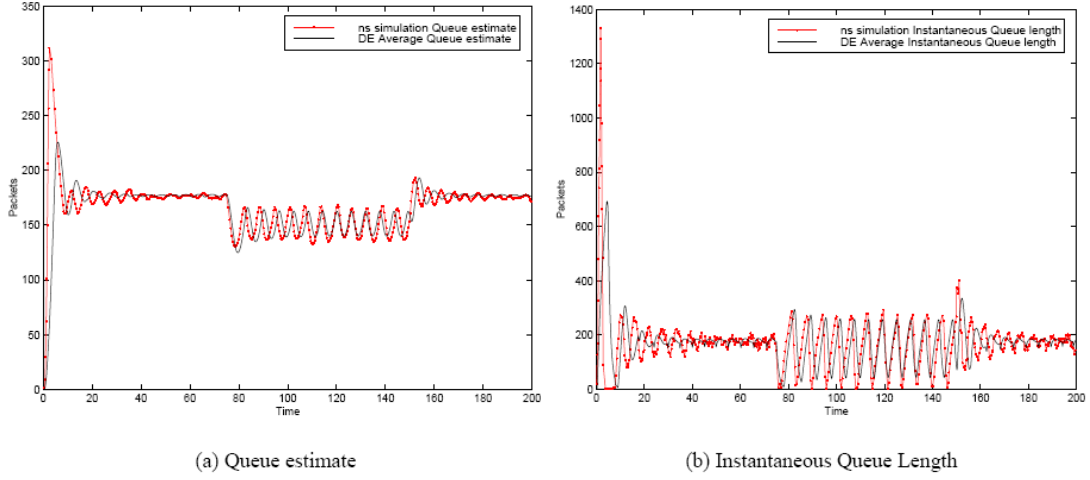


Figure 16: Asymmetric case, plots for Queue 1, Experiment 2

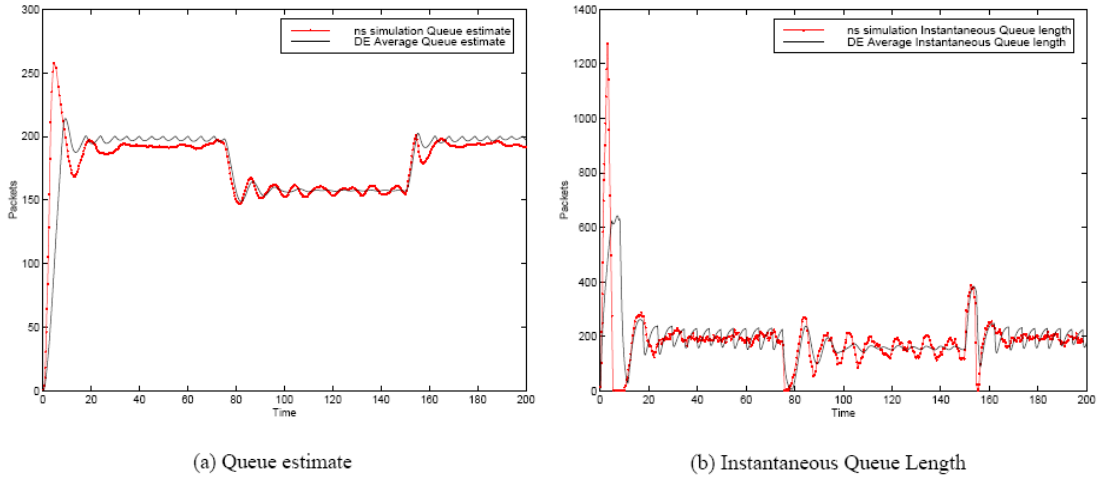


Figure 17: Asymmetric case, plots for Queue 2, Experiment 2

If we focus on the middle portion of the Figures above (when only a fourth of the flows are active), we observe that Queue 1 exhibits more oscillations than Queue 2. The oscillations are not good for the network as they may result in unacceptably large queue lengths and hence a large variability in delays for the flows going through. Even if the mean delay may turn out to be the same, these oscillations add considerable jitter to the delays. If the buffer is not large enough, then the effect of the oscillations will be to cause buffer overflows. They also cause periodically high RED loss rates and affect the throughput adversely. The question is: why? The larger bandwidth capacity of Queue 1 certainly plays a part, reducing stability margins, however we would like to point out another, hidden, cause. The queue averaging module is essentially a first order low pass filter. The original design goal of RED was to track the average queue size (low frequency signal), and to filter out bursts (high frequency signal). The input to this filter is the instantaneous queue length, and the output is the average queue estimate. Asymptotically, the frequency response of this filter is described by the magnitude Bode plot shown in the following Figure, where K is inversely proportional to the sampling period. It allows frequencies smaller than K to pass through, while damping inputs at a frequency higher than K . In simple terms, K

determines the responsiveness of the filter. The higher the value of K , the faster it will respond to a sudden change. If we maintain a high value of K , then the AQM function starts tracking the instantaneous queue length closely resulting in sustained oscillations.

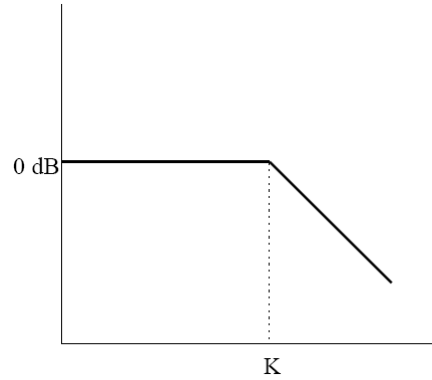


Figure 18: Magnitude Bode plot of first order averaging filter

A2.4. Experiment 3

In this experiment, all settings are unchanged except both link capacities are set at 15 Mb/s. The Figure below illustrates the Queue estimates and Instantaneous Queue Lengths for Queue 1 in the time interval [0 75] (i.e. when all flows are active).

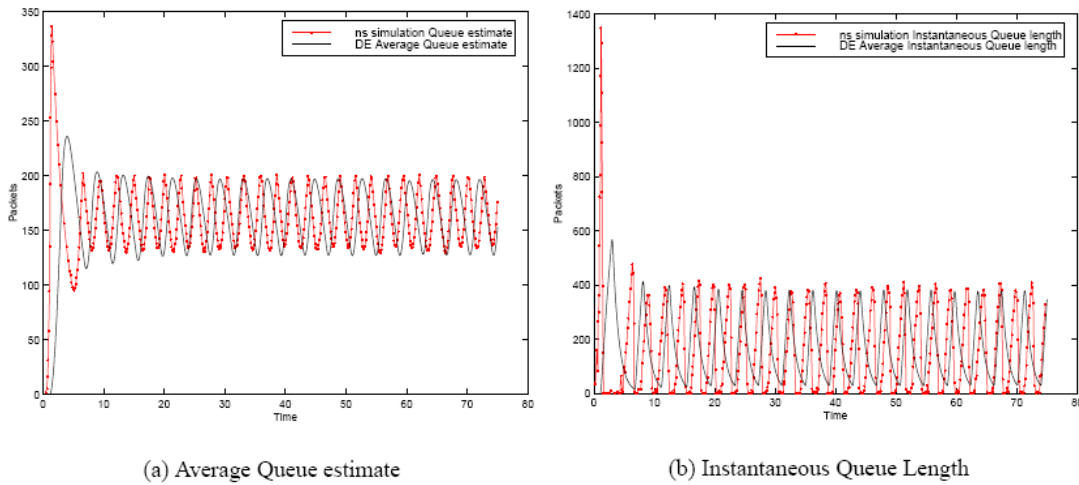


Figure 19: Link speed 15 Mb/s, packet size 500 bytes, Experiment 3

We observe the presence of large oscillations. In the scenario where the link capacity was 5Mb/s and the packet size 500 Bytes, the effective sampling period was 8×10^{-4} (the link capacity being 1250 packets of 500 bytes per second, and the sampling period being the inverse of the link capacity in packets per second). With the increase in link capacity to 15Mb/s, the value of the effective sample period is reduced to 2.66×10^{-4} and K becomes 3 times the earlier value. As the link capacity increases, the RED average queue estimate tracks the instantaneous queue length more closely, essentially resulting in sustained oscillations. This hidden artifact of the RED algorithm, the *adaptive* nature of the sampling period, is, in our opinion, a significant cause of

the "tuning problem" with RED. We can modify our DE solution system by using a static value for the sampling period, larger than what the link capacity would normally dictate, with the hope of reducing K and thereby reducing oscillations. The figure below illustrates the behavior of the Queue Estimate and Instantaneous Queue Length for different values of the sampling period δ . As we observe, increasing the value of δ results in increasing stability, with both the average queue estimate as well as average instantaneous queue length settling down. However, care should be taken that δ not be kept too large, as increasing values of δ result in increasing rise times of the average queue estimate and increasing initial overshoot of the average instantaneous queue length.

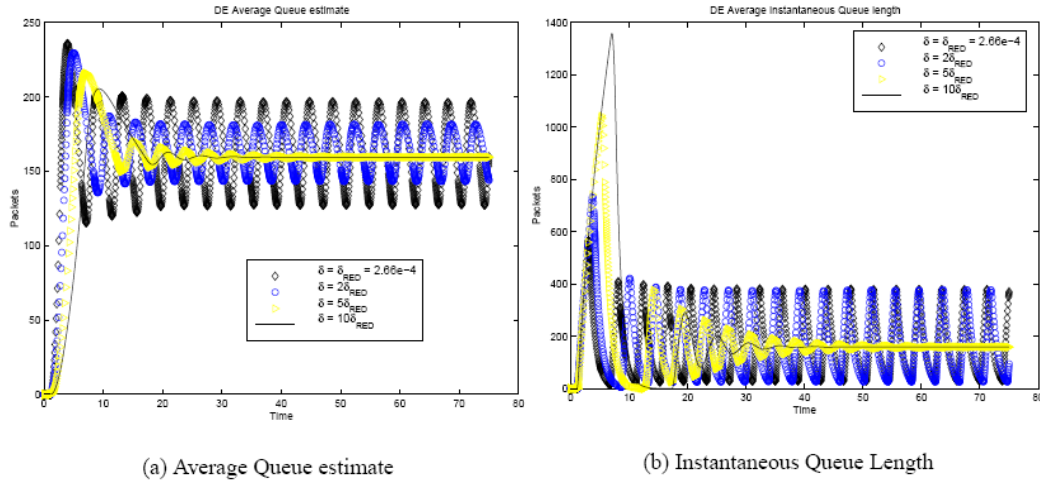


Figure 20: Behavior for different sampling periods, Experiment 3

A2.5. Experiment 4

Returning to the problem of tuning RED parameters, not only does the performance of the mechanism depend on link bandwidth, but also on the average packet size of the flows. We now illustrate this via experiment 4. Consider the two queue setting with the link capacities set to 15Mb. This time we increase the packet size from 500 Bytes to 1500 Bytes. This results in a δ that is approximately the same as the one where link capacity was 5 Mb/s with 500 Byte packets (the number of packets that are being processed remains a constant). Looking at the following Figure, we indeed observe that the system is stabilized.

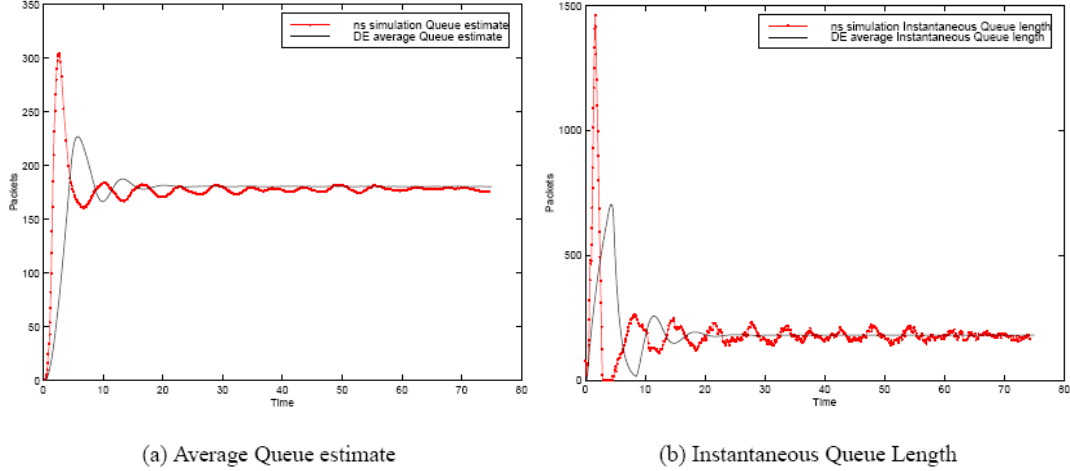


Figure 21: Link speed 15 Mb/s, packet size 1500 bytes, Experiment 4

Since the average packet size is not something network designers can control, the only hope to stabilize the RED algorithm is to make δ a parameter whose value is independent of link capacity and packet size. Implementing an algorithm whose stability is influenced by external factors (user packet sizes) is also not good from a security point of view. A malicious user could conceivably influence RED behavior by sending very small or very large packets.

In [6] Firoiu et. al. suggest that δ should be made equal to the smallest round trip time of the flows going through the link. They also suggested that the δ thus selected is "good enough" and finer sampling won't improve things. The authors propose some guidelines for choosing RED parameters. We are currently investigating "tuning" RED parameter values via a different, control theoretic viewpoint.

A2.6. Experiment 5

Further, in [6] the authors also suggested that the discontinuity in the RED drop function (the jump from p^{max} to 1 at t^{max}) is a cause of oscillations in RED. It has been suggested that making the drop function continuous via the *gentle*- [7] should improve things in that regard. Our experiments show that the discontinuity is *not* the *only* cause of oscillations, and that a simple fix by making the drop function continuous won't remove them. To illustrate that, we repeat the previous experiment with $p^{max} = 1$, thereby removing the discontinuity. We perform experiment 5, and the results are illustrated in the following Figure, showing that the oscillations *persist*.

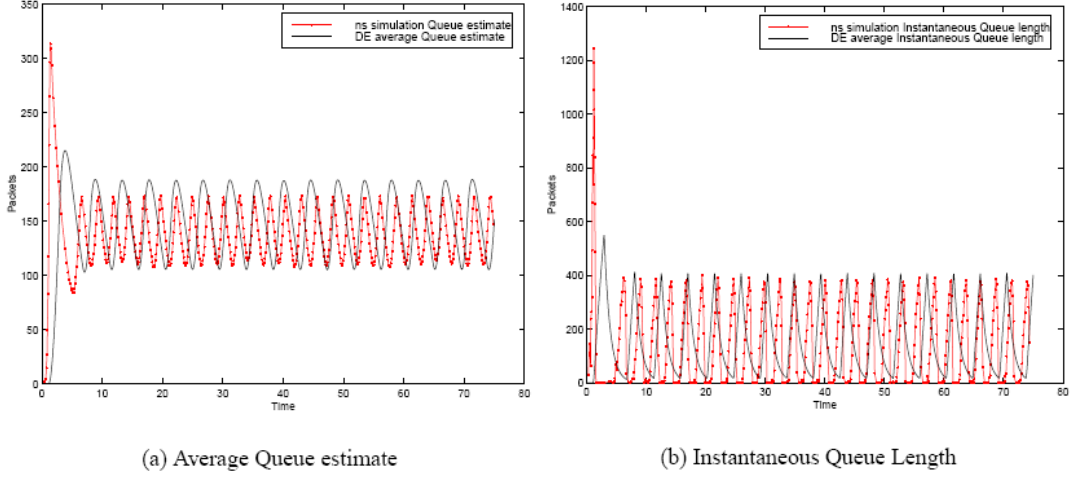


Figure 22: $p_{max}=1$, Experiment 5

A quick investigation with our differential equation tool reveals that the system can be stabilized with a carefully selected weighting function. However, on the downside, the system becomes very sluggish in its response time to changes in the load. Thus, this tradeoff between responsiveness and stability unfortunately cannot be avoided with the RED control mechanism. Thus, it is a combination of the link bandwidth, average packet size, weighting factor, sampling period, and load levels which make the system stable. Summarizing, our main observations with RED are

- The adaptive nature of the sampling interval is harmful and can lead to oscillations.
- The averaging algorithm needs to be modified, to make the sampling period a static value independent of packet sizes or arrival rates.
- The presence of oscillations depends on many factors including packet size, link bandwidth and load levels.

A3. CONCLUSIONS

We developed a methodology to model and obtain (numerically) expected transient behavior of networks with Active Queue Management routers supporting TCP flows. We applied our techniques to analyze networks where the AQM policy was RED. Our results match well with simulation results, and are able to scale up well to large flows. We are able to get a qualitative understanding of the behavior of such networks quickly with our tool. Our modeling technique enables us to spot a possible problem with the RED averaging mechanism, which we verify via simulations. This technique is quite general and can be easily extended to model and analyze other AQM mechanisms. A very important by-product of our formulation is that we can map the differential equation based TCP+AQM system into a classical control systems model. Standard techniques can be used to analyze various mechanisms and propose improvements to algorithms as well as guidelines for choosing parameters of the algorithms.

Appendix B: Scalable Fluid Models and Simulations for Large-Scale IP Networks

B1. INTRODUCTION

In this Appendix we show how scalable modeling can be used to speed up "simulations". The idea is to abstract the behavior of IP networks into analytical models. Solving the models numerically then yields performance metrics close to those of the original networks, thereby enabling an understanding of the key aspects of the performance of networks. The starting point is the differential equation based model described in Section 3.1. The differential equations represent the expected or mean behavior of the system. Recent results [1,6] indicate that with appropriate scaling the differential equations in fact represent the *sample path* behavior, rather than the expected behavior. Hence, our solutions gain in accuracy as the size of the network is increased, a somewhat surprising result. We solve the differential equations numerically, using the Runge-Kutta method, and our simulations show speedups of orders of magnitude compared to packet level discrete event simulators such as ns. Additionally, the time-stepped nature of our solution mode lends itself to a particularly simple parallelization. We also perform optimization to identify links that are not bottlenecks to speed up the simulations.

We also incorporate topological information in the set of differential equations in this Appendix. Our model developed in Section 3.1 of the report (henceforth referred to as the MGT00 model) defined a traffic matrix, which described the set of routers through which a particular set of flows traverses. However, the *order* in which the flows traverse the routers is ignored, and this information is potentially of critical importance. In Section B.2.2 we exemplify this with a pathological case wherein the model described in this report yields misleading results corrected by the refined model presented here. We also model the behavior of a number of variants of TCP such as SACK, Reno and New-Reno. Going beyond RED [2], we incorporate other modern AQM mechanisms such as AVQ [4] and PI control [3].

A related approach is that of [5]. In that approach, the authors exploit the model in Section 3.1 and the ideas of [1,6] to demonstrate that the behavior of the network is invariant if the flow population and the link capacities are scaled together. Their approach to simulating large populations of flows over high capacity links is to scale down the system to a smaller number of flows over smaller capacity links, thereby making the simulation tractable for discrete event simulators. The idea is appealing (and in fact applies to our approach too), however the technique only explores the scaling in the population and capacity of the links. Our methodology enables exploring a wider dynamic range of parameters. We can solve larger networks (links and routers) than is possible using discrete event simulators. Preliminary results indicate that the computational requirement of our method grows linearly with the size of the network, whereas the growth of the computational requirement of discrete event simulators is super-linear in the size of the network.

B2. FLUID MODELS OF IP NETWORKS

In this section we present a fluid model of a network of routers serving a population of TCP flows. We focus on persistent TCP connections working in congestion avoidance stage.

B2.1. Network Model

The network is modeled as a directed graph $G = (V, E)$ where V is a set of routers and E is a set of links. Each link l has a capacity of C_l bps. In addition, associated with each link is an AQM policy, characterized by a probability discarding/marketing function $p_l(t)$, which may depend on link state such as queue length. We develop models for AQMs with both marking and dropping. For the clarity of presentation, we focus on AQMs with packet dropping unless explicitly specified. The queue length of link l is $q_l(t)$, $t > 0$. Traffic propagation delay on link l is a_l .

The network G serves a population of N classes of TCP flows. We denote by n_i the number of flows in class i , $i = 1, \dots, N$. TCP flows within the same class have the same characteristics, follow the same route and experience the same round trip propagation delays.

B2.2. Deficiencies in the MGT00 Model

The second equation, (4) in Section 3.1, models the average queue length behavior as the accumulated difference between packet arrival rate at the queue, which is approximated by $\sum_{i \in N} n_i A_i(t)$, and the link capacity C_l . Observe that the approximation arises in replacing the aggregate arrival rate at the queue at time t with the aggregate sending rate of the TCP flows traversing that queue at t . These two quantities can significantly differ for two reasons: (1) flows are shaped as they traverse bottleneck queues; and (2) the arrival rate at time t at a queue is a function of the sending rate at a time $t - d$, where d is the sum of the propagation and queueing delays from the sender up to the queue. This delay varies from queue to queue and from flow class to flow class. An extreme example consists of one TCP class which traversing two identical RED queues with bandwidth C in tandem. The TCP traffic will be shaped at the first queue so that its peak arrival rate at the second queue is less than or equal, which equals the service rate of the second queue. Clearly, there won't be any congestion in the second queue. However, from (4), we will get identical equations for those two queues. Therefore the model predicts the same queue length and packet dropping probability for them as shown in the Figure below.

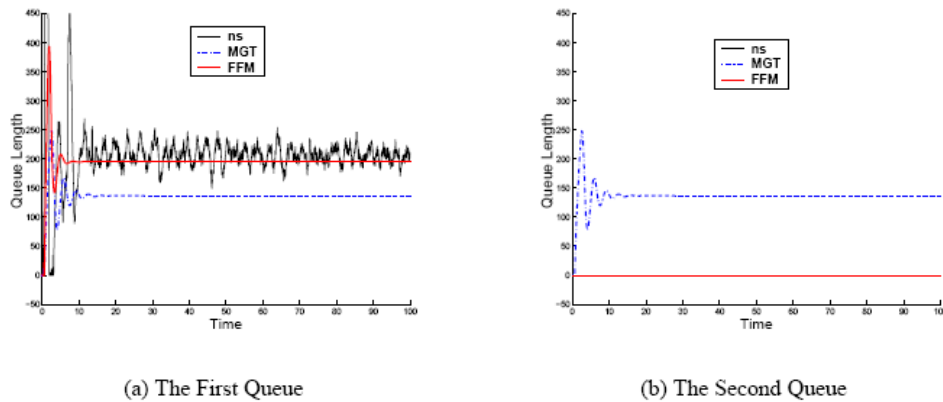


Figure 23: Importance of topology order

B2.3. A Topology Aware Model

We have observed in the previous subsection the importance of accounting for the order in which a TCP flow traverses the links on its path. In this section we present a model that takes into account how flows are shaped and delayed as they traverse the network. This is achieved by explicitly characterizing the arrival and departure rate of each class at each queue. The following expressions relate the departure and arrival process at a queue along the forward path.

