

Role Based Trust Management Security Policy Analysis

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Abstract

We use symmetric keys to encrypt and decrypt attribute values. These keys are distributed only to the brokers that are trusted with the attribute values. The system will never deliver these keys to clients. This reduces the number of nodes that are trusted with sensitive keys, and that take part in key management protocols. This does not affect security since local brokers encrypt and decrypt attribute values on behalf of connected clients, and deliver events to over secure links. Τo support cryptographic properties such as key freshness, and forward and backward secrecy [22], the system requires key management service(s). A representation of quantified trust relationships, the trust dependency graph, and a sample QuanTM application specific to the KeyNote trust management language, are also proposed.

Keywords- Symmetric keys, Trust Management

I. INTRODUCTION

Creating software which is flexible and highly customizable to adapt to fast changing business needs has moved into the main focus of software developers. Enterprises demand a seamless communication between applications independent from the platform on which they run and even boundaries. Service-oriented domain Architectures and XML Web Services have been designed to meet these concerns, allowing a flexible integration of services provided by independent business partners. However, the seamless and straightforward integration of cross-organisational services conflicts with the need to secure and control access to these services. The traditional approach to restrict service access is based on user authentication performed by the service provider itself, cf. [18]. Since credentials (e.g. user name and password) needed to access a service are issued and managed by the service provider, this approach is referred to as isolated identity management as stated in [13]. It requires service users to register a digital identity at each involved service provider and to authenticate separately for each service access. Federated Identity Management as a new identity model provides solutions for these problems by enabling the propagation of identity information to services located in different trust domains. It enables service users to access all services in a federation using the same identification data. Several frameworks and standards for Federated Identity

Management have been specified (e.g. WS-Federation [1] and Liberty Identity Web Services Framework (ID-WSF) 2.0 [31]). The key concept in a federation is the establishment of trust whereby all parties in a federation are willing to rely on asserted claims about a digital identity such as SAML assertions [24]. As Service-oriented Architectures move from an isolated identity management scheme to a federated identity management, service providers are exposed to new risks. In a federation the authentication of a user is not necessarily performed within the service provider's domain, but can be done within the user's local domain. Consequently, the service provider has to trust the authentication performed by the user's identity provider. In terms of security this is a critical situation since authorization and access control of service are highly dependent on the authentication results. A weak authentication jeopardises the dependent service's security by increasing the risk that a user can personate as someone else and gain improper access. OASIS considers this as a serious risk [23] and recommends to agree on a common trust level in terms of policies, procedures and responsibilities to ensure that a relying party can trust the processes and methods used by the identity provider. Jøsang et. al. [13] describe the usage of such a common trust level as a symmetric trust relationship, since all parties are exposed to an equal risk in the case of failure. As opposed to this, having different trust requirements and mechanisms is referred to as an asymmetric trust relationship. They argue that asymmetric trust relationships are hard to establish, since the parties are exposed to different risks in the case of failure. However, with regard to complex SOA – that might be based on the dynamic selection of services and service providers - defining and enforcing a common trust level is disadvantageous: symmetric trust relationship between the providers in a federation would require a trust level, which is sufficient for the service with the strongest authentication requirements. These requirements, however, might not be necessary for all services within the federation and might change if this service is dynamically replaced. Consequently, users are forced to authenticate by a predefined strong even though authentication method, authentication would be sufficient for the service they want to access. Likewise, when users are fixed to a predefined authentication method according to the specified trust level, access will be denied even though the user might be able to verify his identity

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in an even more trusted way. Altogether, there is a demand for more flexibility authentication processes in SOA. To achieve this flexibility, a way to rate the trust relationship between identity provider and service provider is needed in order to restrict the service access based on an individual trust level. The general idea of classifying authentication methods according to their level of trustworthiness is not new. Especially in the field of e-Government, various countries have launched e-authentication initiatives in order to secure access to critical e-Government services [26. 11, 17, 5]. All of these initiatives have in common that they define authentication trust levels – mostly four different levels - in a way that covers the main use cases, reaching from "no security needed" to "critical application". For each level, requirements for the authentication process are defined. This means, authentication methods are always assigned to predefined levels, but not the other way around. To provide authentication in a truly flexible manner, we present in this paper:

- A formal definition of trust levels to quantify the trust that is established by using a particular authentication method. This definition is globally applicable and not restricted to a specific use case setting requiring specific bootstrapping algorithms. This way, the meaning of a trust level based on our approach is clear and can be applied to any use case without the need to know any further set up or environment parameters.
- A mathematical model to combine different authentication methods as used in a two-factor authentication and to calculate their combined authentication trust level.
- An example calculation that demonstrates the applicability of our mathematical model to existing authentication methods.