—*Departure Rate* When the queue length, $q_l(t)$, is zero, the departure rate at time t equals the arrival rate. When the queue is not empty, the service capacity is divided among the competing flows in proportion to their arrival rates. Thus, for each flow of class i and l , we have

$$D_i^l(t) = \begin{cases} A_i^l(t), & q_l(t) = 0 \\ \frac{A_i^l(t-d_l)}{\sum_{j \in N_l} A_j^l(t-d_l)} C_l, & q_l(t) > 0 \end{cases}$$

where d_l is the queueing delay experienced by the traffic departing from l at time t . d_l can be obtained as the solution of the following equation

$$\frac{q_l(t-d_l)}{C_l} = d_l.$$

—*Arrival Rate* For each flow of class i , its arrival rate at the first queue on its route is just its sending rate. On any other queue, its arrival rate is its departure rate from its upstream queue after a time lag consisting of the link propagation delay. It is summarized in the following equation:

$$A_i^l(t) = \begin{cases} A_i(t), & l = k_{i,1} \\ D_i^{b_i(l)}(t - a_{b_i(l)}), & \text{otherwise} \end{cases}$$

The evolution of the system is governed by the following set of differential equations:

(1) *Window Size.*

The window size $W_i(t)$ of a flow of class i satisfies

$$\frac{dW_i(t)}{dt} = \frac{1(W_i(t) < M_i)}{R_i(t)} - \frac{W_i(t)}{2} \lambda_i(t)$$

where M_i is the maximal TCP window size, $R_i(t)$ and $\lambda_i(t)$ denote the round trip time and the loss indication rate at time t (in other words, as seen by the sender at time t).

(2) *Queue Length*

For each queue l , let N_l denote the set of TCP classes traversing it. Then:

$$\frac{dq_l(t)}{dt} = -1(q_l(t) > 0)C_l + \sum_{i \in N_l} n_i A_i^l(t)$$

We repeat the tandem queue experiment in Section B2.2. The revised topology order aware model gives accurate queue length results at both queues as shown in Figure 23.

B2.4. Model Reduction

In most operating networks, congestion only occurs at a small set of links, such as access points and peering points of ISPs. Most network links, especially in backbone networks, operate at low utilization levels. Queues at those links will be always empty and no packet will be dropped. Therefore there is no need to model queueing and RED behavior and maintain TCP states on those links. The network model can be reduced so that we only solve queueing and RED equations for potentially congested links and those *uncongested* links are transparent to all TCP classes except for introducing propagation delays. This can greatly reduce the computation time of the fluid model solver.

B3. EXPERIMENTAL RESULTS

We have performed extensive experiments to evaluate the accuracy and computation efficiency of our fluid models. We present several representative experiments here.

For all the experiments in this section, we use TCP Newreno and RED with ECN marking as the AQM policy. The TCP maximal window size is set to be 128. The step-size of the fluid model solver is fixed at 1 ms. We start with a single bottleneck topology time varying TCP workload. The fluid model's accuracy is tested by comparing its solution with simulation results obtained in ns when the network operates in both stable and unstable regions. In Section B3.2, the fluid model's scalability is demonstrated on a two bottleneck topology. The results show that the fluid model is scalable in the link bandwidth and flow populations. In addition, its accuracy improves as the link bandwidth scales up. In the last experiment, we test the capacity of our fluid model based simulation on a large topology with more than 1,000 nodes and thousands of TCP classes consisting up to 176,000 TCP flows. Computation results show that the fluid model approach is promising for simulating large IP networks.

B3.1. Accuracy of Fluid Model

The first experiment demonstrates the accuracy of our fluid model. As shown in the following Figure, there are 3 TCP classes. Each TCP class consists of 20 homogeneous TCP flows. The bottleneck link is between B_1 and B_2 . It has bandwidth of 10Mbps and propagation delay of 25ms. All other links in the network have bandwidth of 100Mbps and propagation delay of 20ms. There are a total of 14 queues. After model reduction, the fluid model only needs to simulate 4 queues which potentially have congestion. TCP classes 1 and 2 start at time 0. After 40 seconds, class 2 stops sending data. The number of TCP flows on the bottleneck link reduces from 40 to 20. The system enters an unstable region. At 70 second, both class 2 and class 3 become active. TCP workload increases by a factor of 3. The system eventually settles around a stable operation point.

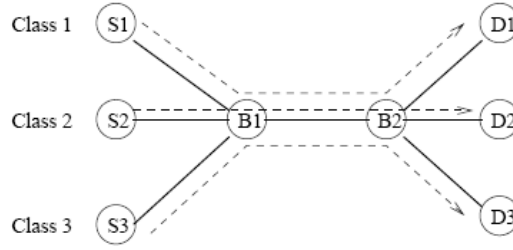


Figure 24: Single bottleneck network with dynamic workload

We compare the fluid model solution with results obtained from ns. Figure 25(a) plots one TCP connection's window sample path and the average window size we obtained from both ns and fluid model. The fluid model captures the average window behavior very well both when the system is stable and unstable. Figure 25(b) plots instantaneous queue length from ns and the average queue length predicted by the fluid model. We also observe a good match, which implies the fluid model calculates RED packet marking probability accurately.

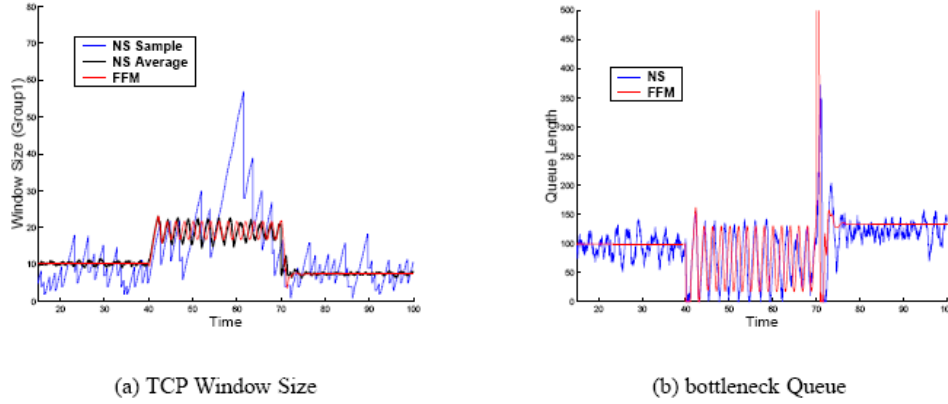


Figure 25: Results for single bottleneck topology

B3.2. Model Scalability with Link Bandwidth

The second set of experiments demonstrates the fluid model's scalability with link bandwidth and flow populations. We set up 3 TCP classes on a network of 8 links as in the Figure below.

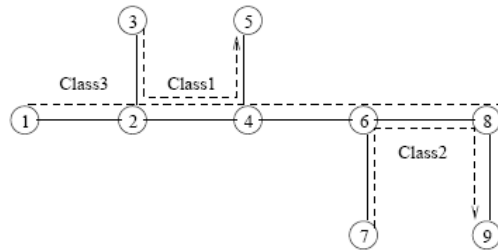


Figure 26: Network with two bottlenecks

Link bandwidth and flow population within each class are set to be proportional to a scale parameter K , which ranges from 1 to 100. The link between nodes 2 and 4, and the link between nodes 6

and 8 have bandwidths of $K * 10Mbps$. Other links have bandwidths of $K * 100Mbps$. Each TCP class consists of $K * 40$ TCP flows. In order for simulation results at different scales to be comparable, we make RED thresholds t^{min} and t^{max} proportional to K and its queue averaging weight α is inversely proportional to K . There are a total of 16 queues in the network. Our model reduction algorithm identifies 12 of them as *uncongested* that don't need to be simulated. For each K , we simulate the network for 100 seconds using both ns and the fluid model solver. The Figures below show simulation results for $K = 1$ and $K = 10$ respectively.

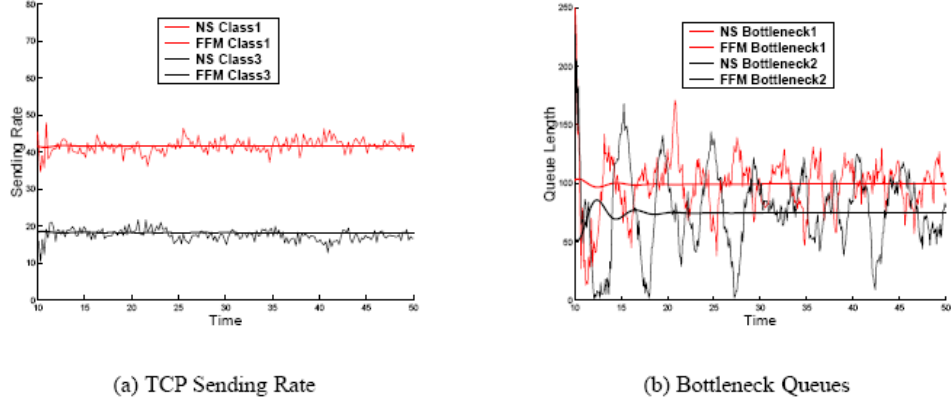


Figure 27: Simulation results when K=1

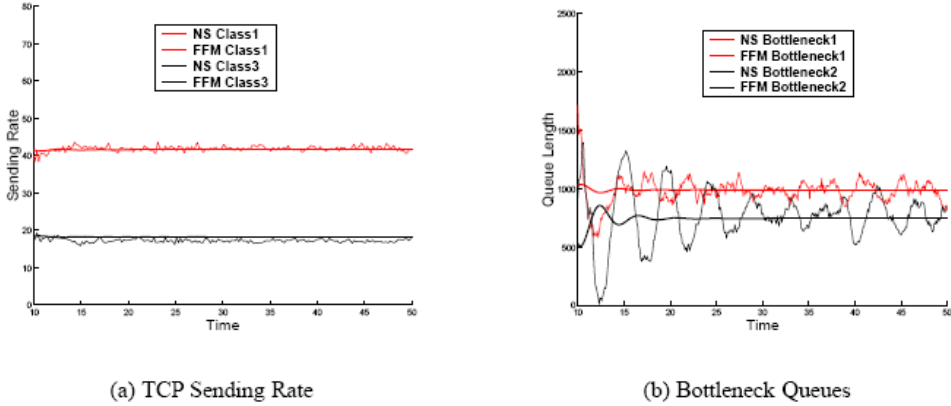


Figure 28: Simulation results when K=10

Tables I and II list simulation statistics of queue lengths and throughputs, including the mean obtained from both ns simulation and fluid model, the standard deviation of ns results and the absolute difference between ns and fluid model results, ns simulation results eventually converge to fluid model solution as K gets larger.

Table I. Sending Rate of Class1				
K	1	10	50	100
ns mean	42.0	41.6	42.5	41.8
ns std. dev.	0.59	0.28	0.766	0.21
FFM	41.6	41.6	41.6	41.6
abs. diff.	1.41	0.47	0.90	0.22

Table II. Queue Length at Bottleneck1

K	1	10	50	100
ns mean	100.4	995.4	4,875	9,942
ns std. dev.	18.7	59.5	100	162
FFM	99.1	990.5	4,953	9,905
abs. diff.	14.5	46.4	91.8	134

Because the fluid model is scalable with both link bandwidth and flow population, the computation cost to obtain model solution is invariant to the scale parameter K . On the other hand, the number of packets need to be processed in ns grow as link bandwidth and number of flows scale up. It takes much longer for ns simulation to finish when $K = 100$ than $K = 1$. Table III lists pure computation costs in unit of second of ns and the fluid model, both without dumping data. The larger the scale, the bigger the computation savings are for the fluid model.

Table III. Computation Cost of ns and Fluid Model

K	1	10	50	100
ns	12.5	122	983	1,676
FFM	0.766	0.766	0.766	0.766
Speedup	16.32	159.3	1,283	2,188

B3.3. Experience with Large IP Networks

In this experiment, we test our fluid model's capacity to simulate large networks. We use a structured network topology adapted from a baseline network model posed as a challenge to large network simulators by the DARPA Network Modeling and Simulation program.

At a high level, the topology can be visualized as a ring of N nodes. Each node in the ring represents a campus network and is connected to its two neighbors by links with bandwidth of 9.6 Gbps and random delays uniformly distributed in the range of 10-20ms. In addition, each node is connected to a randomly chosen node other than its neighbors through a chord link. The Figure below on the left illustrates a ring structure generated for $N = 20$.

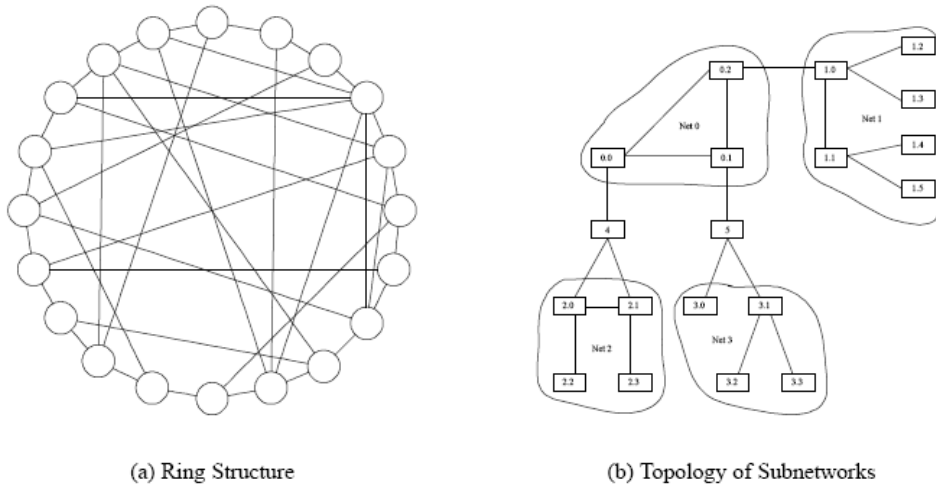


Figure 29: Topology of a large IP network

The campus networks at all nodes share the same topology shown in Figure 29. Each campus network consists of 4 sub networks: Net 0,1,2,3. All the links in campus networks have bandwidth of 2.5 Gbps.

Node 0.0 in Net 0 acts as the border router and connects to border routers of other campus networks. The links within Net 0 have random delays uniformly distributed in the range of 5 - 10ms. Links connecting 0.x to other sub-networks have random delays uniformly distributed in the range of 1 - 5 ms. All links in Net 1, Net 2 and Net 3 have random delays of 1 - 2 ms. Net 1 contains two routers and four application servers. Net 2 and Net 3 each contains four routers connecting to client hosts.

The traffic consists of persistent TCP flows. From each router in Net 2 and Net 3, there are 8 TCP classes. Four of them are destined to servers in its neighboring campus network. The other four classes are destined to servers in the campus network that connects to it through a chordal link. Each TCP class contains K homogeneous TCP flows.

In total, the entire network has $19N$ nodes, $44N$ queues, and $64N$ TCP classes. Our experiment is carried on a Dell Precision Workstation 530, which is configured with two Pentium IV processors (2.2 GHz) and 2 GB memory. However, as our program is not parallelized, only one processor is utilized. We fix the flow population of each TCP class at 50 and vary the number of campus networks on the ring from 5 to 55. Each topology is simulated for 100 seconds. Our model reduction algorithm identifies nearly 60% queues as *uncongested*. The Figure below illustrates simulation times that grow almost linearly with the number of campus networks. The simulation of the largest topology, which consists of 1,045 nodes and 176,000 TCP flows, completed after 74 minutes and 7.2 seconds when the step-size is set to be 1 ms. This simulation time can be reduced linearly when we use a larger step-size.

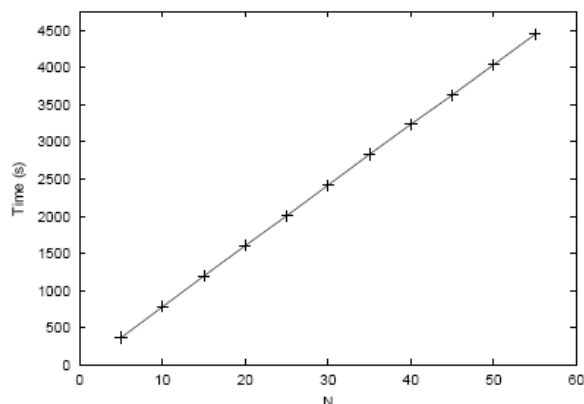


Figure 30: Computation cost as a function of N

B4. CONCLUSIONS AND FUTURE WORKS

In this appendix, we have developed a methodology to obtain performance metrics of large, high bandwidth IP networks. We started with the basic fluid model developed in Section 3.1 and made considerable improvements and enhancements to it. Most importantly, we made the model *topology aware*. That contribution alone is of independent interest in terms of theoretical (fluid)

studies of such networks, as topology awareness can play a critical part in conclusions regarding stability and performance as we demonstrated by a simple tandem queue example. Our solution methodology is computationally extremely efficient, and the scalable model enables us to obtain performance metrics of high bandwidth networks that are well beyond the capabilities of current discrete event simulators. Our technique also scales well with the size of the network, displaying a linear growth in computational complexity, as opposed to a super-linear one observed with discrete event simulators. The time stepped nature of our solution lends itself to a straightforward parallel implementation, pointing to another possible avenue of "simulating" large networks.

Appendix C: Monitoring and Early Detection of Internet Worms

C1. Introduction

Currently, some organizations and security companies, such as the CERT, CAIDA, and SANS Institute [7] [8] [32], are monitoring the Internet and paying close attention to any abnormal traffic. When they observe abnormal network activities, their security experts immediately analyze these incidents. Given the fast-spreading nature of Internet worms and their severe damage to our society, it is necessary to set up a nation-scale worm-monitoring and early-warning system. (The US Department of Homeland Security has launched a "Cybersecurity Monitoring Project" in October 2003 [40]).

A straightforward way to detect an unknown (zero-day) worm, is to use various anomaly detection systems. There are many well-studied methods or systems in the anomaly "intrusion detection" research area. For example, the "IDES" [13], "NIDES" [5] and "eBayes" [39] from SRI International; the anomaly intrusion detection method [15] based on "sequences of system calls"; the automatic model-construction intrusion detection system based on data-mining of audit data [24], etc.

Anomaly intrusion-detection systems usually concentrate on detecting attacks initiated by hackers. In the case of Internet worm detection, we find that we can take advantage of the difference between a worm's propagation and a hacker's intrusion attack. A worm code exhibits simple attack behaviors; all computers infected by a worm send out infection traffic that has similar statistical characteristics. Moreover, a worm's propagation in the Internet usually follows some dynamic models because of its large-scale distributed infection. On the other hand, a hacker's intrusion attack, which is more complicated, usually targets one or a set of specific computers and does not follow any well-defined dynamic model in most cases.

Based on this observation, we present a new detection methodology, "*trend detection*", by using the principle "detecting monitored traffic *trend*, not *burst*" [45]. Our "*trend detection*" system attempts to detect the dynamic *trend* of monitored traffic based on the fact that, at the early stage, a worm propagates exponentially with a *constant, positive* exponential rate. The "*trend*" we try to detect is the exponential growth trend of monitored traffic.

Based on worm propagation dynamic models, we detect the presence of a worm in its early propagation stage using the *Kalman filter* estimation algorithm, which is robust to background noise existing in the monitored data. The Kalman filter is activated when the monitoring system encounters a surge of illegitimate scan activities. If the infection rate estimated by the Kalman filter, which is also the exponential growth rate of a worm's propagation at its early stage, *stabilizes* and *oscillates* slightly around a *constant positive* value, we claim that the illegitimate scan activities are mainly caused by a worm, even if the estimated worm infection rate is still not well converged. If the monitored traffic is caused by non-worm noise, the traffic will not have the exponential growth trend, and the estimated value of the infection rate would converge to zero or oscillate around zero. In other words, the Kalman filter is used to detect the presence of a

worm by detecting the *trend*, not the *burst*, of the observed illegitimate traffic. In this way, the noisy illegitimate traffic in the Internet we observe everyday will not cause too many false alarms in our detection system.

In addition, we present a formula to predict a worm's vulnerable population size when the worm is still at its early propagation stage. We also present a formula to correct the bias in the number of infected hosts observed by a monitoring system. This bias has been mentioned in [10] and [29], but neither of them has presented methods to correct it. In this way, we can know how many computers in the global Internet are really infected based on local monitored data. Furthermore, we point out that in designing a worm monitoring system, the address space covered by a monitoring system should be as distributed as possible in order to monitor and detect non-uniform scan worms, especially a sequential scan worm such as Blaster.

C2. Related Work

In recent years, people have paid attention to the necessity of monitoring the Internet for malicious activities. Symantec Corp. has an "enterprise early warning solution" [1], which collects IDS and firewall attack data from the security systems of thousands of partners to keep track of the latest attack incidents. The SANS Institute set up the "Internet Storm Center" [17], which could gather the log data from participants' intrusion detection sensors distributed around the world.

In the academic research area, Moore *et al.* [29] presented the concept of "network telescope" to use a small fraction of unused IP space for observing security incidents in the global Internet. Pang *et al.* [30] called the abnormal traffic to unused IP space as "background radiation", and presented detailed measurement analysis and characterization of such monitored traffic. From another perspective, Berk *et al.* [6] proposed a monitoring system by collecting ICMP "Destination Unreachable" messages generated by routers for packets to unused IP addresses. In "honeypot" research, "Honeynet" [16] is a network of honeypots to gather comprehensive information of attacks; "Honeyd" presented by Provos [31] is a virtual honeypot framework to simulate many virtual computer systems at the network level.

The monitoring system we present in this appendix can be incorporated into the current monitoring systems such as the SANS "Internet Storm Center". Our contribution in this context is to point out the infrastructure specifically for worm monitoring, and what data should be collected for early detection of worms. We also emphasize the functionality of egress monitors, which has been overlooked in previous research. Worm monitors can be set up as ingress and egress filters on routers, which cover more IP space and gather more comprehensive information than the log data collected from intrusion detection sensors or firewalls for current monitoring systems.

In the area of worm modeling, Kephart, White and Chess of IBM performed a series of studies from 1991 to 1993 on viral infection based on epidemiology models [21] [20] [22]. Staniford *et al.* [37] used the classical epidemic model to model the spread of Code Red right after the Code Red incident on July 19th, 2001; they also proposed several more vicious worms in the same paper. Zou *et al.* [46] presented a "two-factor" worm model that considered both the effect of human countermeasures and the effect of the congestion caused by extensive worm

scan traffic. Chen *et al.* [10] presented a discrete-time version worm model that considered the patching and cleaning effect during a worm's propagation.

For a fast spreading worm such as Slammer, it is necessary to have automatic response and mitigation mechanisms. Moore *et al.* [28] discussed the effect of Internet quarantine for containing the propagation of a worm. Williamson [42] proposed a general rate-limiting "throttling" method to greatly constrain infection traffic sent out by infected hosts while not affecting normal traffic. Zou *et al.* [47] presented a feedback dynamic quarantine system for automatic mitigation by borrowing two principles used in the epidemic disease control in the real world: "preemptive quarantine" and "feedback adjustment." Staniford [36] presented automatic worm quarantine for enterprise networks by using "CouterMalice" devices to separate an enterprise network into many isolated subnetworks. Weaver *et al.* [41] further improved the CounterMalice quarantine by designing hardware-centered quarantine algorithms. Jung *et al.* [18] [19] proposed a "threshold random walk" algorithm to quickly detect and block worm scans based on the excessive illegal scans sent out by worm-infected hosts. "EarlyBird" in [35] and "Autograph" in [23] detect and block worm spreading through identifying the common characteristics, such as a common bit-string, among all infection network traffic of a worm. Wu *et al.* [43] proposed a victim counter-based detection algorithm that tracks the increased rate of new infected hosts.

Our early detection system tries to detect the presence of a worm in the global Internet. For worm detection in local networks, Staniford-Chen *et al.* [38] presented "GrIDS", which can detect worm-infected hosts in a local network through building the worm's infection graph (based on monitored infection traffic between all hosts); Dagon *et al.* [11] presented a "honeystat" worm detection method by correlating infection statistics provided by local honeypots when a worm tries to infect them. The CounterMalice quarantine device [36] also tries to detect infected hosts in local enterprise networks.

We assume that the IP infrastructure is the current IPv4. If IPv6 replaces IPv4, the vast IP space of the IPv6 would make it futile for a worm to propagate through blindly IP scanning [50]. However, we believe IPv6 will not replace IPv4 in the near future, and worms will continue to use various random scan techniques to spread in the Internet.

C3. Worm Propagation Model

A promising approach for modeling and evaluating the behavior of malware is the use of *fluid models*. Fluid models are appropriate for a system that consists of a large number of vulnerable hosts, which is the case for Internet-scale worm propagation modeling. In epidemiology research, the simple epidemic model [12] assumes that each host resides in one of two states: susceptible or infected. The model further assumes that once infected by a virus or a worm, a host remains in the infectious state forever. Thus any host has only one possible state transition: susceptible $\rightarrow$ infected. The simple epidemic model for a finite population is

$$\frac{dI_t}{dt} = \beta I_t [N - I_t], \quad (1)$$

where I_t is the number of infected hosts at time t ; N is the size of vulnerable population before any of them is infected; and β is called *the pairwise rate of infection* in epidemic studies [12]. At $t = 0$, I_0 hosts are initially infected while the remaining $N - I_0$ hosts are susceptible.

This model captures the basic mechanism of the propagation of a random-scan worm, especially for the initial stage of a worm's propagation when the effect of human counteractions and network congestion is ignorable [46]. A sequential-scan worm (such as Blaster), or a subnet-scan worm (such as Code Red II), propagates differently from a uniform-scan worm. However, through simulation and analysis, [48] showed that the propagation of these worms still closely follows the epidemic model (1).

The discrete-time version of the simple epidemic model (1) can be written as

$$I_t = (1 + \alpha\Delta)I_{t-1} - \beta\Delta I_{t-1}^2,$$

where

$$\alpha = \beta N.$$

We call α the *infection rate* because it is the average number of vulnerable hosts that can be infected per unit of time by one infected host during the early stage of a worm's propagation.

The epidemic model (1) has its limitations. First, the model assumes that all hosts can directly contact each other, which means it is not suitable for a topological worm (such as Morris [33]) or a mass-mailing email virus [49]. Second, if worm-infected hosts collaborate their infection efforts, such as the divide-and-conquer approach or the permutation scan used by the Warhol worm [37], then the worm's propagation will deviate from the epidemic model.

For the epidemic model (1), The following Figure shows the dynamics of I_t as time goes on for one set of parameters. We can roughly partition a worm's propagation into three phases: the slow start phase, the fast spread phase, and the slow finish phase. During the *slow start phase*, since $I_t \ll N$, the number of infected hosts increases exponentially (model (1) becomes $dI_t/dt \ll \beta N I_t$). After many hosts are infected and then participate in infecting others, the worm enters the *fast spread phase* where vulnerable hosts are infected at a fast, near linear speed. When most vulnerable computers have been infected, the worm enters the *slow finish phase* because the few leftover vulnerable computers are difficult for the worm to search out.

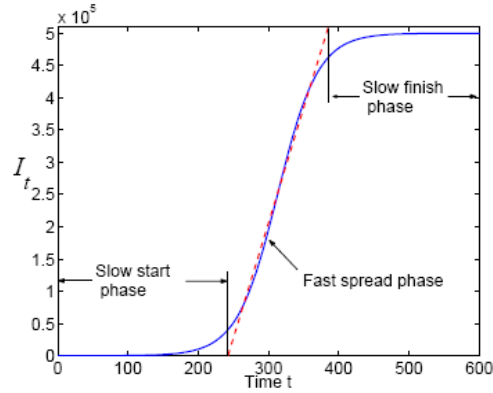


Figure 31: Worm propagation model

C4. Monitoring System

In this section, we propose the architecture of a worm monitoring system. The monitoring system aims to provide comprehensive observation data on a worm's activities for the early detection of the worm. The monitoring system consists of a *Malware Warning Center* (MWC) and distributed monitors as shown in the Figure below.

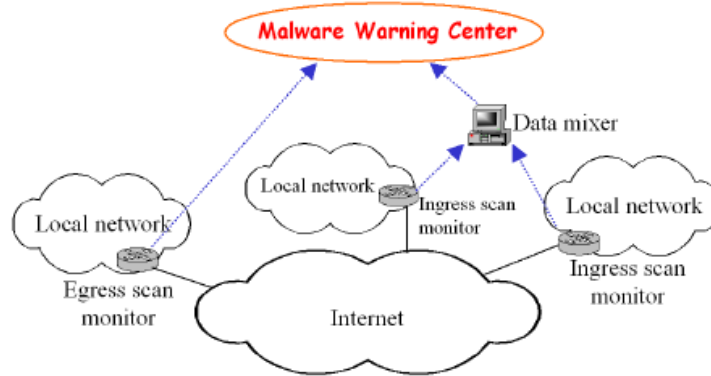


Figure 32: A generic worm monitoring system

A. Monitoring System Architecture

There are two types of monitors: ingress scan monitors and egress scan monitors. *Ingress scan monitors* are located on gateways or border routers of local networks. They can be the ingress filters on border routers of the local networks, or separated passive network monitors. The goal of an ingress scan monitor is to monitor scan traffic coming into a local network by logging incoming traffic to unused local IP addresses. For management reasons, local network administrators know how addresses inside their networks are allocated; it is relatively easy for them to set up the ingress scan monitor on routers in their local networks. For example, during the Code Red incident on July 19th 2002, a "/8" network at UCSD and two "/16" networks at Lawrence Berkeley Laboratory were used to collect Code Red scan traffic. All port 80 TCP SYN

packets coming in to nonexistent IP addresses in these networks were considered to be Code Red scans [27].

An *egress scan monitor* is located at the egress point of a local network. It can be set up as a part of the egress filter on the routers of a local network. The goal of an egress scan monitor is to monitor the outgoing traffic from a network to infer the scan behavior of a potential worm.

Ingress scan monitors listen to the global traffic in the Internet; they are sensors for global worm incidents (called "network telescope" in [29]). However, it is difficult to determine the behavior of each individual infected host from the data collected by ingress scan monitors because such monitors can only capture a small fraction of scans sent out by an infected host. On the other hand, if a computer inside a local network is infected, the egress scan monitor on this network's routers can observe most of the scans sent out by the compromised computer. Therefore, an egress scan monitor is good at observing a worm's scan rate and scan distribution, e.g., uniform scan (such as Code Red), or subnet scan (such as Code Red II and Sasser), or sequential scan (such as Blaster).

In order to provide early warning in real-time, distributed monitors are required to send observation data to the MWC continuously without significant delay, even when a worm's scan traffic has caused congestion to the Internet. For this reason, a tree-like hierarchy of *data mixers* can be set up between monitors and MWC: MWC is the root; the leaves of the tree are monitors. The monitors nearby a data mixer send observed data to the data mixer. After fusing the data together, the data mixer passes the data to a higher level data mixer or directly to MWC. An example of data fusion is the removal of repetitive IP addresses from the list of infected hosts. However, the tree structure of data mixers creates single points of failure, thus there is a trade-off in designing this hierarchical structure.

B. Location of Distributed Monitors

Ingress scan monitors on a local network may need to be placed on several routers instead of only on the border router because the border router may not know the usage of all IP addresses of this local network. In addition, since worms might choose different destination addresses by using different preferences, such as subnet scanning, we need to use distributed address spaces with different sizes and characteristics to ensure proper coverage. Later on, we show that for monitoring non-uniform scan worms such as Blaster, the IP space covered by a monitoring system should be as distributed as possible.

For egress scan monitors, worms on different infected computers may exhibit different scan behaviors. For example, Slammer's scan rate is constrained by an infected computer's bandwidth [26]. Therefore, we need to set up distributed egress filters to record the scan behaviors of many infected hosts at different locations and in different network environments. In this way, the monitoring system could obtain a comprehensive view of the behaviors of a worm. For example, it can get a better observation of the *average* number of scans an infected host sends out per unit of time.

C5. Data Collection and Bias Correction

After setting up a monitoring system, we need to determine what kind of data should be collected. The main task for an egress scan monitor is to determine the behaviors of a worm, such as the worm's average scan rate and scan distribution. Denote η as the "average worm scan rate", which is the *average* number of scans sent out by an infected host in a unit time. Thus in a monitoring interval Δ , an infected host sends out on average $\eta\Delta$ scans. The ingress scan monitors record two types of data: the number of scans they receive, and the source IP addresses of computers that send scans to them.

If all monitors send observation data to MWC once in every monitoring interval, then MWC obtains the following observation data at each discrete time epoch $t, t = 1, 2, \dots$:

- (1) The number of scans monitored in a monitoring interval from discrete time $(t - 1)$ to t , denoted by Z_t ,
- (2) The cumulative number of infected hosts observed by the discrete time t , denoted by C_t
- (3) A worm's scan distribution,
- (4) A worm's average scan rate η .

Let us first focus on worms that uniformly scan the Internet. Let p denote the probability that a worm's scan is monitored by a monitoring system. If ingress scan monitors cover m IP addresses, then a worm's scan has the probability $p = m/2^{32}$ to hit the monitoring system. We assume that in the discrete-time model all changes happen right before the discrete time epoch t , then we have

$$E[Z_t] = \eta\Delta p I_{t-1}$$

In order to detect non-uniform scan worms, it is important to observe a worm's scan distribution since it affects how we should use monitored data in our early detection. For example, if a subnet-scan worm has a higher preference in scanning local "/16" IP space, we can remove these "/16" local scans from monitored data Z_t in order to observe the worm's global scan trend. For a sequential scan worm, we can first apply a low-pass filter on monitored data Z_t to remove its excessive high-frequency noise before using the Kalman filter for early detection.

An egress scan monitor can observe the scan rates of all its internal infected hosts. If egress scan monitors cover many infected hosts, and if the scan rate of the worm does not vary too much, then we can obtain an accurate estimation of η , the worm's *average* scan rate. However, it is hard for the monitoring system to obtain an accurate estimate of η for a bandwidth-limited worm, such as Slammer or Witty, since the worm's scan rate could vary over several orders of magnitude [26] [34].

In [45] we derive the following estimate for the number of infected hosts at time, $\hat{I}_t$

$$\hat{I}_t = \frac{C_{t+1} - (1-p)^{\eta\Delta} C_t}{1 - (1-p)^{\eta\Delta}}$$

Expressions for the error are also given in [45]. The following Figure shows how well this estimator performs when the monitored IP address space is 2^{14} in the case of a Code Red style worm.

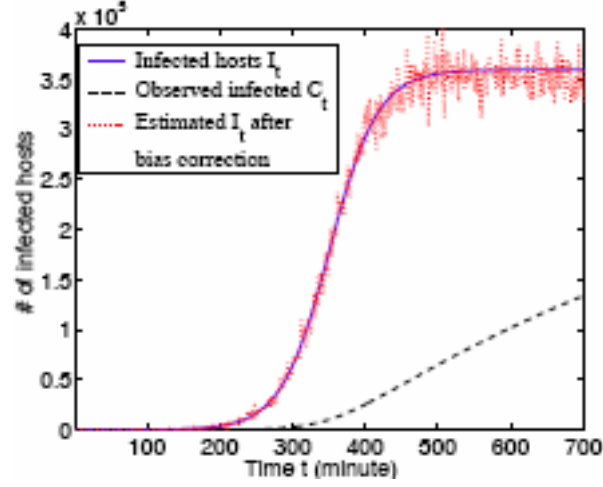


Figure 33. Estimate of number of infected hosts.

C6. Early Detection and Estimation of Worm Virulence

We present estimation methods based on recursive filtering algorithms (e.g., Kalman filters [4]) for stochastic dynamic systems. At MWC, we recursively estimate the parameter α based on observation data at each monitoring interval in order to detect a worm at its early propagation stage.

Let $y_1, y_2, \dots, y_t$, be the measurement data used by a Kalman filter estimation algorithm. Suppose the observations have one monitoring interval delay:

$$y_t = \delta I_{t-1} + w_t,$$

where w_t is the observation error. δ is a constant ratio: if we use Z_t as y_t , then $\delta = \eta \Delta p$; if we use $\hat{I}_t - 1$ derived from C_t by the bias correction, then $\delta = 1$. We use the following recurrence for I_t

$$\ln I_t = t \Delta \alpha + \ln I_0,$$

We take as the state of the Kalman filter $X_t = \begin{bmatrix} \Delta \alpha \\ K \end{bmatrix}$. Now $H_t = [t - t_0 \ 1]$ and the system is

described by

$$\begin{cases} X_t &= X_{t-1} \\ \ln y_t &= H_t X_t + v_t^n \end{cases}$$

The Kalman filter estimating X_t is

$$\begin{cases} H_t &= [t - t_0 \quad 1] \\ K_t &= P_{t-1}H_t^r / (H_t P_{t-1} H_t^r + 1/\tau_t) \\ P_t &= (I - K_t H_t) P_{t-1} \\ \hat{X}_t &= \hat{X}_{t-1} + K_t [\ln y_t - H_t X_{t-1}] \end{cases}$$

where τ_t is the weight of the t -th error term in the Least Square (LS) estimation algorithm [25]. We can use it to adjust whether our estimation should rely more on recently monitored data (τ_t increases as t increases) or equally on all monitored data (τ_t is a constant).

For a uniform-scan worm, we present an effective way to predict the population size N based on the observation data η and the estimate α . In this way, we can know how many computers are vulnerable in the Internet when a worm is still in its slow start phase. A uniform-scan worm sends out on average η scans per unit time; each scan has the probability $N/2^{32}$ to hit a host in the population under consideration. Hence, at the beginning when most hosts in the vulnerable population N are still vulnerable, a worm can infect on average $\eta N/2^{32}$ hosts per unit time. (The probability of two scans sent out by a single infected host hitting the same target is negligible). From the definition of infection rate α , we have $\alpha = \eta N/2^{32}$. Therefore, the population N is

$$N = \frac{2^{32}\alpha}{\eta},$$

where the average worm scan rate η is directly estimated from monitored data generated by egress scan monitors. When we use one of the Kalman filters above to estimate α , we can use the equation above to estimate N along with the Kalman filter estimation. In this way, the estimation of N has similar convergence properties to the estimation of α from the Kalman filter.

MWC collects and aggregates reports of worm scans from all distributed monitors once in every monitoring interval in real-time. For each TCP or UDP port, MWC has an alarm threshold for monitored illegitimate scan traffic Z_t . The observed number of scans Z_t , which contains non-worm noise, is below this threshold when there is no global spreading worm. This threshold can be chosen based on observations on normal days when no wide-spreading worm exists in the Internet. If the monitored scan traffic is over the alarm threshold for several consecutive monitoring intervals, e.g., Z_t is over the threshold for three consecutive times, a Kalman filter will be activated. Then MWC begins to record C_t and calculates the average worm scan rate η from the reports of egress scan monitors. Because C_t is a cumulative observation data that could cumulate all non-worm noise, MWC begins to record data C_t only after the Kalman filter is activated. The Kalman filter can either use C_t or Z_t to estimate all the parameters of a worm at discrete time t ($t = 1, 2, 3, \dots$).

The recursive estimation will continue until the estimated value of α shows a trend: if the estimate stabilizes and oscillates slightly around a *positive constant* value, we have detected the presence of a worm; if the estimate converges to or oscillates around zero, we believe the surge of illegitimate monitored traffic is caused by non-worm noise.

C7. Simulation Experiments

In this section, we describe the extensive simulations we used to study: (1) how a random-scan worm and a sequential-scan worm propagate; and (2) the performance of our Kalman filter-based early detection system. In addition, we show that the address space covered by the worm monitoring system should be as distributed as possible in order to better monitor non-uniform scan worms, especially a sequential-scan worm such as Blaster.

In our simulation experiments, we do not simply use the epidemic model (1) to numerically generate a worm's propagation curve. Instead, we have programmed discrete-time worm propagation simulators, which can be downloaded from [44], to simulate the detailed scanning behaviors of scans sent out by each infected host during each discrete time interval. In this way, we can accurately simulate the detailed propagation of a worm that uses any kind of scanning strategy.

A. Simulation Settings

We have simulated Code Red [2], SQL Slammer [26], and a sequential-scan worm similar to Blaster [3]. First, we explain how we choose the simulation parameters. In the case of Code Red, more than 359,000 Code Red infected hosts were observed on July 19th 2001 by CAIDA [27]. Thus in our simulation we set the Code Red vulnerable population $N = 360,000$. Staniford *et al.* [37] used a different format but the same epidemic model as (1) to model Code Red, where their model's parameter K is actually $K = \beta N = \alpha$ [46]. They determined that $K = 1.8$ for the time scale of one hour. Therefore, for the discrete time unit of one minute in our simulation, $\alpha = 1.8/60 = 0.03$. From (32) we can reversely derive $\eta = 2^{32}\alpha/N = 358$ per minute, i.e., Code Red sends out on average about 358 scans per minute per infected host.

Because different infected hosts have different scan rates, we assume that each infected host has a constant scan rate x , a rate that is independently predetermined by a normal distribution $N(\eta, \sigma^2)$, where $\sigma = 100$ (the scan rate x is bounded by $x \geq 1$). In our simulation, ingress scan monitors cover 2^{20} IP space. We also assume $I_0 = 10$ at the beginning.

Because of the sequential scan used by Blaster, people do not have a good estimation of how many computers were really infected by Blaster within the days following the worm's outbreak. We will explain the reason for this later in our experiments (shown in Fig. 10). In addition, there is no authoritative study of this worm's scan rate η . Therefore, in this appendix we simulate and study a sequential-scan worm that has the same "local preference" as Blaster [3], which is called a "Blaster-like" worm in this appendix. Since we want to understand how the sequential scan affects a worm's propagation and our early detection system, we give this Blaster-like worm the same parameters as Code Red in order to compare it with Code Red, i.e., we set the Blaster-like worm to have $N = 360,000$, $\eta = 358$ per minute. Each worm's scan rate x follows normal distribution $N(358, 100^2)$ with the bound $x \geq 1$, and $I_0 = 10$ at the beginning.

For a uniform-scan worm, such as Code Red, the distribution of vulnerable hosts in the Internet will not affect the worm's propagation. However, this distribution may affect the propagation of Blaster [48] because of its sequential scan. Since we do not know the true distribution of vulnerable hosts in the Internet, in our simulations of the Blaster-like worm, we assume

vulnerable hosts are uniformly distributed in the IP space defined by BGP routing prefixes, which is less than 30% of the entire IPv4 space [50].

We should choose an appropriate monitoring interval Δ in the Kalman filter estimation. Δ should not be too big in order to obtain enough sampling points in a worm's slow start phase for the Kalman filter. On the other hand, a too small monitoring interval puts more pressure on the MWC data collection, and introduces more monitoring statistical error w_t in (17) (because we can observe fewer worm scans in a smaller monitoring interval). In the discrete-time simulations in this appendix, the monitoring interval Δ is set to be one minute for Code Red and the Blaster-like worm. SQL Slammer propagates much faster and can finish infection in about 10 minutes [26]. Hence its monitoring interval should be much shorter in order to catch the dynamics of this worm. For this reason, the monitoring interval for Slammer is set to be one or several seconds.

B. Background Noise Consideration

We need to consider background non-worm noise in our simulations. Fortunately, Goldsmith [14] provided simple data of the background noise for Code Red activities monitored on a "/16" network (covers 2^{16} IP addresses). He recorded TCP port 80 SYN requests from Internet hosts to any unused IP addresses inside his local network. Such data are exactly the monitored data collected by ingress scan monitors in our proposed monitoring system. His monitored data showed that the background noise was small compared to Code Red traffic and the noise did not vary much. If we use normal distribution to model the background noise, then for each hour the number of noise scans follows $N(110.5, 30^2)$ and the number of source hosts that send noise follows $N(17.4, 3.3^2)$.

We try to hold the statistics of the observed background noise in our experiments: we monitor 2^{20} IP space, which is 16 times larger than what Goldsmith monitored, so the number of noise scans or noise sources should be enlarged by 16 times. We use one minute instead of one hour as the monitoring interval, thus we should decrease the number of noise scans or noise sources by 60. In this way, in our simulations of Code Red and the Blaster-like worm, the noise added into the observation data at each monitoring interval follows $N(29.5, 8^2)$ for Z_t and $N(4.63, 0.893^2)$ for C_t . Of course, this kind of extension of noise is very rough, but it is the best we can do based on the data available. Currently, we are trying to obtain detailed log data on previous worms from other researchers in order to have more realistic experiments.

In the simulation experiments, the alarm threshold for Z_t is set to be two times as large as the mean value of the background noise, i.e., the alarm threshold is $29.5 \times 2 = 59$. The Kalman filter we use in early detection will be activated when the monitored scan traffic Z_t is over the alarm threshold for three consecutive monitoring intervals. In this way, the Kalman filter will not be frequently activated by the surge of background noise traffic in the normal days.

C. Code Red Simulation and Early Detection

We simulate Code Red propagation for 100 simulation runs with the same input parameters but different seeds for random number generator. The following Figure shows the number of infected hosts as a function of time for three cases: the average value, the 95th percentile, and the 5th

percentile. The curve of 95th percentile means that in 95 out of our 100 simulation runs, Code Red propagates no faster than this curve represents.

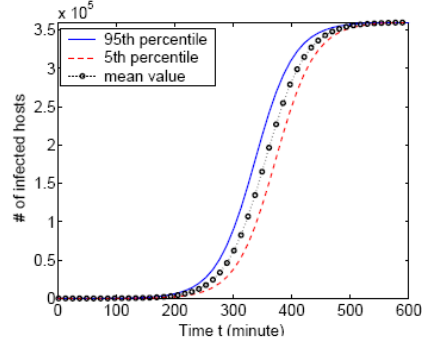


Figure 34: Code Red propagation and its variability (100 simulation runs)

This figure shows that a worm propagates slightly differently in different sample runs. The propagation speed difference is mainly caused by a worm's spreading at the beginning, when only several infected hosts scan and attempt to infect others. In fact, we have chosen $I_0 = 1$ and run Code Red propagation for another 100 simulation runs. It shows that Code Red in the $I_0 = 1$ case propagates more variously than the one shown in the figure above, where $I_0 = 10$.

For one simulation run of Code Red propagation, The Figure below shows the estimation of the worm infection rate α as a function of time by using three Kalman filters based on three discrete-time models: epidemic model, AR exponential model, and transformed linear model, respectively. This figure shows the estimates by using the processed monitored data Z_t after subtracting the average value of background noise from it. We can obtain the average value of noise based on the observations before activating the Kalman filter. We can use either the monitored data Z_t or the data C_t after bias correction to estimate α for Code Red. They provide the similar estimation results [45]. Later, when we study the early detection of the Blaster-like worm, because of its non-uniform scan, we cannot use the bias correction for the monitored data C_t and have to rely on the monitored data Z_t in our early detection. Therefore, in this appendix we will only discuss early detection by using the monitored data Z_t .

In this simulation run, Z_t at time 126, 127, and 128 minutes are over the alarm threshold 59, thus the Kalman filter is activated at time 128 minutes. The Figure below shows that the Kalman filter estimation based on the transformed linear model provides a much better estimation result than the other two because the noise v'_t introduced by the transformed linear model (7) is much smaller than the noise v_t and v_l introduced by the other two models.