This paper is organized as follows. Section 2 provides an overview about related work and current efforts in this area. In Section 3 we present our approach for assessing and quantifying trust in authentication methods. This section gives a definition for an authentication trust level and shows how this level can be determined. Section 4 introduces a mathematical model to calculate the trust value for the combination of two authentication methods taking into account the similarity of two mechanisms. To demonstrate the effect of the similarity on the combined trust level, an example calculation is presented in Section 5. Finally, Section 6 concludes this paper and highlights some The emergence of distributed future work. topologies and networked services has resulted in applications that are stored, maintained, and accessed remotely via a client/server model. The advantages of such a setup are many, but the challenges of access control and identity management must be addressed. Trust management and reputation management are two differing approaches to the problem. While effective with regard to explicit declarations, trust management applicability when relationships characterized by uncertainty. management is useful in enforcing existing trust relationships but ineffective in the formation of partially trusted ones. Reputation management provides a means of quantifying trust relationships dynamically, but lacks access enforcement and delegation mechanisms. To address this divide we introduce the notion of Quantitative Trust Management (QTM), an approach that merges concepts from trust and reputation management. It (QTM) creates a method for specifying both policy and reputation for dynamic decision making in access control settings. A system built upon QTM can not only enforce delegated authorizations but also adapt its policy as partial information becomes more complete. The output is a quantitative trust value that expresses how much a policy-based decision should be trusted given the reputations of the entities involved. Further, to make this novel concept concrete, we propose QuanTM, an architecture for supporting QTM. In this application of QuanTM, we use the KeyNote [8, 7] (KN) trust management language and specification, due to its well defined delegation logic and compliance system. Summarily, a KN evaluator checks a user's access credentials against local policy to produce a compliance value from a finite and predefined set of values. The compliance value is then used to make access decisions. KN allows principals to delegate access rights to other principals without affecting the resulting compliance value. Further, KN is monotonic: If a given request evaluates to some compliance value, adding more credentials or delegations will not lower that value. We argue that credentials should not be explicitly trusted, nor should the trustworthiness of delegating principals be ignored. Furthermore, the result of evaluation for a given access request may need to be dynamic [9]. Service providers may find it desirable to arrive at different opinions based on local constraints, policies, and principals for the same request. In QuanTM, this is easily expressed. We address these issues in the following two ways: (1) It includes a means to dynamically assign reputation to principals and their relationships within a request, and (2) It provides a mechanism for combining this information to produce a trust value. In QuanTM, a trust value (often a real number) is used to represent the the trustworthiness of a given compliance value and how it was reached. Our proposed QuanTM architecture (see Fig. 1) consists of three subsystems

1. Trust management consists of a trust language evaluator that verifies requests meet policy constraints, and a trust dependency graph (TDG) extractor that constructs a graph representing trust relationships.



- 2. Reputation management consists of two modules. First, a reputation algorithm to dynamically produce reputation values by combining feedback. These reputation values weigh TDG edges. Second, a reputation quantifier computes the trust value for a given request by evaluating the weighted TDG.
- 3. Decision management is composed of a decision maker that arrives at an access determination based on a trust

value, context, and an application specific metapolicy that encodes a cost-benefit analysis. The design of QuanTM has been guided by the requirement that the individual components will be application specific, and thus, we have designed QuanTM modularly. QuanTM provides a simple interface by which different trust management languages, reputation algorithms, and decision procedures may be included. In this paper, we propose a QuanTM design instance that utilizes the KeyNote language and TNA-SL [11, 12] reputation algorithm. This instance's implementation and evaluation is the subject of future work.

A. Background

Several approaches to define levels of trustworthiness for authentication mechanisms have been proposed in recent years indicating the importance of such a concept. In the area of e-Government, the UK Office of the e-Envoy has published a document called "Registration and Authentication e-Government Framework Policy and Guideline" [26]. In this document the initial registration process of a person with the system as well as the authentication process for a user's engagement in an e-Government transaction are defined. Depending on the severity of consequences that might arise from unauthorized access, four authentication trust levels are defined, reaching from Level 0 for minimal damage up to Level 3 for substantial damage. The IDABC [11] (Interoperable Delivery of European eGovernment Services to public Administrations, Businesses and Citizens) is a similar project managed by the publishes European Commission. It recommendations and develops common solutions in order to improve the electronic communication within the public sector. Its Authentication Policy Document [7] defines four assurance levels as well, which are also associated with the potential damage that could be caused. For each of the four levels the document defines the requirements for the registration phase and for the electronic authentication. The e-Authentication Initiative is a major project of the e-Government program of the US. The core concept is a federated architecture with multiple e-Government applications and credential providers. The intention is that the e-Authentication Initiative provides an architecture which delivers a uniform, government-wide approach for authentication while leaving the choice of concrete authentication technologies with the

individual government agencies. In this context, the initiative has published a policy called "EAuthentication Guidance for Federal Agencies" [5] to assist agencies in determing the appropriate of identity assurance for transactions. The document defines four assurance levels, which are based on the risks associated with authentication error. Which technical requirements apply for each assurance level is described in a recommendation of the National Institute of Standards and Technology (NIST). which is called

II. ROLE BASED TRUST POLICY ANALYSIS

Policy analysis [12] as we consider it here examines whether the specified relationships between roles hold in all reachable policy states. We explain reachable policy states below. The relationships, called queries, are set containments and take the form % w in which % and are each either roles or explicit (constant) sets of principals. For instance, X.u w A.r holds if every member of A.r is a member of X.u in every reachable policy state P0, i.e., [[A.r]]P0 $_$ [[X.u]]P0 . Queries of this form can be used to express many important security properties such as availability, safety, liveness and mutual exclusion. For instance, a safety property might be that everyone in the role that has access to the secret database is in the employee role. In general, any policy state can evolve into any other policy state by having principals issue new policy statements and revoke old ones. In security analysis we ask whether queries hold in all policy states that differ from a given current policy state only by changes to roles outside some trusted set. Intuitively, this corresponds to the fact that we expect certain principals to cooperate with us in our goal of preserving certain desired security properties. Specifically we assume that to this end these principals agree not to add or remove statements defining certain roles that they control. Other roles are not be assumed to be managed in cooperation with our goals. This intuition leads [12] to the defined two sets of roles that are used to determine the reachable policy states, the set of growth-restricted roles GR and the set of shrinkrestricted roles SR. Such a pair is called a restriction rule and is denoted by R = (GR, SR). Growthrestricted roles (GR) are not allowed to have new statements defining them added to the state. Shrinkrestricted roles (SR) are not allowed to have statements defining them removed. We write P 7!R P0 to indicate that P0 GR P and P0 P SR. It is important to note that these restrictions are not actually enforced. They are simply assumptions under which the analysis is performed. Their presence enables the analysis to provide us with assurances of things like, "So long as the people I trust do not make policy changes without first

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running the analysis, only company employees will be able to access the secret database." Queries of certain restricted forms can be analyzed and verified in polynomial time. These include queries in which at least one of % and _ is an explicit set. They also include situations in which only Type I and Type II statements are allowed in the policy state. However, when both % and _ are roles and all forms of statements are allowed, the decision problem is EXPTIME complete [20]. This is unfortunate because such properties are extremely useful. For instance, suppose we want to determine whether only employees could ever get access to a company's secret database. This can be determined efficiently if the set of employees is enumerated explicitly in the query. However, this does not consider the effect of employee turnover. By identifying employees and those users with access to the database both as roles, we can determine whether the desired property will continue to hold as new employees are added. Thus we seek techniques that can solve queries of this general form as often as possible.

We consider the following anycast field equations defined over an open bounded piece of network and /or feature space $\Omega \subset R^d$. They describe the dynamics of the mean anycast of each of p node populations.

$$\begin{cases} (\frac{d}{dt} + l_i)V_i(t,r) = \sum_{j=1}^p \int_{\Omega} J_{ij}(r,r)S[(V_j(t-\tau_{ij}(r,r),r)-h_{|j})]dr \\ + I_i^{ext}(r,t), & t \ge 0, 1 \le i \le p, \\ V_i(t,r) = \phi_i(t,r) & t \in [-T,0] \end{cases}$$
 (1)

We give an interpretation of the various parameters and functions that appear in (1), Ω is finite piece of nodes and/or feature space and is represented as an open bounded set of R^d . The vector r and r represent points in Ω . The function $S:R \to (0,1)$ is the normalized sigmoid function:

$$S(z) = \frac{1}{1 + e^{-z}} \tag{2}$$

It describes the relation between the input rate v_i of population i as a function of the packets potential, for example, $V_i = v_i = S[\sigma_i(V_i - h_i)]$. We note V the p-dimensional vector $(V_1, ..., V_p)$. The p function $\phi_i, i = 1, ..., p$, represent the initial conditions, see below. We note ϕ the p-dimensional vector $(\phi_1, ..., \phi_p)$. The p function $I_i^{ext}, i = 1, ..., p$, represent external factors from other network areas. We note I_i^{ext} the I_i^{ext}