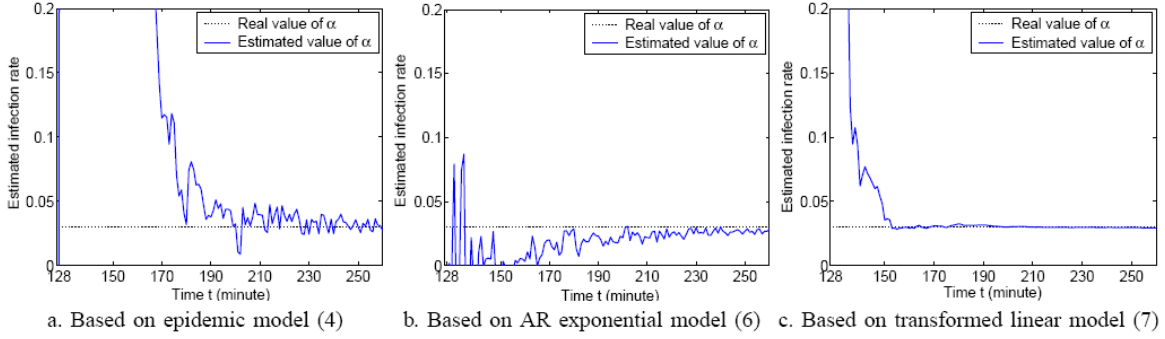


Figure 35: Kalman filter estimation of Code Red infection rate (for one simulation run)

In the Code Red simulation run shown in the Figure above, the worm infects 0.3% of vulnerable computers in the Internet at time 157 minutes. If we use the transformed linear model in our early detection, Plot (c) in the Figure above shows that the estimate of α has already stabilized at a positive constant value by that time. Therefore, we can detect the presence of Code Red when it has only infected 0.3% of all vulnerable population in the Internet. For the remaining 99 Code Red simulation runs, we have done such early detection by using Kalman filters and have achieved the similar early detection performance. In [45], we have shown that the early detection system can achieve a similar detection performance — detect a worm when it infects a similar fraction of the vulnerable population — no matter whether this worm propagates faster or slower in those 100 simulation runs.

The Kalman filter based on epidemic model is still useful since it is able to estimate worm infection rate α during the whole propagation period of a worm. On the other hand, because the transformed linear model is derived from the exponential-growth model (7), its Kalman filter will underestimate α when the worm enters its "fast spread phase". The Figure below shows the estimation results from these two Kalman filters before the worm infects 80% of vulnerable hosts at time 400 minutes. It shows that we should use these two Kalman filters together in the early detection of a worm.

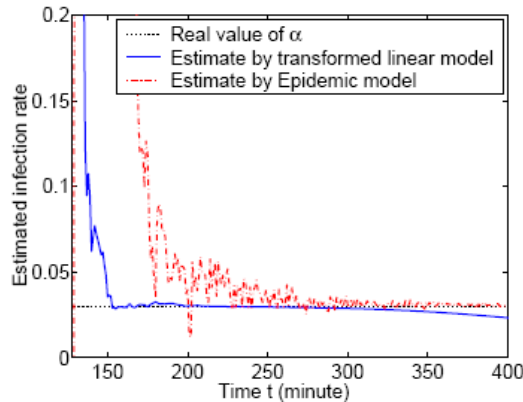


Figure 36: Long-term Kalman filter estimation

We predict the vulnerable population size N at each discrete time when we update the estimate of α from Kalman filters. The following Figure shows the estimated value of N as a function of

time based on the Kalman filters of transformed linear model and epidemic model, respectively. Because the estimate of N is proportional to the estimate of α , this figure has the same pattern as the Figure above. In a real implementation, we should combine both estimation curves shown in the Figure below to predict the vulnerable population size N .

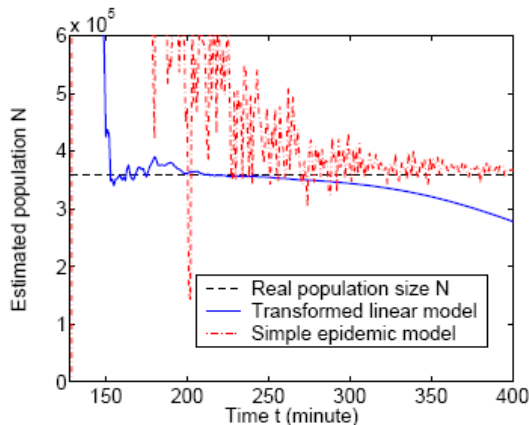


Figure 37: Estimate of the vulnerable population size N of Code Red

Because Slammer propagates in the same way as Code Red — by uniformly scanning the Internet — its propagation and its early detection are very similar to the methods used for detecting Code Red [45]. (We choose $\eta = 4000/\text{sec}$ as explained in [26] and $\eta = 1$ second). Therefore, we do not repeatedly show the early detection of Slammer here.

D. Blaster-like Worm Simulation and Early Detection

Each Blaster infected host scans the entire IP space sequentially from a selected starting point. To select this starting IP address, each worm copy has a 40% probability to choose the first address of its "Class C"-size subnet ($x.x.x.0$), and a 60% probability to choose a completely random IP address [3]. In our simulations, we let the Blaster-like worm to have the same local preference in selecting its starting point.

Since we select the same parameters for simulations of both Code Red and the Blaster-like worm, we can compare them to study how the sequential scan affects a worm's propagation. Again, we run the simulation of the Blaster-like worm for 100 simulation runs. The Figure below shows the 95th percentile and 5th percentile of the worm's propagation compared with the previous Code Red simulations shown earlier.

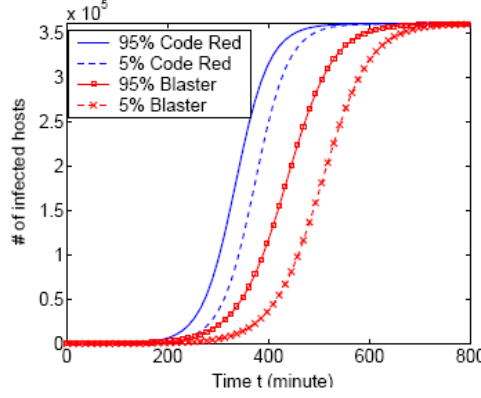


Figure 38: Worm propagation comparison between Code Red and Blaster-like worm

Even though the simple epidemic model (1) is derived based on uniform scanning, our simulation experiments show that the Blaster-like worm can still be accurately modeled by the simple epidemic model (1), and thus the worm can be modeled by the three discrete-time models presented in this appendix. This is consistent with a conclusion in [48], which shows that a sequential-scan worm has the same propagation dynamics as a uniform-scan worm when the vulnerable hosts are uniformly distributed.

However, we should keep in mind that a worm's propagation is in fact a stochastic process; the epidemic model (1) is accurate only when both the number of vulnerable hosts and the number of infected hosts are relatively large. For example, no ordinary differential equation models are suitable to model the very end of a worm's propagation when the worm finishes infecting the last several vulnerable hosts, which can only be modeled accurately by a stochastic model. Since we study an Internet-scale worm's propagation that involves hundreds of thousands or even millions of computers, the epidemic model (1) is a good abstract model for modeling a worm's dynamics except the very beginning and the very end of the worm's propagation.

The following Figure shows that the Blaster-like worm propagates slower than Code Red. Zou *et al.* [48] pointed out that this is because the Blaster-like worm selects its starting scanning point with a local preference, not because of its sequential scan mechanism.

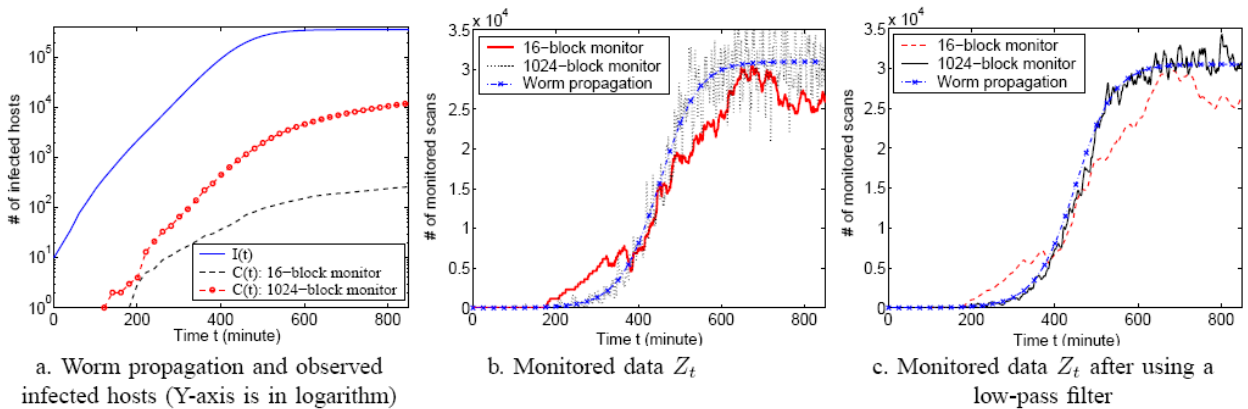


Figure 39: Blaster-like worm propagation and its monitored data

Because of its sequential scan, when monitoring the Blaster-like worm, we cannot let the monitoring system cover only one big block of IP address space — such a monitoring system can only observe a very small fraction of infected hosts in the Internet. For example, if a sequential scan worm has the same fast scan rate $\eta = 4000$ per second as Slammer [26], each infected host will take $2^{32}/\eta = 12.4$ days to finish scanning the entire IPv4 space. Therefore, most hosts infected by the Blaster-like worm will take days before their scans hit the big block IP address space monitored in such a monitoring system.

For this reason, a good worm monitoring system should cover as distributed as possible an IP address space in the Internet. We simulate two monitoring systems. Both monitoring systems cover the same 2^{20} IP addresses (the same as the monitoring system in previous Code Red study), but they consist of a different number of monitored IP blocks: one monitors 16 "/16" networks; the other monitors 1024 "/22" networks. All monitored address blocks in a monitoring system are evenly distributed in the entire IPv4 space. We call these two monitoring systems as "the 16-block monitoring system" and "the 1024-block monitoring system," respectively.

The previous Figure shows one simulation run of the Blaster-like worm. Plot (a) in that Figure shows the number of infected hosts $I(t)$ in the entire Internet as a function of time t . It also shows the cumulative number of observed infected hosts, $C(t)$, from both monitoring systems. Because observed $C(t)$ is very small compared with $I(t)$, we plot this figure by taking logarithmically on the Y-axis.

Plot (a) shows that, during the worm's propagation period, we can observe less than 0.1% of infected hosts in the Internet from the 16-block monitoring system. Even if we use the 1024-block monitoring system, we can only observe less than 4% of infected hosts in the Internet during the worm's propagation period. This is the reason why researchers have not derived an accurate estimate of how many computers were really infected by the Blaster-like worm.

Plot (b) in that Figure shows the monitored data $Z(t)$, the number of worm scans observed within each minute. Compared to the 16-block monitoring system, The 1024-block monitoring system gives noisier observation $Z(t)$. This is because as time goes on, an infected host will enter or leave one of the monitored IP blocks. It happens more frequently in the 1024-block monitoring system than in the 16-block monitoring system.

Although noisier than the data from the 16-block monitoring system, the monitored data from the 1024-block monitoring system represents more accurately the propagation of a sequential-scan worm. From the monitored data sets, we want to know the worm's propagation pattern in the global Internet, i.e., the curve of I_t shown in Plot (b). Such a growth pattern of I_t is a low frequency signal compared with the high frequency noise presented in the observed data Z_t . Therefore, we can use a low-pass filter to filter out high frequency noise from Z_t without changing the worm's propagation pattern. Plot (c) in that Figure shows the observation data Z_t after being filtered by a first-order low-pass filter. This Figure clearly shows that the monitored data from the 1024-block monitoring system can better represent the worm's propagation pattern in the entire Internet.

Based on the filtered monitored data Z_t from the 1024-block monitoring system as shown in Plot (c), we run the Kalman filter estimation based on the transformed linear model. The estimated

value of α is shown in the following Figure as a function of time. In this simulation run, the Blaster-like worm infects 1.3% of vulnerable population at time 240 minutes, by which time the estimate of α has already stabilized and oscillated slightly around a positive, constant value. Hence our early detection system can detect the Blaster-like worm before it infects 1.3% of vulnerable population in the Internet.

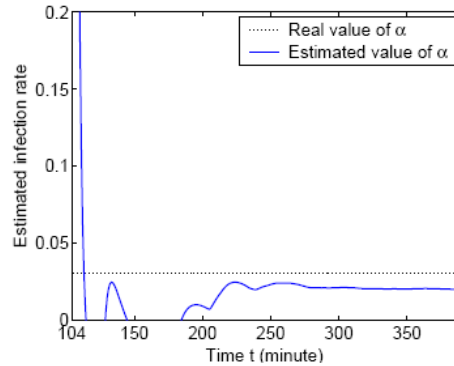


Figure 40: Kalman filter estimation of worm infection rate for Blaster-like worm based on transformed linear model

Worm propagation in other simulation runs of the Blaster-like worm gives results similar to those shown in the previous two Figures. On occasion the 16-block monitoring system provides as good observation as the 1024-block monitoring system. However, the 1024-block monitoring system always provides stable and good observations, while the 16-block monitoring system provides poor observations in many instances.

C8. Discussion and Future Work

We have used the simple epidemic model and the exponential model for the estimation and prediction. While these models give good results so far, we need to develop more detailed models to reflect a future worm's dynamics. For example, if a worm spreads through a topology, or spreads by exploiting multiple vulnerabilities, or is a meta-server worm, then its propagation may not follow the models used here.

The monitoring interval Δ is an important parameter in the system design. For a slow-spreading worm, it could be set to be long, but for a fast-spreading worm such as Slammer, the time interval should be in the order of seconds to catch up with the worm's dynamics. How can we select the appropriate Δ before we know a worm's presence and its speed? We need to do further research on designing a recursive estimation algorithm that uses adaptive sampling rate. Currently, one way we contemplate is to tag the time stamp with each observed scan. Then at MWC, several estimators run in parallel with different monitoring intervals. From the tagged time stamp the correct C_t or Z_t for every estimator can easily be restored.

It could be useful to develop distributed estimation algorithms so as to reduce the latency and traffic for the report to a central server. Distributed estimators may also reduce the impact of noise when a few monitors experience larger than normal noise-to-signal ratios. In addition, we want to use a continuous version of the Kalman filter. This approach would reduce the

significance of the monitoring interval selection, and would work nicely with the distributed estimation setting.

The worm detection method presented here assumes that only worm scans can cause exponentially increased traffic to monitors, while other background scan noise cannot. We believe this is a reasonable assumption. If we want to further improve the detection accuracy, however, we can add some other rule sets in the detection system. For example, in order to distinguish a worm attack from a DDoS attack, we can exploit the differences between them: a DDoS attack has one or several targets while a worm's propagation has no specific target.

We have presented several major issues in designing an Internet Malware Warning Center. However, there are still many challenges in building such a system, such as cooperation mechanism among a large number of communities; the privacy concern in monitored data sharing, and the robustness of the monitoring system itself towards attacks by worms or hackers.

C9. Conclusions

We propose a monitoring and early detection system for Internet worms to provide an accurate triggering signal for mitigation mechanisms in the early stage of a future worm. Such a system is needed in view of the propagation scale and the speed of the past worms. We have been lucky that the previous worms have not been very malicious; the same cannot be said for the future worms. Based on the idea of "detecting the *trend*, not the *burst*" of monitored illegitimate scan traffic, we present a "trend detection" methodology to detect the presence of a worm in its early propagation stage by using the Kalman filter and worm propagation models. Our analysis and simulation studies indicate that such a system is feasible, and the "trend detection" methodology poses many interesting research issues.

Appendix D: Network Resilience: Exploring Cascading Failures within BGP

D1. Introduction

The Internet is a large collection of Autonomous Systems (AS). There are various glues that hold together this massive system, but the most important of these is the routing infrastructure. The Internet routing protocols maintain connectivity between and within AS's, and are designed to automatically reconfigure and recompute routing tables when they detect a link failure. This computation starts locally around the failure point, and then the information propagates through the Internet. While telephone networks recover from failures on the order of milliseconds [2], the Internet routing table convergence has been observed to take a much longer time [3]; it is not uncommon to have tens of minutes of downtime before recovery has taken place. Although empirical observations of the dynamics have been made in [4, 5], formal models of the process are not very common (there has been recent work on an analytical study of Route Flap Damping [6], and another study of BGP in congested networks [7]). There are other models of BGP that verify correctness, satisfiability etc. [8, 9, 10]; here we explore the dynamical behavior of BGP. We develop and analyze models in this paper that reflect the behavior of the protocols. Our analysis confirms the observations made in [1], that it is indeed possible to have cascading failures in systems like the BGP routing infrastructure. Our results indicate the presence of phase transitions in these systems, and the presence and intensity of the phase transitions are strongly dependent on system parameters. We use the term *phase transitions* in the sense used by Erdos and Renyi in their work on random graphs, that is "an abrupt change in a global system property". The phase transition is to be interpreted as a sharp threshold rather than the definition used in statistical mechanics. In our analysis, we observe that the propensity for phase transitions increases as clique size increases, and additionally also increases as the processing capacity of the routers decreases.

D2. Background

An interesting vulnerability of BGP came to light in July and September of 2001. This was the incident where CODE RED and Nimda viruses disrupted the Internet routing infrastructure and caused widely reported BGP storms [1]. We briefly describe the incident, discuss possible explanations and then elucidate how modeling can help in this scenario. A direct quote from [1] is reproduced below

On July 19th, we observed an exponentially growing eight-fold increase in the advertisement rate, over a period of about eight hours (all times are in GMT; subtract 4 hours for EDT). This BGP surge faded over the same time scale as it arrived. When one considers the conventional wisdom about BGP convergence times (seconds to minutes), it is more than a little disturbing to see a fundamental quantity like BGP advertisement rate exhibiting exponential growth for eight hours.

This event coincided with the spread of the CODE RED virus on the Internet. Then again, on September 18th, the day the NIMDA virus started to spread, the following was observed:

On Tuesday, September 18, simultaneous with the onset of the propagation phase of the Nimda worm, we observed another BGP storm. This one came on faster, rode the trend higher, and then, just as mysteriously, turned itself off, though much more slowly. Over a period of roughly two hours, starting at about 13:00 GMT (9am EDT), rrc00 aggregate BGP announcement rates exponentially ramped up by a factor of 25, from 400 per minute to 10,000 per minute, with sustained "gusts" of more than 200,000 per minute. The advertisement rate then decayed gradually over many days, reaching pre-Nimda levels by September 24th.

These two events clearly demonstrated that an application layer event (the virus attack) caused problems at lower layers of the Internet infrastructure. What caused this unexpected event? Briefly, [1] hypothesized the following chain of events

- The viruses started random IP port scanning
- Most of these random IP addresses were not in the cached entries of the routing table, causing....
- Frequent cache misses, and,
- In the case of invalid IP addresses, generation of ICMP (router error) messages.
- Both of the above causes led to router CPU overload, causing routers to crash
- Router failure led to withdrawal announcements by the peers, generating a high level of advertisement traffic.
- When the router came back on, it required a full state update from its peers, creating a large spike in the load of its peers who provided the state dump
- Once the restarted router obtained all the dumps, it dumped *its* full state to all its peers, creating another spike in the load..
- Frequent full state dumps led to more CPU overload, leading to more crashes, and the propagation of the cycle...

An immediate question of concern is: how self-sustaining is this process? On the two particular days, the BGP storms lasted for hours, causing major disruptions on the Internet. Although there is some debate ranging over the findings of [1], e.g. [11], our interest is answering the question if this process (or something similar) can result in cascading failures. In the next two sections, we develop simple models to analytically understand this behavior.

Before we begin our modeling, we give a brief overview of BGP behavior. BGP is the inter-autonomous system (AS) routing protocol. At the boundary of each autonomous system, peer border routers exchange network reachability information with other autonomous systems through BGP. BGP uses TCP as its transport protocol. Two BGP speakers form a transport protocol connection between one another, and they exchange messages to open and confirm the connection parameters. When a connection is first established, a BGP speaker sends its entire routing table to

the peer (a full state dump). During the following BGP session, incremental updates are sent as the routing table changes. Two types of BGP messages are important to BGP operation. First is Keep Alive message, which is sent periodically to ensure the connection is live. If the peers can't receive KeepAlive messages in a preset period of time, the BGP connection has to be closed. Physical connectivity failure (link failure, router crash), transient connectivity problems due to congestion, or even manual reboots, may result in the delay of KeepAlive message to the peers. When BGP sessions restart, the peers have to send the full routing table again. Update messages are used to exchange routing information change between two peers. Route withdrawals are sent when a router makes a new local decision that a network is no longer reachable. We study the clique topology (fully connected mesh) in subsequent sections. The clique topology provides the most complex interactions, and in the core of the Internet the major Autonomous Systems form a clique.

D3. Simple Fluid Model

Let's build a very simple model of the interactions in the system and explore its properties. We consider a simple scenario with a finite set of N routers, that are all connected to each other, i.e. form a clique. Let the number of down routers at any instant be $D(t)$. We define a "down" router to be one that does not have a functioning routing table, so a router that is in the process of rebooting and obtaining state dumps is also defined to be "down". Now, we study the system of the number of down routers. The arrival and departure process to the system is defined as:

$$\alpha(t) \doteq \text{Number of arrivals in } [0, t) \quad (1)$$

$$\delta(t) \doteq \text{Number of departures in } [0, t) \quad (2)$$

Consider the process $\delta(t)$. The down routers come up with the help of the routers that are currently up ($N(t) - D(t)$). We define the service rate of an up router as k_s , where k_s is the average number of down routers a functioning router restores per unit time. Now, if $N(t) - D(t)$ servers are up and providing service, the service received by a single down node is its share of the total service capacity of the system. Thus, the share received by a single router is $(N(t) - D(t))/D(t)$. To account for the boundary condition $D = 0$, the denominator should actually be the term $(D(t) + k_a)$, where k_a represents the ambient load on the servers, representing for instance processing of normal route advertisements, and prevents $D \rightarrow \infty$. However, for simplicity of exposition we ignore the term as it does not affect the main observation we obtain later. Hence, the number of departures in an infinitesimal time dt , $d\delta(t)$, is defined by

$$d\delta(t) = D(t) \frac{N(t) - D(t)}{D(t)} k_s dt = (N(t) - D(t)) k_s dt \quad (3)$$

Resetting of BGP sessions lead to two kinds of messages, withdrawal announcements and subsequently full state updates when the BGP session is restored. We model the rate at which a functioning router goes down due to the load imposed by the resetting of a single BGP session as k_l . Typically, we expect $k_s \gg k_l$, as BGP resets are not uncommon and restoring a single session is unlikely to cause a peer router to go down. Now, the average arrivals in an infinitesimal time dt , $d\alpha(t)$, is given by the product of three quantities: the constant k_l , the

number of routers (BGP sessions) that are currently down $D(t)$ (denoting the total load offered), and finally the number of routers currently up (that can go down), $(N(t) - D(t))$, i.e.

$$d\alpha(t) = k_l D(t)(N(t) - D(t))dt \quad (4)$$

Now $D(t)$ is $\alpha(t) - \delta(t)$, hence combining (3) and (4) we obtain the drift equation

$$dD(t) = D(t)(N(t) - D(t))k_l dt - k_s D(t) \frac{N(t) - D(t)}{D(t)} dt \quad (5)$$

Simplifying and dividing by dt both sides, we obtain the following relation

$$\frac{dD}{dt} = -k_l D^2 + (k_s + k_l N)D - k_s N \quad (6)$$

This is a Riccati equation, and without going into the actual solution of the equation, we immediately observe that the dynamical system described by this model exhibits a *phase transition*: If the initial state $D(0)$ of the system is above a certain threshold, then as $\lim_{t \rightarrow \infty} D(t) = N$, else $\lim_{t \rightarrow \infty} D(t) = 0$. In other words, if by some exogenous process (e.g. CODE RED) we manage to bring a certain number of the routers down, thereby resetting the BGP sessions, then depending upon that number the system either fully recovers or there is a *cascading failure*. This is seen in plot (a) in the Figure below, where we plot the RHS of (6) as a function of D . This is very similar to results obtained in epidemic modeling, see [12]. For a value of D above k_s/k_l , the derivative of D remains uniformly positive, making the system drift to the state where all routers are down. For D below that number the derivative remains uniformly negative, bringing the system back to recovery. An interesting fact here is that this threshold is an *absolute* quantity rather than a fraction of N . Hence, given the right parameter set this phase transition may not be exhibited at all. A simulation of the system with different initial conditions is shown in plot (b) in the Figure below, where $k_s/k_l = 20$ and we plot two trajectories, one with $D(0) = 21$ and another with $D(0) = 19$. The presence of phase transition in the fluid model piques our interest enough to attempt a refinement of the model, and observe the behavior. In the next section, we move away from the fluid model and develop a birth-death model for the system.

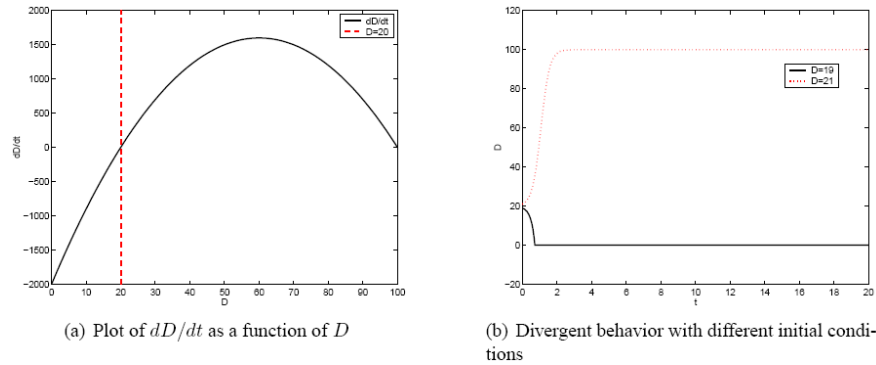


Figure 41: Phase transition in model

D4. A Birth-Death model

Now we construct and analyze the discrete analog of the basic fluid model we studied in the previous section. Again, we assume there are N identical routers in the system, forming a clique. A state in the birth-death represents the number of down routers in the system. The model is depicted in the following Figure.

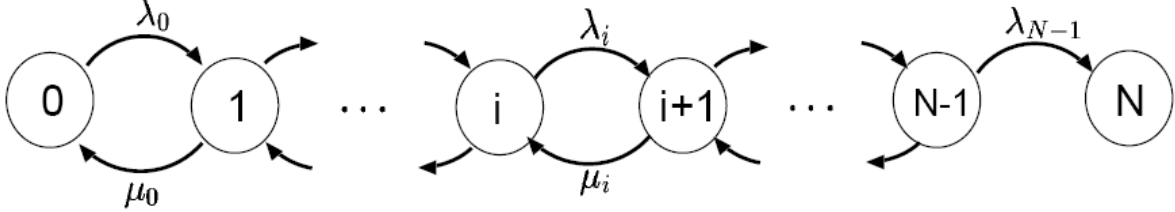


Figure 42: Birth-Death Process - state i represents number of down nodes

The transition rate from state i to $i + 1$ is given by $\lambda_i = (N - i) \times (i) \times k_l + k_a$, $i = 0 \dots N-1$; and similarly the transition from $i + 1$ to i is given by $\mu_i = (N - i) \times k_s$, $i = 0 \dots N-2$. μ_{N-1} is defined to be zero, as there is no repair when all routers are down, hence state N is an *absorbing* state. To understand the behavior of this model, we have to perform a transient analysis, as the system ends up in state N with probability 1. The mean time to absorption from state i to N , W_i , is a good indicator of the behavior of the system. Physically, W_i is to be interpreted as roughly the time it takes for the system to collapse if i routers are currently down. Next, we compute W_i . In state i , the mean time to the next transition out of state i is $1/(\lambda_i + \mu_{i-1})$, $0 < i < N$. At that transition, the expected remaining time to absorption is W_{i+1} with probability $\lambda_i/(\lambda_i + \mu_{i-1})$ and W_{i-1} with probability $\mu_{i-1}/(\lambda_i + \mu_{i-1})$. Then

$$W_i = \frac{\mu_{i-1}}{\lambda_i + \mu_{i-1}} W_{i-1} + \frac{\lambda_i}{\lambda_i + \mu_{i-1}} W_{i+1} + \frac{1}{\lambda_i + \mu_{i-1}} \quad (7)$$

Next, we apply the boundary conditions that are $\mu_{-1} = 0$, and $\lambda_{-1} = 0$. The first condition yields

$$W_0 = \frac{1}{\lambda_0} + W_1 \quad (8)$$

while the second boundary condition yields

$$W_{N-1} = \frac{1}{\lambda_{N-1} + \mu_{N-2}} + \frac{\mu_{N-2}}{\lambda_{N-1} + \mu_{N-2}} W_{N-2} \quad (9)$$

Combining equations (7), (8) and (9), we obtain for $i > 1$

$$W_i - W_{i+1} = \sum_{j=0}^{i-1} \frac{\prod_{k=j}^{i-1} \rho_k}{\lambda_i} + \frac{1}{\lambda_i} \quad (10)$$

Where $\rho_k = \lambda_k/\mu_k$. With $W_N = 0$, (10) yields a way to compute W_i as a function of i .

D5. Model Analysis

The phase transition that we encountered in the fluid model is manifested in the form of a sharp decrease in W_i in the birth-death model. We explore the parameter space of N , k_i , k_s and k_a to identify possibilities of cascading failures. First we assume that the "ambient" load, k_a , is a small fraction of the full state update load k_i . We choose the fraction to be $1/100$. Now we vary the ratio k_i/k_s and the size of the system N and observe the behavior. For k_i we assume a value $0.01/s$, i.e., the rate of crashing of a router under the load of a full state update is 0.01 per second, and for k_s we assume a value of $1/s$, in other words a working router brings up a crashed router at the rate of 1 router per second. First we assume a small clique, of 20 routers. The mean time to absorption is shown in plot (a) of the Figure below. As we can observe the mean time to absorption is of the order 10^{23} seconds, as long as even one router is up. Hence, this system is unlikely to undergo cascading failures. Next, we increase the size of the clique, from 20 to 100 , and plot the results in (b) of the Figure below.

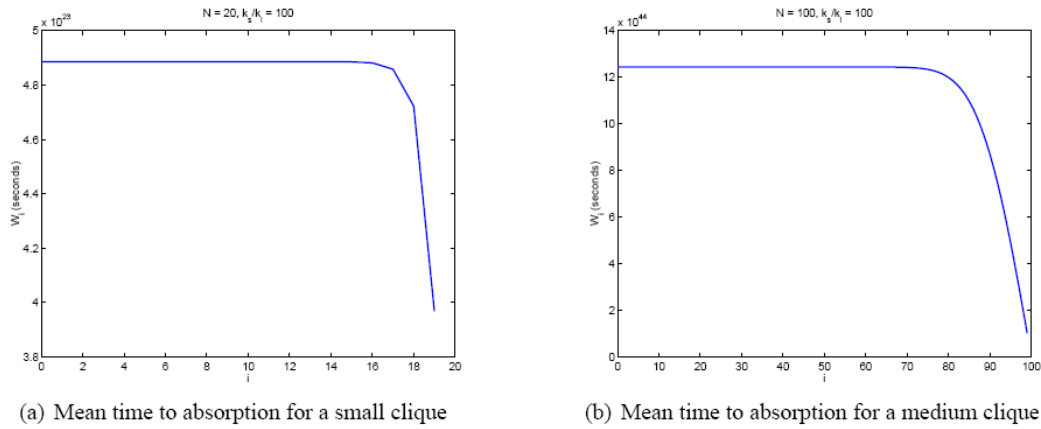


Figure 43: Stable behavior exhibited by small to medium cliques

The mean time to absorption now starts to show a decrease as i increases, but is still relatively quite high throughout. It is of the order 10^{44} throughout, hence when the system appears to have become more stable, as the meantime to absorption has increased by several orders of magnitude. This makes intuitive sense, as increasing the size of the clique increases the redundancy in the system and there are more up routers "available" to bring up crashed routers. Now, we increase the clique size to 200 , and show the results in plot (a) of the Figure below. Now, we observe the first appearance of *the phase transition*. At roughly $i = 100$, the mean time to absorption starts decreasing rapidly, and in a matter of few states, the time falls from an order of 10^{44} to close to 0 ! Thus, an apparently highly stable system becomes extremely vulnerable and can cascade into a collapsed state, where all routers are down in no time. This behavior is again observed in a clique of size 400 , shown in plot (b) of the following Figure. The transition appears at around the same spot, around state 100 .

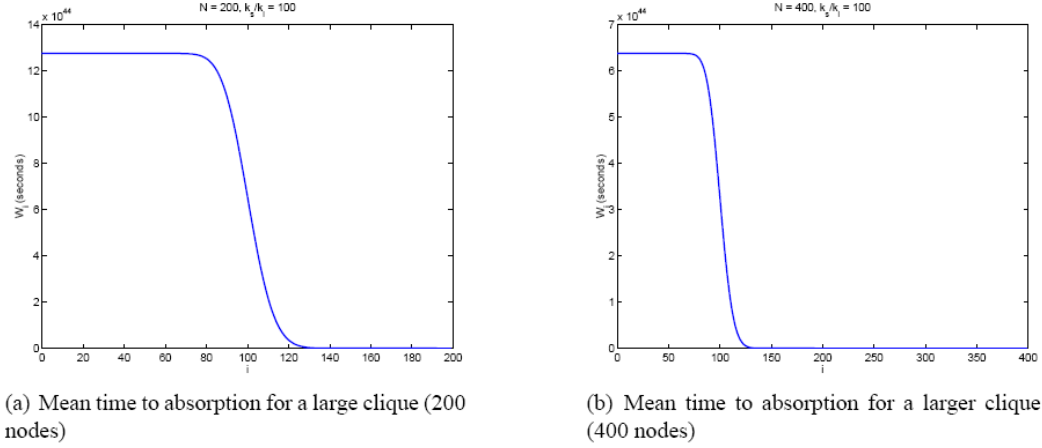


Figure 44: Phase transitions observed in larger cliques

The state 100 around which the phase transition starts to take effect is roughly the ratio k_s/k_l . Note that this is similar to what we observed in the fluid model. This can be observed in the Figure below, where we plot the mean time to absorption for an identical clique size of 200, but vary the ratio k_s/k_l .

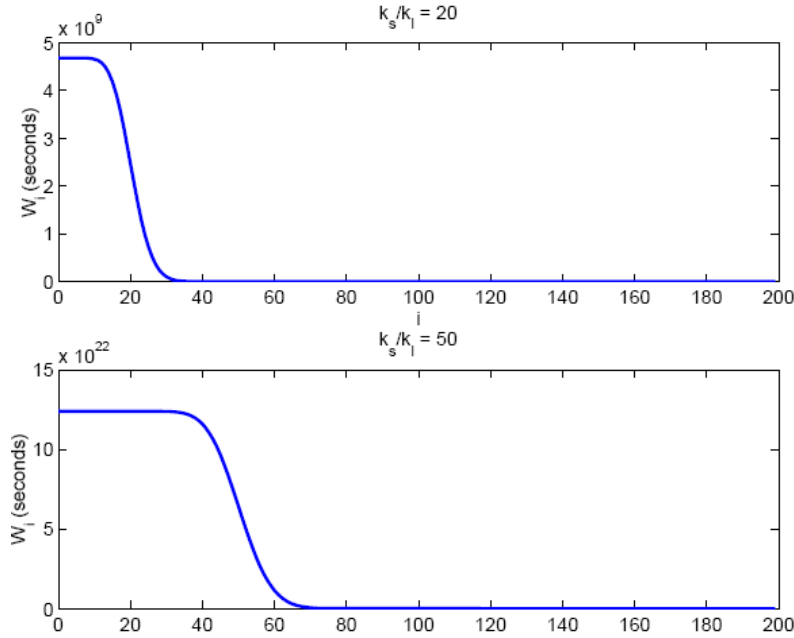


Figure 45: Phase transition point as a function of k_s/k_l

The effect of the clique size on cascading failures is also interesting. We now fix the ratio k_s/k_l , and vary N . The results are plotted in the following Figure. Increasing N does not change the location of the phase transition, but does affect the relative stability before the transition. As we increase N , beyond the size where the system shows a phase transition, the relative stability *decreases*. This is in contrast to the scenario we earlier saw, where increasing N increased the relative stability for small to medium clique sizes.

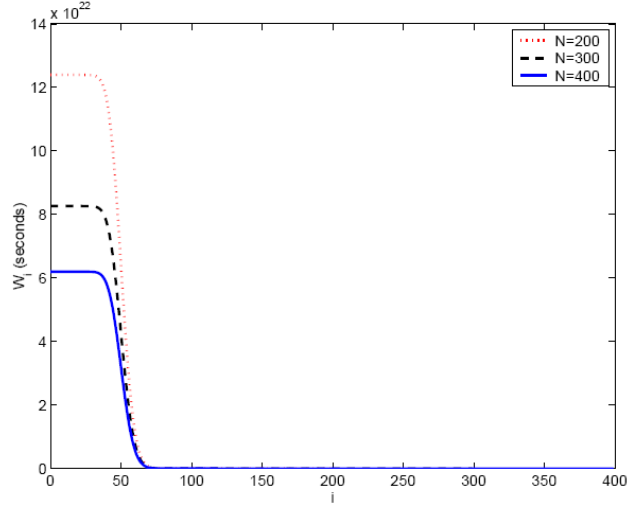


Figure 46: Relative stability and phase transition point for $N=200, 300, 400$

This behavior is more clearly observed in the Figure below, where we plot the mean time to absorption as N goes from 80 to 100 in steps of 10. We observe that the relative stability increases as N increases.

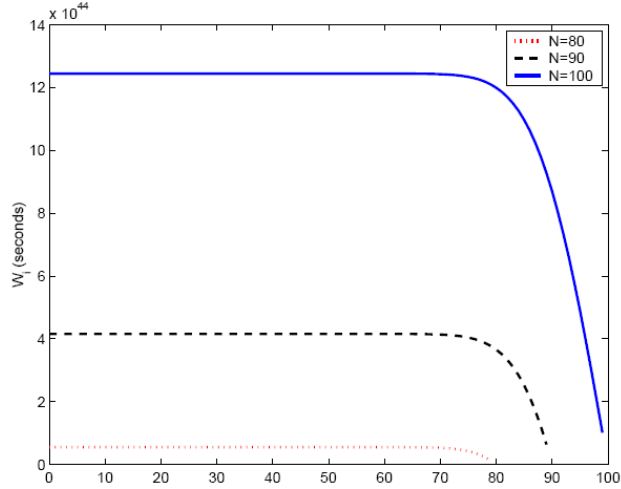


Figure 47: Relative stability and phase transition point for $N=80, 90, 100$

D6. Conclusions

We have developed simple analytical models to capture the interaction of systems like BGP routing. We specifically develop models for the clique topology (the one that is used at the AS level on the global Internet) and discovered phase transitions with respect to cascading failures in the models. Our findings confirm the measurement based inferences made in [1] on possibilities of cascading failures in the routing infrastructure. Our models indicate that both the size of the clique as well as the capacity of the nodes in the clique is an important consideration

for the phase transitions. The size of the clique acts as a threshold for the phase transitions, given other parameters, the clique must be large enough for the transition to appear. Increasing the clique sizes beyond the threshold does not change the location of the phase transition, but does have an effect on relative stability. On the other hand, if the clique size is large enough, then the capacity of the nodes in the system decides the location where the phase transition occurs.

In terms of future work, we are working on a more detailed model for a single router to understand individual node failures. A better understanding of individual node behavior would then permit us to model more complex topologies, moving beyond the homogeneous cliques that we have studied here. An interesting question that our models can answer is to what kinds of interactions make the system most resilient? In other words, the transition rates in our model assume a certain kind of interaction. If that interaction is modified, does that lead to more resilient networks? That question of course has to be asked in conjunction with the issue of the performance of system under normal circumstances, and these are questions that we are trying to answer.

Appendix E: Analysis and Design of Controllers for AQM Routers Supporting TCP Flows

E1. Introduction

The development of new Active Queue Management (AQM) routers will play a key role in meeting tomorrow's increasing demand for performance in Internet applications. Such applications include voice over IP (VoIP), class of service (CoS) and streaming video where packet size and session duration exhibit significant variations. In this context we have three objectives. First, to relate key network parameters to the AQM problem. Secondly, to analyze the present *de facto* AQM standard: *random early detection* (RED) and finally, to recommend alternative AQM schemes. The uniqueness of our approach comes from the use of a recently developed dynamic model of the Transmission Control Protocol (TCP) which enables application of control principles to address the basic feedback nature of AQM.

To begin, we first consider a simple sender-receiver connection passing through a bottleneck router as shown in the Figure below and schematic in the subsequent Figure. Under TCP, a sender probes the network's available bandwidth by linearly increasing its rate until data packets are lost. Upon packet loss, the receiver signals the sender to reduce its rate. Some drawbacks in this packet-dropping scheme include flow-synchronization and performance degradation due to excessive time-outs and restarts. Motivated by these network inefficiencies, the RED scheme was introduced in [2] to allow the router to assist TCP's management of network performance. Rather than waiting for packet loss to occur, RED acts preemptively by measuring the router's queue length and throttling the sender's rate accordingly. Since TCP is an end-to-end protocol, RED achieves this feedback indirectly by randomly dropping/marking packets and routing them to the receiver.

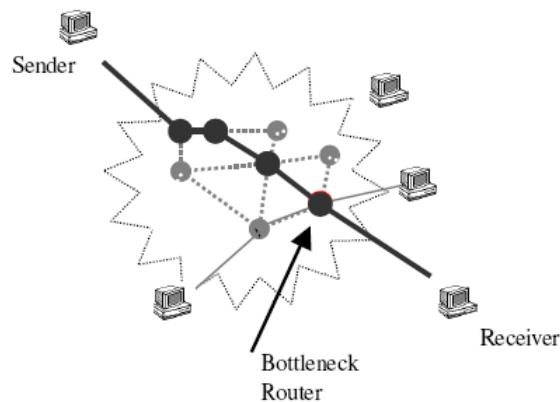


Figure 48: A single sender-receiver connection

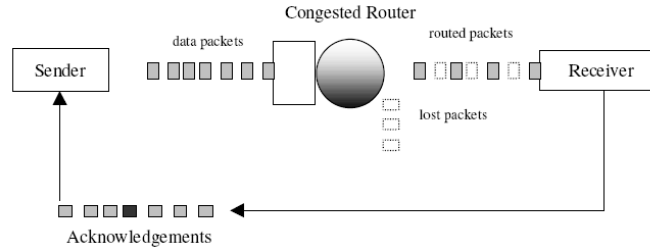


Figure 49: A schematic of a sender-receiver connection

The receiver, in turn, completes the feedback by acknowledging the receipt of marked packets to the sender; this is depicted in the following Figure where we emphasize the implicit, delayed, feeding-back of acknowledgment packets. Upon receipt of such acknowledgments, the sender increases or decreases its rate according to the TCP algorithm. The randomness in RED's packet-marking scheme is meant to eliminate flow-synchronization while queue-averaging was introduced to attenuate the effects of bursty traffic on the feedback signal. A drawback in deploying RED stems from its apparent tuning difficulties (see [4] and [5]) and the research community has responded with modifications such as in [6], [7], [8], [9], [10] and [11]. It has also motivated our research. Central to our approach is the recognition that AQM schemes, such as RED, use feedback (evident in the following Figure) to regulate queue efficiency. Consequently, feedback control principles appear to be an appropriate tool in the analysis and design of AQM strategies. While such principles can be found in the analysis of ATM networks (see for example [12] - [14] and the references cited therein) they have not been applied to TCP-controlled flows. This is apparently due to a lack of analytical model of TCP's congestion-avoidance mode. Recently, there has been progress in modeling of TCP; see [15] and [16].

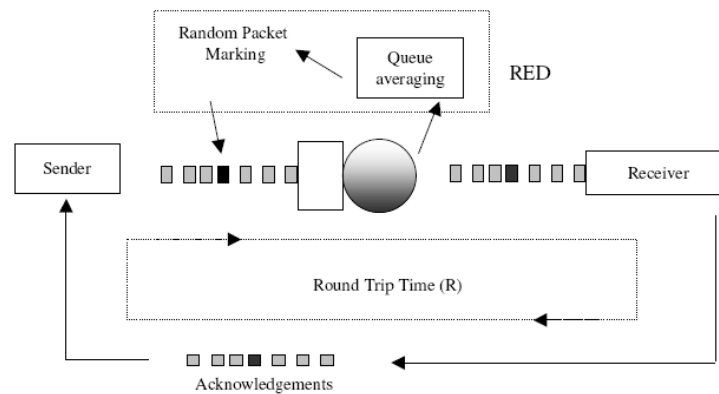


Figure 50: RED randomly marks packets to anticipate congestion

Our launching point is the fluid-flow model of TCP behavior presented in Section 3 (see ref. [17]). This model expresses TCP in a language that allows control engineers to analyze and design AQM schemes. To be more specific, this model enables us to do several things. First, to relate key network parameters such as TCP load, link capacity and round-trip time to the underlying feedback control problem. Secondly, to analyze RED and recommend that *averaging* is not beneficial, and finally, to suggest alternative AQM schemes which amount to classical proportional (P) and proportional-integral (PI) control. The benefits of these schemes are illustrated through ns simulations. One of the most promising outcomes of this work is the impact PI control has on

queue utilization. The key feature is that PI control allows one to explicitly set the network queuing delay - independent of network parameters.

E2. Dynamics of TCP's congestion-avoidance flow-control mode

In [17], a dynamic model of TCP behavior was developed using fluid-flow and stochastic differential equation analysis. Simulation results demonstrated that this model accurately captured the dynamics of TCP. We use a simplified version of that model which ignores the TCP timeout mechanism. This model relates the average value of key network variables and is described by the following coupled, nonlinear differential equations:

$$\begin{aligned}\dot{W}(t) &= \frac{1}{R(t)} - \frac{W(t)}{2} \frac{W(t - R(t))}{R(t - R(t))} p(t - R(t)) \\ \dot{q} &= \begin{cases} -C + \frac{N(t)}{R(t)} W(t), & q > 0 \\ \max\left\{0, -C + \frac{N(t)}{R(t)} W(t)\right\}, & q = 0. \end{cases} \end{aligned} \quad (1)$$

where x denotes the time-derivative and

$$\begin{aligned}W &\doteq \text{average TCP window size (packets);} \\ q &\doteq \text{average queue length (packets);} \\ R(t) &\doteq \text{round-trip time} = \frac{q(t)}{C} + T_p \text{ (secs);} \\ C &\doteq \text{link capacity (packets/sec);} \\ T_p &\doteq \text{propagation delay (secs);} \\ N &\doteq \text{load factor (number of TCP sessions);} \\ p &\doteq \text{probability of packet mark.}\end{aligned}$$

The first differential equation in (1) describes the TCP window control dynamic. Roughly speaking, the $1/R$ term on its right-hand side models the window's *additive increase*, while the $W/2$ term models the window's *multiplicative decrease* in response to packet marking p . The second equation in (1) models the bottleneck queue length as simply an accumulated difference between packet arrival rate NW/R and link capacity C . We illustrate these differential equations in the block diagram of the following Figure which highlights TCP window-control and queue dynamics.

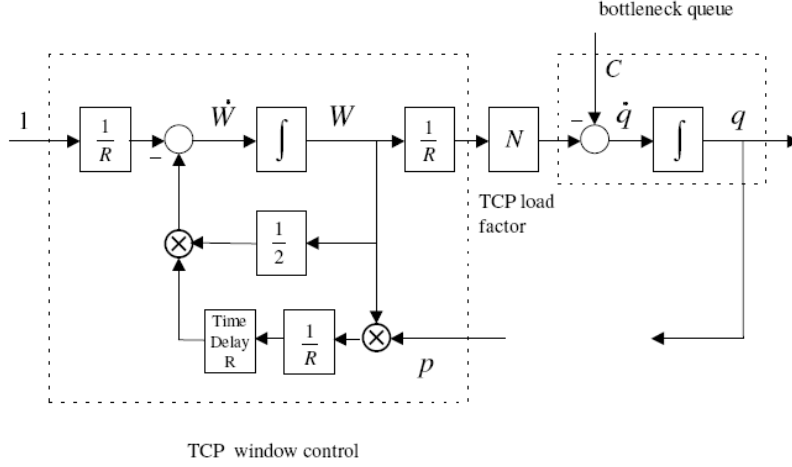


Figure 51: Block-diagram of TCP's congestion-avoidance mode

We now approximate these dynamics by their small-signal linearization about an operating point to gain insight for the purposes of feedback control (AQM). A block-diagram representation of these linearized dynamics is given in the following Figure where the TCP window-control and queue dynamics are explicitly identified.

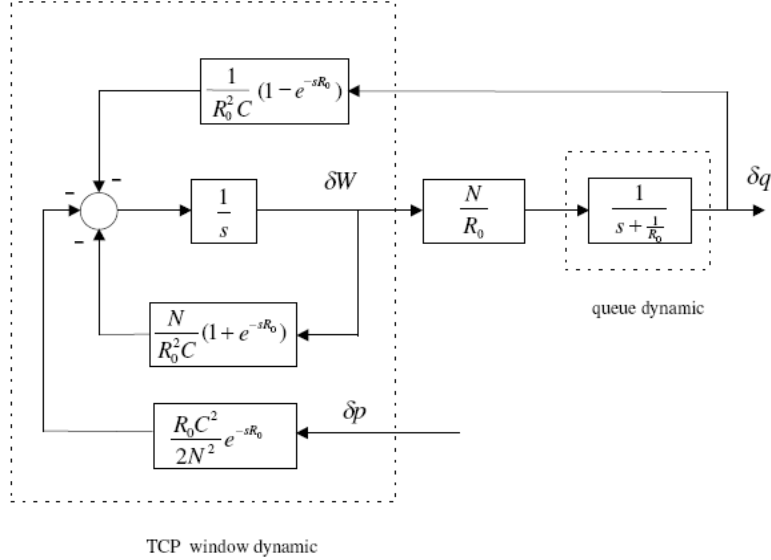


Figure 52: Block-diagram of the linearized TCP connection

A main reason for modeling and linearization of window and queue dynamics is for the purpose of the design and analysis of AQM schemes. To this end, we continue to simplify these dynamics by focusing on the nominal (low-frequency) behavior of the window dynamic and accounting for the residual behavior into a high-frequency parasitic. In the following Figure, we perform block-diagram manipulation to isolate this dynamic as well as identifying the high-frequency residual $\Delta(s)$.

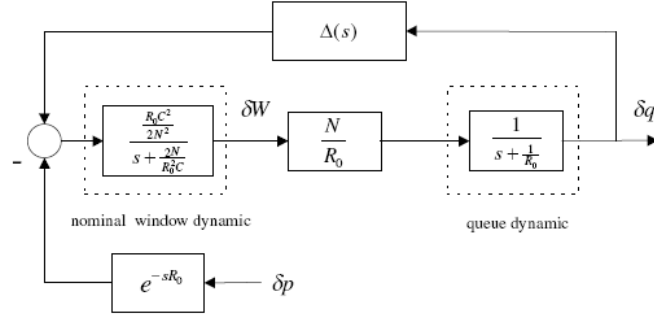


Figure 53: Linearized dynamic illustrating nominal window dynamic and high-frequency parasitic

E3. The AQM Control Problem

In the following Figure we give a feedback control system depiction of AQM. The action of an AQM control law is to mark packets (with probability p) as a function of measured queue length q . The plant dynamics, denoted by the transfer function $P(s)$, then relates how this packet-marking probability dynamically affects the queue length. The transfer function $\Delta(s)$ represents high-frequency window dynamics.

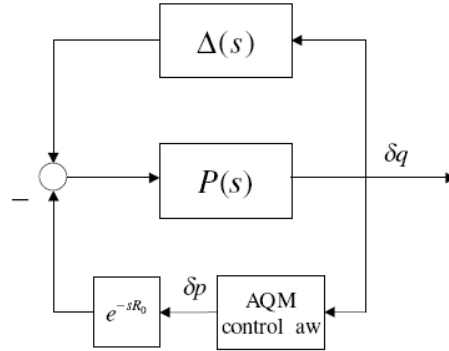


Figure 54: AQM as feedback control

The magnitude Bode plots for these transfer functions are shown in the following Figure. The Bode plots of P reveal the low-pass nature of the TCP-queue dynamics as well as the inverse dependence of loop gain on the number of TCP sessions N . The frequency response of residual Δ shows its influence at higher frequencies. One objective of an AQM design is to stabilize these residual dynamics.

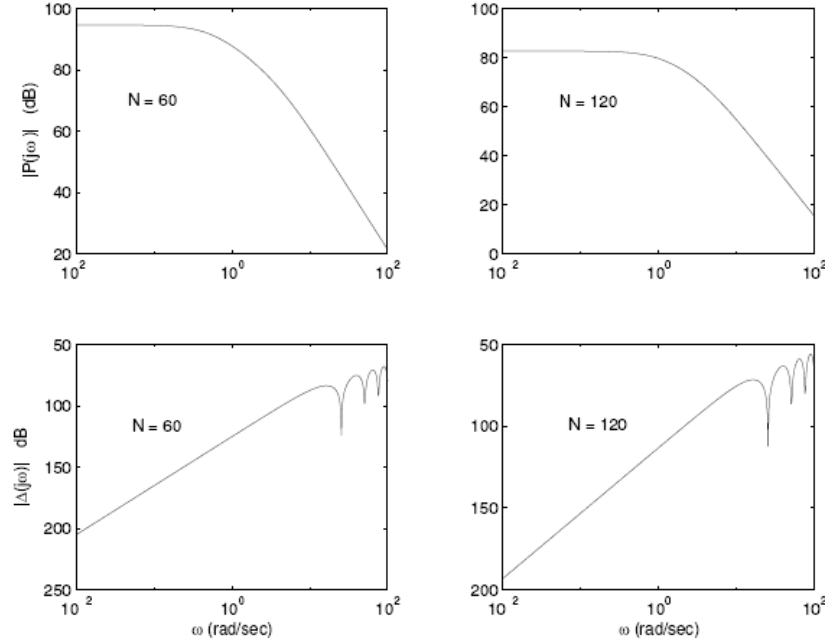


Figure 55: Magnitude Bode plots for $P(s)$ and $\Delta(s)$ for TCP loads of 60 and 120 sessions

As in any control system design, a first step is to pose performance objectives. For AQM, performance objectives include efficient queue utilization, regulated queuing delay and robustness.

- 1) *efficient queue utilization*: For efficient use, the queue should avoid overflow or emptiness. The former situation results in lost packets and undesired retransmissions, while an empty buffer underutilizes the link. Both of these extremes should be avoided in both transient and steady-state operation.
- 2) *queueing delay*: The time required for a data packet to be serviced by the routing queue is called the queuing delay and is equal to Cq . This time, together with the propagation delay T_p , accounts for the network's queuing delay and it is desirable to keep small both the queuing delay and its variations. This calls for regulating to small queue lengths; however, doing so may result in link underutilization and this limitation presents a fundamental tradeoff to AQM design.
- 3) *robustness*: AQM schemes need to maintain closed-loop performance in spite of varying network conditions. These conditions include variations in the number of TCP sessions N , variations in the propagation delay T_p .

Consider the linear control system in the following Figure where transfer function $C(s)$ represents a linear AQM control law. Closed-loop stability is fundamental in meeting the above performance objectives.

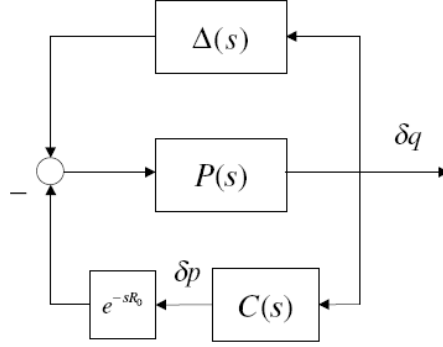


Figure 56: Block diagram of a linearized AQM control system

E4. AQM using RED

The simplest of congestion-avoidance scheme is the so-called *drop-tail* law which signals TCP sources to reduce window sizes whenever the queue overflows, which amounts to an "on-off" AQM control law. This on-off mechanism leads to queue-length oscillations, flow synchronization and performance degradation due to excessive time-outs and restarts.

Motivated by these network inefficiencies, the RED AQM scheme was introduced in [2] to allow the router to assist TCP's management of network performance. Rather than waiting for packet loss to occur, RED acts preemptively by taking an *average* measure of the router's queue length and throttling the TCP window accordingly by *randomly* marking packets. This randomness in RED's packet-marking scheme was meant to eliminate flow-synchronization, while queue-averaging was introduced to attenuate the effects of bursty traffic due to restarts and time-outs on the feedback signal. A drawback in deploying RED stems from its apparent tuning difficulties, see [4] and [5]. As we now show, we believe this difficulty stems in large part to RED's use of average queue length. Indeed, introduction of a low-pass filter into the feedback system in can lead to sluggish, oscillatory closed-loop behavior.

The RED active queue management control law computes the packet-marking probability p as a function of measured queue length q as depicted by the AQM control law in the block diagram of Figure E7. Specifically, RED consists of a low-pass filter (for queue averaging) and packet-marking profile as shown in the following Figure. Tuning RED amounts to selection of the low-pass filter pole K , threshold q_{min} , level p_{max} and gain L_{red} . The apparent motivation for introducing low-pass filtering in the AQM control law was to attenuate the effect of bursty, non-TCP controlled traffic on packet-marking. While this rationale for introducing low-pass filtering has some intuitive merit, it ignores the effect on closed-loop stability which amounts to introducing additional phase lag into a loop already containing time delay and two low-pass filter dynamics (associated with the TCP window and queue).

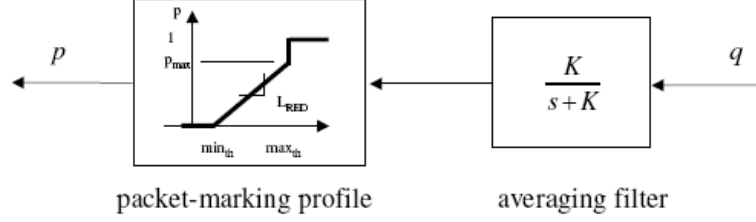


Figure 57: RED as a cascade of low-pass filter and nonlinear gain element

We illustrate our RED design via simulations using the ns simulator. Although the preceding analysis was carried out using the linearized model, the ns simulator captures the stochastic, nonlinear nature of the network dynamic. We considered a single bottlenecked router running RED and, in addition to the TCP flows addressed in our model, we also introduced short-lived http flows into the router to generate a realistic traffic scenario. The http flows were simulated using the http module provided with ns. The effect of these short-lived flows was to introduce an exogenous, noisy flow into the queue. In all of our plots the horizontal axis measures time (secs) while the vertical axis displays instantaneous queue length q (packets).

In the first simulation, we introduced 60 TCP flows and 180 http sessions. The capacity C is 15 Mb/s and the propagation delay T_p ranges uniformly between 160 and 240 ms. To provide a queuing delay of around 50-70 ms we set min_{th} and max_{th} of the packet-marking profile 10 to 200 and 250 respectively. The average packet size was set to 500 bytes. RED's averaging weight α and p_{max} were taken to be "vanilla;" i.e., the default values in ns. The buffer capacity was 800 packets. We set the *gentle* - parameter in RED to "on". The result is shown in Plot (a) of the following Figure which shows the oscillating nature of the queue length. The link is underutilized whenever the queue length goes to zero. Also, the large queue oscillation results in considerable variation in the round-trip times of packets.

Now we use the RED design in (11) and take the averaging weight α to be 1.33×10^{-6} , p_{max} to 0.1 and the dynamic range (min_{th} , max_{th}) as (150,700) packets. The results are plotted in Plot (b) of the following Figure. We see that the system response is stable, with fluctuations about an operating level of the queue. The larger oscillations experienced in the previous experiment are absent in this RED design. The slow response time is related to a low value of crossover frequency ω_q in our design. To improve the transient response, we can design for larger bandwidth ω_q ; however this would come at the expense of lower stability margins.

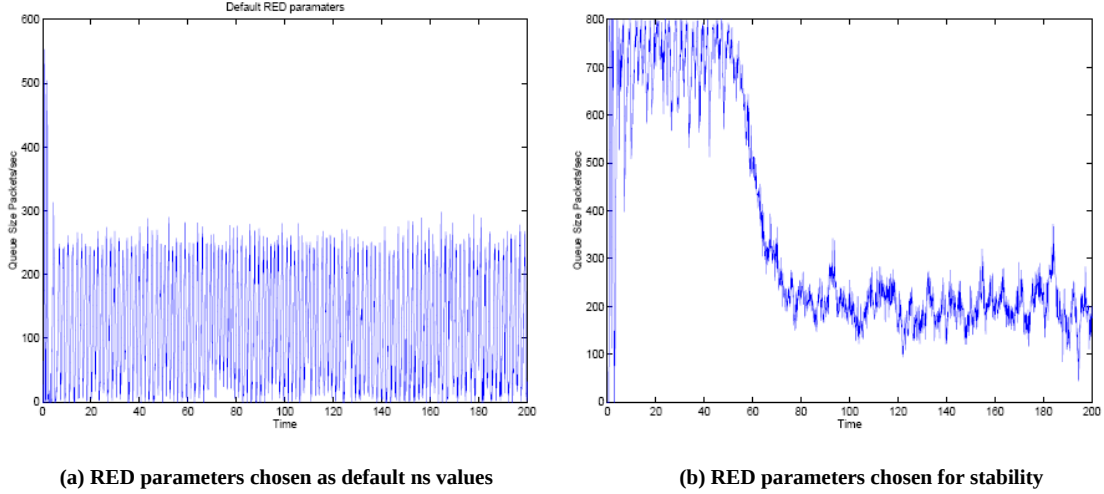


Figure 58: ns simulations comparing performance of RED controllers

E5. AQM using Proportional Control

The preceding RED controller resulted in small closed-loop bandwidths and thus sluggish behavior. One way to improve the response time of the system is to remove the low-pass filter completely. In doing so, we arrive at a classical, proportional controller. In proportional control the feedback is simply a scaling of the queue length and, in context of RED, this amounts to obtaining the packet-marking probability based on the instantaneous queue length rather than averaged queue length. Again, there exists a tradeoff between loop bandwidth and stability. However, the tradeoff is more favorable in the proportional case.

In this simulation, we consider a queue with 60 TCP flows and 180 http sessions. The link bandwidth is 15 Mb/s, and the propagation delays for the flows range uniformly between 160 and 240 ms, with average packet size being 500 Bytes. The buffer size is 800 packets. We also introduce a time-varying TCP load $N(t)$ to compare the speed of response between the RED and proportional controllers. At time $t = 100$, 20 of the TCP flows drop out and at time $t = 140$ they return. For the proportional controller, we set the averaging weight $\alpha = 1$ thereby removing the low-pass filter. We set the slope of the packet-marking profile to be the gain K_P , varying the loss linearly from 0 at queue length 100 with the slope specified by gain. Note that the buffer size of 800 puts an upper limit on the marking probability, which is approximately 0.04. We'll return to this issue following the experiment. The queue length plots are shown in Plot (a) of the following Figure. As evident from the traces, the proportional controller performs better, responding more quickly to load variations. However, this was to be expected since the closed-loop bandwidth for the proportional design exceeded that of the RED design by almost a factor of thirty.

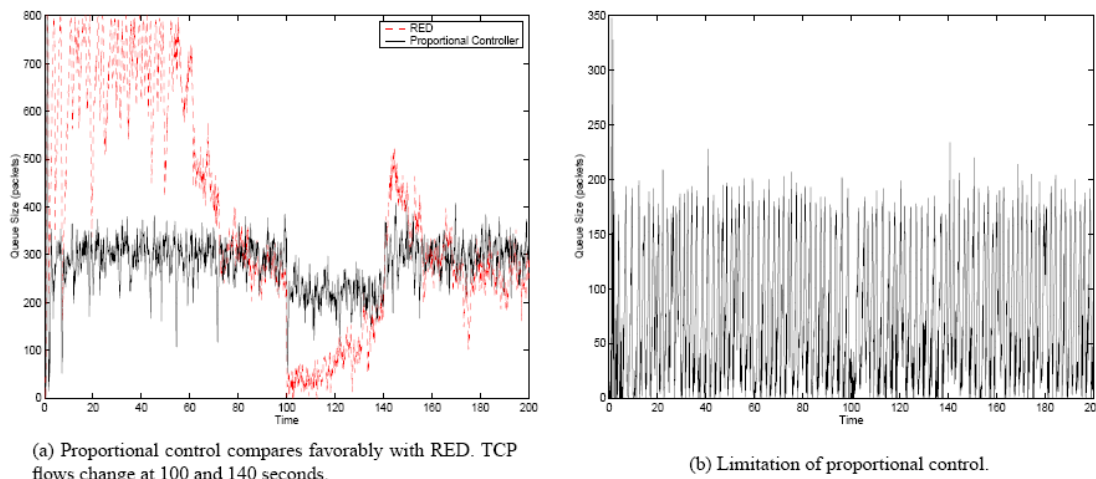


Figure 59: ns simulations of proportional control

While the proportional controller exhibits more responsive behavior than RED, it also suffers from a limitation which makes it impractical to implement under certain situations. For example, the steady-state buffer length is commensurate to the proportional controller's gain. Hence, buffer-size limitations could require gains outside the region of stabilizing proportional gains; such observations are also made in [11]. To illustrate, we repeat the previous experiment but change p_{max} from 0.04 to 1 for the proportional controller, to reflect a desire to keep the steady-state buffer length small. The result in Plot (b) of the previous Figure shows significant oscillations.

Increasing the buffer size to work around this problem is not an option since this could lead to unacceptably large queuing delays. The problem arises due to the coupling between queue length and marking probability. The two can be decoupled if we use integral control [22] in the AQM controller. Both the proportional and RED controllers result in control systems having (network-dependent) steady-state errors. For stable closed-loop systems, integral control drives the steady-state error to zero. Thus, we can design an integral controller for AQM which regulates the queue to a given operating level, independent of the load N . The simplest integral controller is the proportional-integral (PI) controller which appears appropriate for AQM context since, in comparison to RED, yields larger closed-loop bandwidths without sacrificing stability margins.

E6. AQM using Proportional-Integral Control

A PI controller has a transfer function of the form

$$C(s) = K_{PI} \frac{\left(\frac{s}{z} + 1\right)}{s}.$$

A desired consequence of the integral term is that δ_q in Figure E9 asymptotically converges to zero if $C(s)$ stabilizes. In the following Figure we show implementation of the PI control law with the nonlinear TCP/AQM dynamic (1) emphasizing the role of the queue's operating point q_0 .

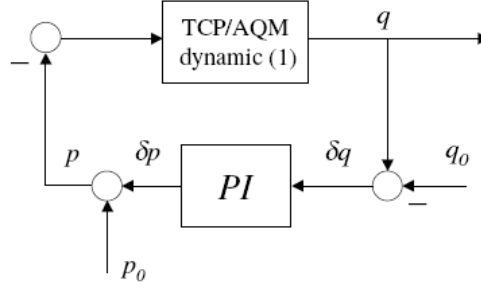


Figure 60: Implementation of PI controller emphasizing role of operating point q_0

The open-loop transfer function using PI control is

$$L(s) = \frac{K_{PI} C^2 \left(\frac{s}{z} + 1 \right) e^{-sR_0}}{s \left(s + \frac{2N}{R_0^2 C} \right) \left(s + \frac{1}{R_0} \right)}.$$

To validate the performance of the PI controller, we implemented it in ns with a sampling frequency of 160 Hz. The operating point was chosen as $q_0 = 200$ packets. We repeat the simulation shown in Figure E12(a), using the PI controller in lieu of proportional. The queue length plots for the RED and PI controllers are plotted in Plot (a) of the Figure below. The faster response time as well as the regulation of the output to a constant value by the PI controller is clearly observed. The PI controller is less sensitive to the load level variations and regulates the queue length to the operating point of 200 packets.

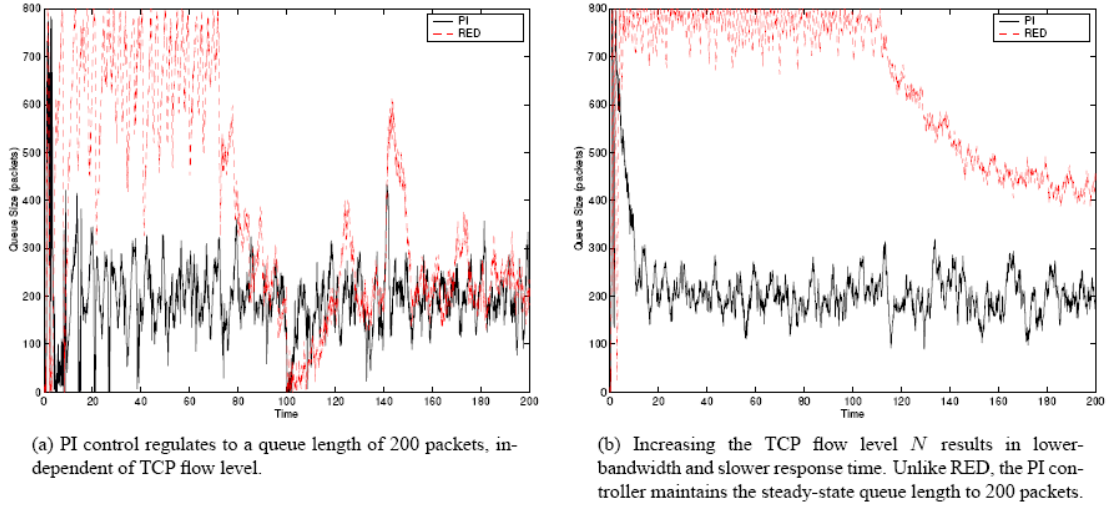


Figure 61: Comparison of RED and PI control under time-varying and heavy TCP loads

We now increase the number of TCP flows to 180 and http flows to 360. The response should be slower for this higher load level N . The queue lengths are plotted in Plot (b) of the previous Figure and we observe significantly better performance from the PI controller. The RED controller takes a long time to settle down, with the steady-state queue length quite large compared to the preceding simulation. The PI controller on the other hand is still regulating the queue length to 200 packets. The PI controller regulates to q_0 independent of TCP load.

In the next simulation we exercise the controllers at the other end of the load spectrum by reducing the TCP flows to 16 sessions. As observed in Plot (a) of the Figure below, the responses are more oscillatory corresponding to reduced phase margins. Finally, we stretch the controllers to the limit by increasing the number of TCP flows to 400. In Plot (b) of the Figure below the PI controller continues to exhibit acceptable performance, although its response has slowed. The RED controller, on the other hand, keeps the buffer in overflow. At such high load levels, the loop gain has decreased to a point where (large) regulating errors have pushed the steady-state queue length beyond the buffer size. This simulation illustrates the benefit of integral control in an AQM system with a finite buffer.

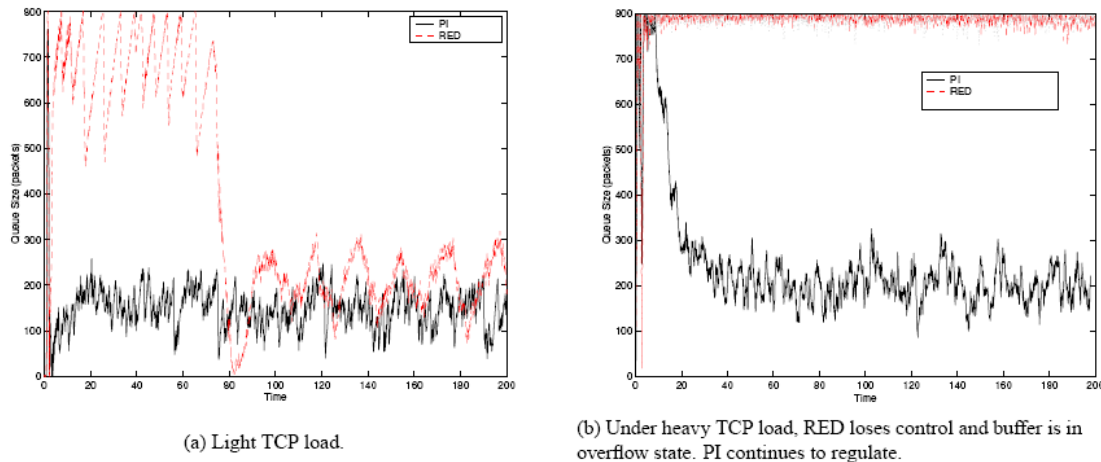


Figure 62: Comparison of RED and PI control under a light and very TCP loads

An important consideration in designing AQM systems is the tradeoff between queuing delay and utilization. Intuitively, larger buffers lead to higher utilizations of the link, but they also result in larger queuing delays. With the PI controller, the delay is essentially tunable with a single parameter q_0 . Larger values of q_0 give larger delay and utilization. In contrast, with RED, the delay is a function network conditions such as load level and packet-marking profile parameters min_{th} , max_{th} and p_{max} . We performed simulations to study this tradeoff as illustrated in the following three Figures where both pure ftp and mixed ftp and http flows are considered. As we observe in Figure E16 (a), small q_0 yields nearly full utilization in the case of pure ftp flows, whereas a larger q_0 is needed to reach this same level of utilization when both ftp and http are considered. The corresponding queuing delays are shown in Figure E16 (b) indicating a nearly linear relationship with q_0 . The corresponding delay-utilization curves are shown in Figure E17. We repeated these experiments with RED attempting to control delay through parameter min_{th} . We kept the range $max_{th} - min_{th}$ constant throughout. We ran the first experiment using a dynamic range of 550 and then repeated with a range of 55. We compare the performance with the PI design in Figure E18

where both long and short-lived flows were used. In the first of these figures, RED yields high utilization at the expense of large delays. When we reduced the queuing delay by lowering RED's dynamic range, utilization suffered. The PI design was capable of operating at both low delay and high utilization.

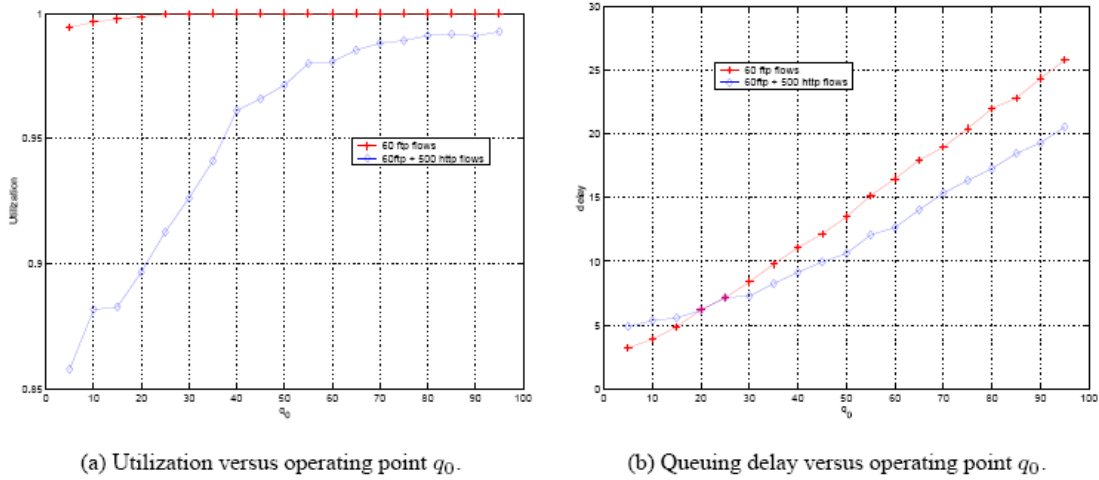


Figure 63: Utilization and queuing delay of the PI controller

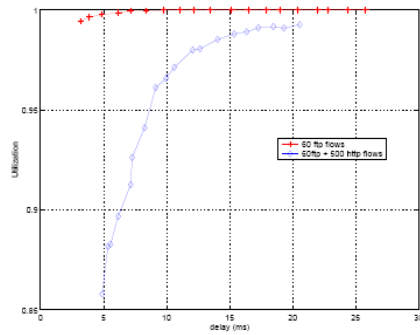


Figure 64: Utilization versus queuing delay: PI controller

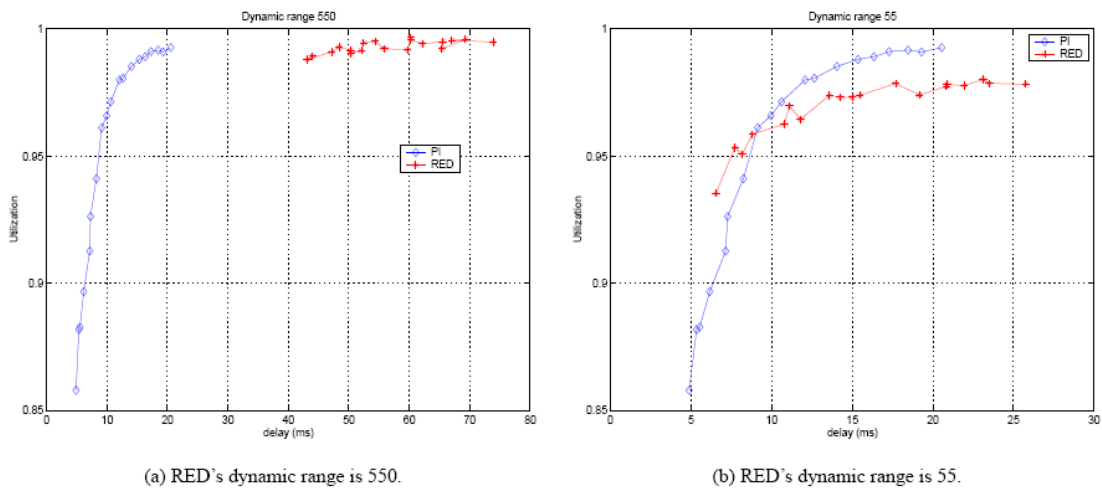


Figure 65: Queuing delay - utilization tradeoff: comparison between RED and PI control

E7. Conclusion

We analyzed a combined TCP and AQM model from a control engineering standpoint. We used linearization to analyze a previously developed nonlinear model of TCP. We performed the analysis on an AQM system implementing RED and presented design guidelines for choosing parameters which lead to local stability. We performed nonlinear simulations using ns which verified our analysis. In doing the analysis, we uncovered limitations of the averaging algorithm in RED. In addition we have proposed and designed two alternative controllers. The resulting control systems had faster response than the RED controller. The first of the designs, a proportional controller, displayed good transient response but suffered steady-state errors in queue regulation. This restricts its usefulness in systems where the buffer size is limited. Motivated by that limitation, we designed a classical PI controller which exhibited zero steady-state regulation error and acceptable transient behavior. The PI controller was simple to implement in ns which we compared under various scenarios with RED. The PI controller exhibited better performance under all cases considered. We also demonstrated the practical impact of the PI controller in managing queue utilization and delay. We have concentrated on simple and classical designs for AQM control. Modern control methodologies could be used; however, going this route may have obfuscated one of our main objectives which is to relate AQM control objectives directly to network parameters. Finally, there are a number of different areas in which the techniques presented here could be extended. Examples include networks with heterogeneous round-trip times, multiple congested routers and uncertain routing topologies.

Appendix F: On Integrating Fluid Models with Packet Simulation

F1. Introduction

Networks, and the Internet in particular, have seen an exponential growth over the past several years. This growth is likely to continue for the foreseeable future, and understanding the behavior of such systems is critically important. A number of discrete event-driven simulators [1], [2], [3], [4] have been developed for this purpose. These simulators provide accurate information for every simulated packet. They provide an important tool for designing new protocols, improving existing protocols, and verifying new observations. However the simulation capabilities of these simulators have fallen far behind the scale of the Internet today and this gap is growing as the size and speed of the network is growing. As an alternative, fluid models have been proposed in recent works [5], [6], [7] to analyze the performance of networks. The fluid models can predict the behavior of large networks both accurately and efficiently. In [6], networks consisting of hundreds of routers and thousands of high bandwidth links supporting millions of flows can be simulated in minutes on a desktop PC, a feat that is unachievable by current discrete event-driven simulators. However the fluid model provides no detailed information regarding individual packets, and its application is hindered by this limitation.

A simulation method that combines the best features of the above fluid and packet-level approaches is desirable for several reasons. First, by efficiently simulating large networks and providing detailed information for selected individual traffic flows, it is possible to study the performance of communication protocols deployed at end hosts across a wide area high speed network. Research on existing and future communication protocols for the Internet can potentially benefit from such a simulation method. Second, it would permit the performance study of small edge networks, such as wireless networks that exchange traffic with wide area networks. A small edge wireless network would best be simulated as a packet-level network whereas the WAN would be more efficiently modeled as a fluid network. Finally, introducing non-TCP traffic into the fluid model provides an alternative method to study wide area traffic, which may lead to better models of real Internet traffic. Such models could be useful in creating more efficient simulation methods and studying a large class of network problems.

In this appendix, using the topology-aware fluid model presented in Appendix B (also [6]), we develop a hybrid simulation approach that consists of a set of packet flows and fluid flows. We present two models of the interaction between the packet flows and the fluid flows traversing a network. Algorithms are presented for deriving the behavior of the packets, e.g., end-to-end latency and loss. The simplest of these models ignores the effect that the packets have on the fluid network and simply uses the solutions to the fluid network model to determine the outcome of each packet. The second approach accounts for the load introduced by the packet flows by transforming them into fluid flows that feed the fluid network. The solution of this augmented network is then used to determine the outcomes of the packets. These approaches require one and two passes, respectively and will henceforth be referred to as the one pass and two pass interaction models. The following Figure shows an example of such a hybrid simulation.

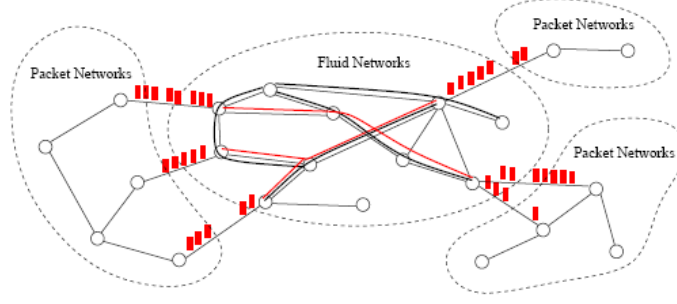


Figure 66: An example hybrid simulation

For both approaches, we establish the correctness of the resulting simulators and evaluate their accuracy and computational speedup through simulation. We have implemented both approaches in several popular simulators, including ns-2. We find that the one pass algorithm is fast and accurate provided that the offered packet load is very small compared to the load offered by fluid flows. Its accuracy degrades as the offered load due to the packet flow increases. On the other hand, the two-pass algorithm is very accurate, independent of the load introduced by the packet flows while incurring a slight slowdown. Henceforth, the word *simulator* refers to the discrete event-driven simulator used in the hybrid simulation unless further clarified.

Several works are particularly related to our study. In [8], a hybrid model is established for fast network simulation. The idea is to model network traffic as fluid flows and generate discrete events for packet losses within the network. [9] outlines the design of a fluid-oriented hybrid simulator admitting both packet flows and fluid flows. The hybrid simulator is implemented using the discrete event simulation framework. Recently, a dynamic simulation backplane was developed for creating distributed, component-based simulations of communication networks by interconnecting models of subnetworks drawn from different network simulation packages [10]. Based on this backplane, a multi-paradigm simulation framework, MAYA, is proposed to integrate three disparate modeling paradigms: discrete event models, analytical models and physical network interfaces. In particular, the fluid model proposed in [5] is integrated with the packet simulator Qualnet [11], [3].

F2. Traffic Interaction Models

In this section, we address the problem of modeling interactions between the foreground packet traffic and the background TCP traffic. When the foreground packet traffic traverses the fluid network, the delay and drop probabilities of the packets are determined by the queue lengths and drop probabilities of the queues in the fluid network, which are computed by the fluid model solver. On the other hand, the queue lengths and drop probabilities of these queues are affected also by the traversing foreground packet traffic, which then affect the background TCP traffic accordingly.

A hybrid model is built based on the fluid model. In the hybrid model, the network G serves not only a population of N classes of TCP flows, but also a population of M packet flows. A packet flow consists of a sequence of packets that share the same path in the fluid network. Each packet

flow i has a packet arrival process, $\lambda_i(t)$, which denotes the number of packets arriving at the fluid network by time t , $\{S_i\}$, which denotes the sequence of packet lengths associated with the flow, and a path in the fluid model defined as $E_i = (l_{i,1}, l_{i,2}, \dots, l_{i,h_i})$.

We present two algorithms for dealing with the traffic interaction problem. The first assumes that packet flows impose a load on the fluid network that can be neglected. The second accounts for the load that the packets impose on the fluid network. The first will require a single pass through time in order to determine outcomes for each packet (loss/no loss, end-end delay) and will be referred to as the one-pass traffic interaction model; whereas the second requires an additional preliminary pass to solve the fluid network accounting for the load introduced by the packet flows and will be referred to as the two-pass traffic interaction model.

A. One-pass Traffic Interaction Model

Since the one-pass traffic interaction model assumes that the packet flows have negligible effect on the background TCP traffic, the queue lengths and drop probabilities seen by the packets are identical to those predicted by the fluid model. Thus, the delay and drop probability experienced by a packet traversing through a fluid network are estimated using the queue lengths and drop probabilities derived directly from the fluid model in [6]. The delay and drop probability for each packet are computed cumulatively along its path in the fluid network. The packet is scheduled to depart from the fluid network according to this delay and drop probability.

The one-pass model is based on the assumption that the packet flows have negligible effect on the fluid network. This should be reasonable provided that the rate of the traversing foreground packet traffic is small enough. However, it is often the case that the foreground flows have considerable throughput such that the accuracy of the simulation results given by the one-pass model can be impaired. One extreme example is a hybrid simulation in which there is one background TCP class traversing two identical queues in the fluid network and there are foreground packets traversing the second queue. If the throughput of the foreground packet traffic is high enough, the second queue becomes the bottleneck. However, in the one-pass model, the second queue is always empty regardless of the rate of the traversing foreground packet traffic.

B. Two-pass Traffic Interaction Model

The two-pass traffic interaction model extends the one-pass model to account for the interaction between the packet flows and the background TCP traffic in the fluid network. In the two-pass model, the behavior of the fluid network is determined during a first pass and the effect on the packets is determined once they traverse the network during the second pass.

In the first pass of the two-pass model, the packet flows are transformed into foreground flows that can be incorporated into the solution of the fluid model. This transformation and the resulting solution of the model constitute the first pass. In the second pass, the delay and the drop probabilities of the traversing packets are estimated using the queue states obtained during the first pass.

F3. Synchronization in Hybrid Simulation

While most existing network simulators, [1], [2], [4] etc., are discrete event-driven simulators, the fluid model is a set of (coupled) ordinary differential equations of TCP network dynamics as functions of time [5], [6]. The states of the TCP network at any time can be directly obtained, in theory, by solving the set of differential equations. In practice, in order to save computational resources, the fluid model is solved incrementally as a function of time and all the network states prior to the current time in the fluid model can be obtained. For example, in [6], a fixed step-size Runge-Kutta algorithm is implemented to solve the fluid model; thus the fluid model is solved using a time-stepped network simulator. Thus, the simulator generating packets arriving to the fluid network and the fluid model solver are likely to have separate time management systems and, if integrated in the context of a hybrid simulation, they will have to be synchronized. For example, in a two-pass hybrid simulation, the fluid model takes packet traffic from the packet simulator. It is essential to have the packet traffic rate available in time to advance the fluid model. On the other hand, the packet simulator needs those packets traversing fluid network to be delivered by the fluid model in time.

Definition 1 (T_f): the simulation clock of the fluid model solver, meaning the states of the fluid network, e.g. queueing delay, packet loss, flow rate, etc., have been obtained for all $t < T_f$.

Definition 2 (T_s): the simulation clock of the simulator, meaning all the packet events before T_s have been processed by the simulator.

Definition 3 (τ): the minimal propagation delay of links that are the last hops on the packet's paths inside the fluid network.

Definition 4 (δ_s): smoothing interval, packet flow during each interval is modeled as a constant rate fluid.

The purpose of synchronization between the fluid model and packet simulator is to avoid out of order events. More specifically, we want to ensure that the fluid model never deliver a packet to the packet simulator with a timestamp $t_0 < T_s$. At the same time, the packet simulator must never inject a packet into the fluid network with timestamp $t_0 < T_f$. This places constraints on the rates that $\{T_s, T_f\}$ can advance.

Our approach for synchronizing the fluid model solver and the simulator aims at satisfying the synchronization requirements for the two-pass model, which is stricter than those for the one-pass model. So when the one-pass model is used in our implementation, its synchronization requirements are naturally satisfied.

In our implementation, the fluid model solver is synchronized to the simulator every smoothing interval δ_s which is chosen to be smaller than τ . At each time when $T_s = k\delta_s$, $k = 1, 2, \dots$, a synchronization event is scheduled to happen. During this event, the fluid model solver evolves from $T_f = (k - 1)\delta_s$ to $T_f = T_s = k\delta_s$. The fluid model schedules out all packets that are supposed to arrive at packet network before $T_f + \tau$. Since $\delta_s < \tau$, we can advance T_s to the next synchronization point $(k + 1)\delta_s$.

F4. Implementation

We have implemented the two traffic interaction models in ns-2[1], pdns[4] and the Backplane[10] by integrating the fixed step-size fluid model solver developed in [6]. In this section, we describe our implementation in the ns-2 network simulator as an example of how the hybrid simulation approach might be implemented.

The design objective of our implementation is to achieve maximum flexibility when dividing the network under simulation into fluid networks and packet networks and deploying foreground packet traffic in the network.

The Figure below on the left shows a network topology under hybrid simulation. Circled by the dotted oval is the fluid network simulated by the fluid model solver. The grey nodes in the fluid networks are points where packets enter or leave the fluid network. These nodes are duplicated in the form of ns node objects and serve as *access points* to the fluid networks. The Figure below on the right is the structure of the corresponding ns-2 network topology which consists of a fluid model solver and six ns node objects. The access points are connected to the fluid model solver via specially designed 'fluid links'. These fluid links are virtual links that have infinite service capacity and no propagation delay. Their only function is passing packets to the fluid model solver and vice versa. With this design, the network can be divided arbitrarily into fluid networks and packet networks and the foreground packet traffic can reach anywhere in the network.

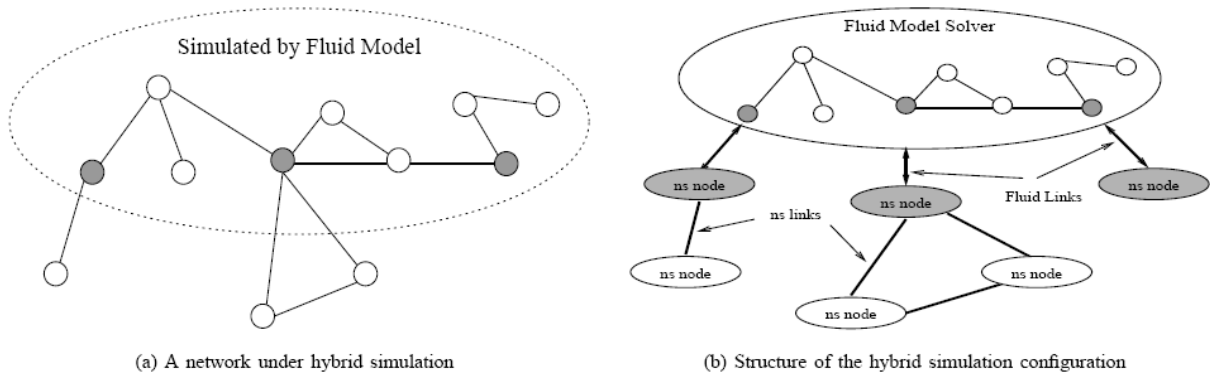


Figure 67: Fluid model in a hybrid simulation in ns-2

In our current implementation of the hybrid simulation, we assume that the paths between two access points in the fluid network are known and statically pre-configured. When a packet is routed to pass across the fluid network, it is first sent to one of the access points. The access point knows from its routing table that it should send the packet to the 'fluid link', which directly passes the packet to the fluid model solver and identifies the ingress point. The fluid model solver is an extension of an ns node object. After receiving a packet from the fluid link, the routing mechanism embedded in the ns node object is able to tell the next packet level node that the packet is going to arrive, which is also the egress point of the packet from the fluid network. Thus the two access points (the ingress point and the egress point) of the packet is known and so is the path of the packet in the fluid network. When the delay and the drop probability of the packet is computed, the packet is scheduled to arrive at the fluid link that

connects to the egress point ns node on the path of the packet. This procedure is shown in the following Figure.

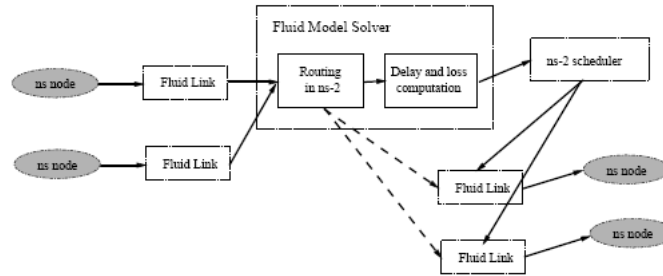


Figure 68: Connecting the fluid model and ns node objects

F5. Experimental Results

We have carried out extensive experiments to test the accuracy and performance of the hybrid simulation approach. The accuracy and efficiency are evaluated by comparing the simulation results given by the hybrid simulation with those obtained in a packet level simulation. Five sets of experiments are presented to answer the following questions:

- How should the smoothing interval length be set?
- How does the one-pass model perform in terms of its accuracy?
- How does the background TCP traffic in the fluid network interact with traversing unresponsive packet flows? These flows might be used to model video or audio flows running on top of UDP.
- How do the TCP flows simulated by the fluid model compete with TCP flows simulated by a packet source when they share the same bottleneck queue? The fluid model has been shown to be accurate in describing the behavior of TCP networks, so the question is to judge its accuracy when the flows interact with real packet TCP flows.
- Does the two-pass model hybrid simulation capture the interactions between multiple traversing foreground flows?
- Do hybrid simulations scale?

In the simulation, the smaller the smoothing interval, the more accurate the packet arrival rate is transformed into sending rate of the foreground flows in the fluid networks, which gives more accurate simulation results. On the other hand, smaller smoothing interval may cause more frequent interactions between the fluid model solver and the simulator. We test this potential cost that small interval may bring to the experiment. The result is that decreasing the smoothing interval doesn't bring any obvious increase in the simulation cost. This can be explained by the fact that, in our implementation, the extra work brought by decreasing the smoothing interval is a proportional increasing of synchronizing events and measuring the packet arrival rate.

The first part grows in reverse proportion to the smoothing interval, but its cost is negligible compared to the computation cost of the fluid model solver and packet simulation. The second part is a cheap division operation. In our case, the smoothing interval is set to be the same length as the fluid model solver step size since a smoothing interval smaller than the model solver step size is meaningless.

The first set of experiments will explore the accuracy of the one-pass model as we increase the number of packet flows sharing a bottleneck link with background TCP fluid flows.

The remaining experiments are all performed using the two-pass model. The second experiment observes a class of background TCP flows in a single bottleneck topology interacting with a UDP source whose sending rate changes with time. The accuracy is tested by comparing the delay and drops of the UDP traffic and the TCP network behavior given by the hybrid simulation and the packet simulation. Then, we present a set of simulation results showing the interaction between the background TCP traffic and different number of traversing foreground packet TCP flows. These results show that when sharing a bottleneck queue, the TCP flows modeled by the fluid model have the same competing capability for bandwidth as those simulated by the packet TCP sources, which further proves the correctness of both the hybrid simulation and the original fluid model. After that, simulation results of a hybrid simulation which consists of multiple foreground flows are given. In the last experiment, we show the capability of the hybrid simulation by simulating a network with more than 3,000 nodes and thousands of TCP classes consisting up to 12,378,340 TCP flows.

All the experiments are performed multiple times and yield similar results. In all the experiments, we use TCP Newreno and RED with ECN marking as the AQM policy. The step size of the fluid model solver is fixed at 1ms except for the last experiment where the step size is set to 5 ms.

F5.1. Accuracy of One-pass Model

The setting of the experiment is shown in the Figure below. The six queues on the top are simulated in the fluid model and the other four queues on the bottom are simulated in packet form. *B1* and *B2* are access points between the fluid network and the packet network. The queue between *B1* and *B2* has a capacity of 100 Mbps and a propagation delay of 10 ms and is the bottleneck. Other queues have a capacity of 200 Mbps and a 10 ms propagation delay. Class 0 and Class 1 are classes of TCP flows from node *S1* to *D1* and from node *S2* to *D2* respectively. Class 0 is the background TCP traffic simulated in the fluid model and Class 1 is the foreground packet traffic.

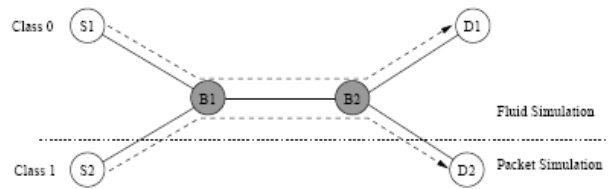


Figure 69: Network with single bottleneck

We perform the experiment 5 times. Each time there are 40 TCP flows going through the bottleneck queue and the 40 flows are divided between Class 0 and Class 1. In the k^{th} time experiment, Class 0 contains $40 - k$ TCP flows and Class 1 contains k TCP flows. The TCP flows start at time 0 and ends at time 100s. We measure the throughput of the flows from 30s to 90s. For comparison, we perform experiments with the same scenario in a packet level simulation.

Table I compares the throughput of the packet TCP flow(s) (Class 1) in a hybrid simulation with that given by the packet simulation. In all cases, the throughput given in the hybrid simulation is larger than that in the packet simulation. However, we can see that when the packet traffic is small, the one-pass model can be pretty accurate and in these cases, generally the error is under 10%. As the fraction of the packet traffic increases, there is an increasing trend in error.

k	1	2	3	4	5
Hybrid (Class 1 Throughput)	18309	36646	57331	75736	104654
Packet (Class 1 Throughput)	18203	35752	54307	71512	92441
Error	0.58%	2.50%	5.57%	5.90%	13.21%

TABLE I
ACCURACY OF ONE-PASS MODEL

We will come back to this same scenario in the coming experiments and show how the two-pass model performs.

F5.2. Interaction with UDP Traffic

This experiment is to test the accuracy of the hybrid simulation when simulating the interaction between the background TCP traffic and a UDP traffic whose sending rate changes with time.

The same network setting in the Figure above is used. This time, Class 0 is a class of 10 TCP flows from node $S1$ to $D1$ and Class 1 is UDP traffic from node $S2$ to $D2$. A 100s simulation is performed. The rate of the UDP traffic is 10 Mbps at time 0 and increases to 20 Mbps at 10s, 30 Mbps at 20s, ..., until 90 Mbps at 80s. The packet size is set to 1000 bytes.

The top plots in the Figure below compare the average delay and drop probability experienced by the UDP traffic in different simulation time slots given by the hybrid simulation and the packet simulation. In both cases, they show a good match. This implies that the hybrid simulation is able to predict the behavior of a UDP foreground flow when it traverses the background TCP traffic. The bottom two plots in the Figure below show the behavior of the TCP window size for Class 0 and the queue behavior of the bottleneck link. In the hybrid simulation, these results are given by the fluid model, which predicts the average behavior. We can also see a close match in these results. This accuracy of the fluid model predicting the average behavior is also reflected by the correct estimation of the average delay and drop probability that the UDP traffic experience. It shows that while the hybrid simulation can provide accurate packet level information for foreground packet flows, it can also estimate the average behavior for background TCP traffic and queues in the fluid network.

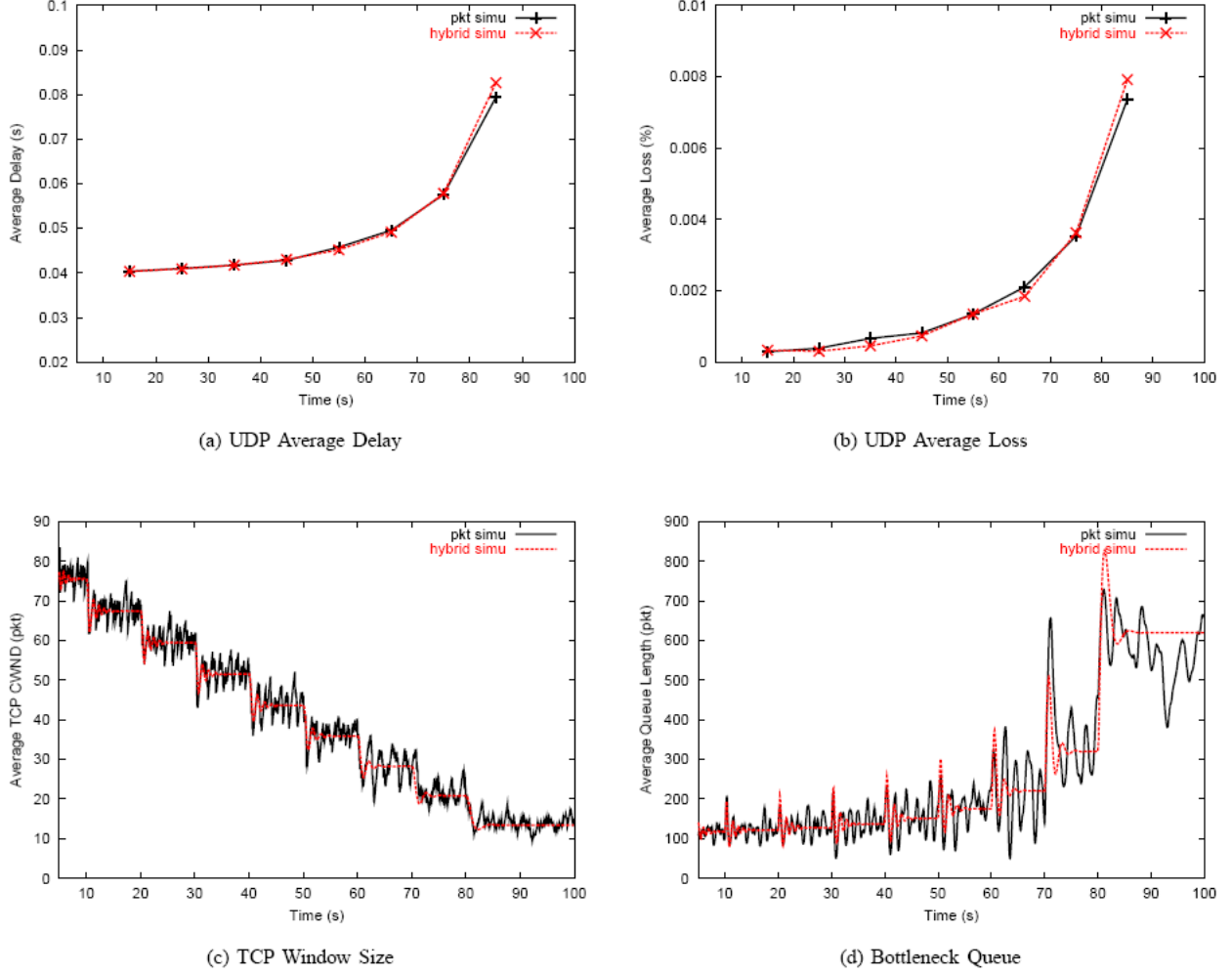


Figure 70: Hybrid simulation between the fluid model and UDP traffic

F5.3. Interaction with TCP Traffic

We come back to the first experiment here. Now, we illustrate the accuracy of the two-pass model with the same setting as in the first experiment except that the packet traffic in this experiment takes more fraction of the total TCP flows.

The experiment is carried out 10 times. Each time there are 40 TCP flows going through the bottleneck queue and the 40 flows are divided between Class 0 and Class 1. In the k^{th} time experiment, Class 0 contains $4 * (11 - k)$ TCP flows and Class 1 contains $4 * (k - 1)$ TCP flows. So the fraction of the packet traffic increases proportionally to k and the experiment is a fluid model simulation at the first time.

We record the throughput of the bottleneck from 30 s to 90 s in simulation and compare the computation time of each simulation. These experiment results are shown in the top two plots in the Figure below and Table II. In Table II, the first row shows the throughput of Class 0, which is the TCP class simulated by the fluid model. The second row shows the throughput of Class 1, which are TCP flows simulated by the packet simulator. The third row is the total

throughput of the two and the forth row is the percentage of the packet throughput in the total throughput. As a reference, the simulation results given by a packet simulation are shown at place where 100% of the flows are packet flows. The bottom three plots in the Figure below show the average TCP window sizes of class 0 and class 1 and the bottleneck queue behavior in both hybrid simulation and peer packet simulation when both Class 0 and Class 1 contain 50% traffic (20 TCP flows) and evenly share the bottleneck service capacity. In (c) and (e) below, the comparison is between solutions of the fluid model used in the hybrid simulation and those given by a packet simulation.

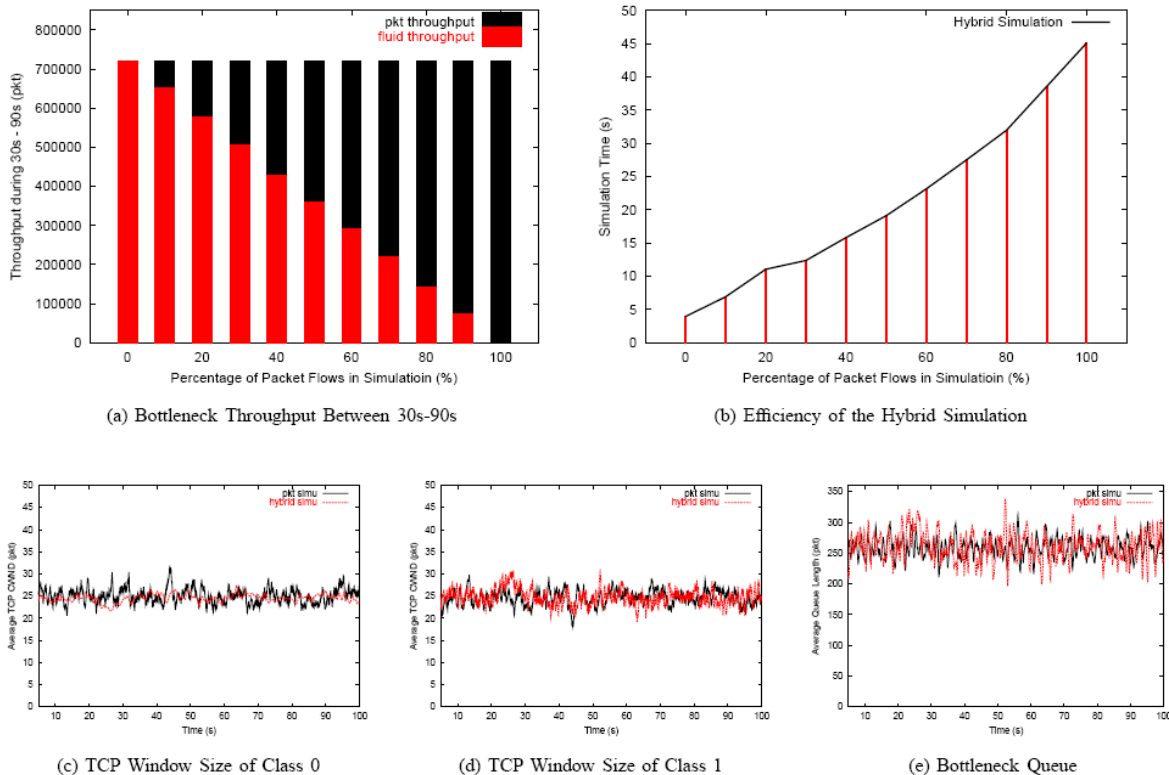


Figure 71: Interaction between fluid TCP flows and packet TCP flows

k	1	2	3	4	5	6	7	8	9	10	pkt
Fluid	721153	652906	577602	505252	428611	361022	290473	218446	142799	72899	0
Packet	0	68211	143526	215763	292512	360166	430703	502700	578332	648228	721168
Total	721153	721127	721128	721015	721173	721188	721176	721146	721131	721127	721168
Pkt/Tot(%)	0	9	20	30	41	50	60	70	80	90	100

TABLE II
INTERACTION BETWEEN FLUID TCP FLOWS AND PACKET TCP FLOWS

From these results, we see that the percentage of the packet throughput increases in proportion to its percentage of the total number of TCP flows, which strongly supports that the TCP flows simulated by the fluid model have the same behavior as those simulated by the packet source when sharing a bottleneck queue in a hybrid simulation. This further proves the correctness of the original fluid model and the hybrid simulation's ability to simulate the interaction between TCP flows in both fluid form and packet form. Also from this experiment, we can see the performance advantage brought by the fluid model from (b) in the Figure above.

Compared with a packet simulation, the speedup of this hybrid simulation depends on the amount of the packet traffic in the hybrid simulation. If C_{pkt} denotes the packet simulation cost, C_{hf} denotes the fluid model solver computation cost in a hybrid simulation and C_{hp} denotes the simulation cost for the packet traffic in the hybrid simulation, there is

$$Speedup = \frac{C_{pkt}}{C_{hf} + C_{hp}}.$$

, and as the simulation cost of the packet traffic in a hybrid simulation with small network topology is almost the same as that in the packet simulation, the fraction of the packet traffic decides the upper bound of the speedup. For example, if 10% of the traffic is packet traffic, the speedup can go up to at most 10. In this experiment, we see that when 10% of the traffic is packet traffic, the hybrid simulation reaches a speedup of 6.53 compared to the packet simulation.

F5.4. Multiple Foreground Flows

In this experiment, we show that the two-pass model hybrid simulation captures the additional interaction between multiple traversing foreground flows.

The network under simulation is shown in the Figure below. Node S2, D2, S3 and D3 are four packet nodes connected to the fluid network via different access points. Class 0 and Class 1 are both classes of 10 TCP flows simulated in the fluid model. Class 2 is a class of 20 TCP flows. Class 3 is a class of 40 TCP flows. Class 0, 1 and 2 start at time 0. Class 3 is only active from 30s to 60s.

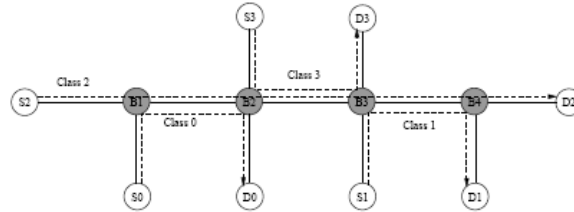


Figure 72: Network with three bottlenecks

The Figure below shows the matching simulation results between the hybrid simulation and the packet simulation. At 30s, the average TCP window size of Class 2 decreases because of the new traffic brought by Class 3, which is another foreground flow, and the queue behavior at each bottleneck also changes accordingly. At 60s, the network behavior changes back because Class 3 stops sending traffic.

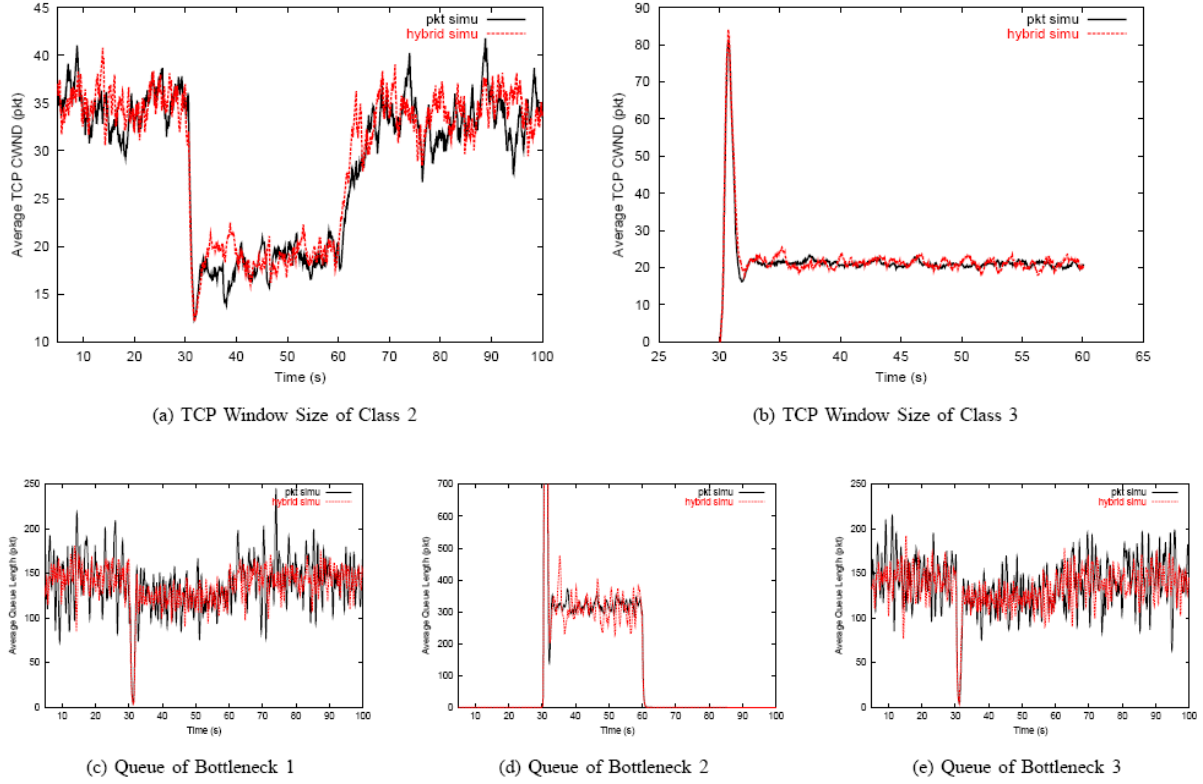


Figure 73: Hybrid simulation with multiple foreground flows

F5.5 Experience with Large IP Networks

In this section, we show the ability of the hybrid simulation to simulate a large IP network. We use the Inet Topology Generator from University of Michigan [14] and generate a network of 3500 nodes and 11334 links, each with a capacity of 2.5 Gbps. Then we randomly create 5000 classes of TCP flows which is a total of 12,378,340 TCP flows in the network. Then we randomly pick two nodes as the source and the destination of a UDP traffic. We set the UDP traffic as a CBR at 50 kbps with a packet size of 50 bytes. A 100 s simulation is performed. Our experiment is carried on a Dell Precision Workstation 530, which is configured with two Pentium IV processors (2.2 GHz) and 2 GB memory. Since our program is not parallelized, only one processor is utilized. We have taken several trials. On average, the simulation takes about 30 minutes and 659M memory. The Figure below shows the time sequence of the UDP packets between time 74 s and 75 s in one of the experiments. This UDP traffic has an average delay of 101.66ms and an average drop probability of 15.70%.

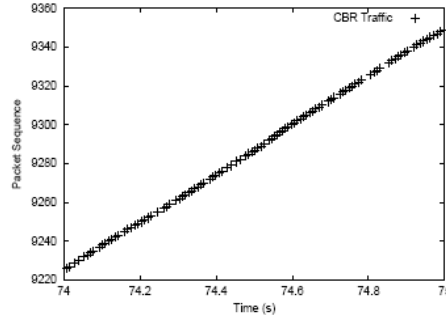


Figure 74: Time sequence of the UDP traffic between 74 s and 75 s

F6. Conclusion & Future Work

We have developed a simulation method that takes the advantage of the highly efficient fluid model and at the same time provides detailed information at packet level for selected individual traffic flows. We achieve this by an effort to simulate the network using fluid model solvers and discrete event-driven simulators, the hybrid simulation. Two models are proposed to describe the interactions between the background TCP traffic and the foreground packet traffic. Synchronization between the fluid model solver and the simulator are analyzed. Simulation results show that our method maintains the performance advantage of the fluid model and generates accurate simulation results comparable to those given by the discrete event-driven simulator. Our work can be reached and downloaded at <http://www-net.cs.umass.edu/fluid/ffm.html>.

As a future work, we will implement our hybrid simulation approach in more simulators and at the same time, incorporate new features of the fluid model as well as other network traffic models. Also, our current implementation in ns-2 could be further improved, such as supporting more functions and increasing its performance. Another exciting future work direction is to further boost the speed of the hybrid simulation such that the hybrid simulation can generate real time traffic information. This information can be used in network emulators to provide more controlled and more realistic delay and loss reference. Our current approach can handle real time traffic to some extent. Further performance boost can be reached by parallelization, for which the time-stepped nature of the fluid model solver is perfectly suitable.