dimensional vector $(I_1^{ext},...,I_n^{ext})$. The $p \times p$ matrix of functions $J = \{J_{ii}\}_{i,i=1,\dots,p}$ represents the connectivity between populations i and j, see below. The *p* real values h_i , i = 1,..., p, determine the threshold of activity for each population, that is, the value of the nodes potential corresponding to 50% of the maximal activity. The p real positive values σ_i , i = 1,...,p, determine the slopes of the sigmoids at the origin. Finally the p real positive values l_i , i = 1, ..., p, determine the speed at which each anycast node potential decreases exponentially toward its real value. We also introduce the function $S: \mathbb{R}^p \to \mathbb{R}^p$, defined $S(x) = [S(\sigma_1(x_1 - h_1)), ..., S(\sigma_p - h_p))],$ diagonal $p \times p$ and $L_0 = diag(l_1,...,l_n)$. Is the intrinsic dynamics of the population given by the linear response of data transfer. $(\frac{d}{dt} + l_i)$ is replaced by $(\frac{d}{dt} + l_i)^2$ to use the alpha function response. We use $(\frac{d}{dt} + l_i)$ for simplicity although our analysis applies to more general intrinsic dynamics. For the sake, of generality, the propagation delays are not assumed to be identical for all populations, hence they are described by a matrix $\tau(r,r)$ whose element $\tau_{ii}(r,r)$ is the propagation delay between population i at r and population i at r. The reason for this assumption is that it is still unclear from anycast if propagation delays are independent of the populations. We assume for technical reasons that τ is continuous, that is $\tau \in C^0(\overline{\Omega}^2, R_+^{p \times p})$. Moreover packet data indicate that τ is not a symmetric function i.e., $\tau_{ii}(r,r) \neq \tau_{ii}(r,r)$, thus no assumption is made about this symmetry unless otherwise stated. In order to compute the righthand side of (1), we need to know the node potential

$$\tau_m = \max_{i,j(r,r\in\overline{\Omega}\times\overline{\Omega})} \tau_{i,j}(r,r)$$
 (3)

factor V on interval [-T,0]. The value of T is

obtained by considering the maximal delay:

A. Mathematical Framework

A convenient functional setting for the non-delayed packet field equations is to use the space $F=L^2(\Omega,R^p)$ which is a Hilbert space endowed with the usual inner product:



$$\left\langle V,U\right\rangle _{F}=\sum_{i=1}^{p}\int_{\Omega}V_{i}(r)U_{i}(r)dr$$
 (1)

To give a meaning to (1), we defined the history space $C = C^0([-\tau_m, 0], F)$ with

 $\|\phi\| = \sup_{t \in [-\tau_m,0]} \|\phi(t)\| F$, which is the Banach phase space associated with equation (3). Using the notation $V_t(\theta) = V(t+\theta), \theta \in [-\tau_m,0]$, we write (1) as

$$\begin{cases} V(t) = -L_0 V(t) + L_1 S(V_t) + I^{ext}(t), \\ V_0 = \phi \in C, \end{cases}$$
 (2)

Where

$$\begin{cases} L_1: C \to F, \\ \phi \to \int_{\Omega} J(., r) \phi(r, -\tau(., r)) dr \end{cases}$$

Is the linear continuous operator satisfying $\|L_1\| \leq \|J\|_{L^2(\Omega^2,R^{p\times p})}$. Notice that most of the papers on this subject assume Ω infinite, hence requiring $\tau_m = \infty$.

Proposition 1.0 If the following assumptions are satisfied.

1.
$$J \in L^2(\Omega^2, \mathbb{R}^{p \times p}),$$

2. The external current
$$I^{ext} \in C^0(R, F)$$
,

3.
$$\tau \in C^0(\overline{\Omega^2}, R_+^{p \times p}), \sup_{\overline{\Omega^2}} \tau \le \tau_m$$
.

Then for any $\phi \in C$, there exists a unique solution $V \in C^1([0,\infty), F) \cap C^0([-\tau_m,\infty,F))$ to (3)

Notice that this result gives existence on R_+ , finite-time explosion is impossible for this delayed differential equation. Nevertheless, a particular solution could grow indefinitely, we now prove that this cannot happen.

B. Boundedness of Solutions

A valid model of neural networks should only feature bounded packet node potentials.

Theorem 1.0 All the trajectories are ultimately bounded by the same constant R if $I \equiv \max_{t \in R^+} \left\| I^{ext}(t) \right\|_F < \infty$.

Proof :Let us defined $f: R \times C \to R^+$ as $f(t, V_t) \stackrel{def}{=} \left\langle -L_0 V_t(0) + L_1 S(V_t) + I^{ext}(t), V(t) \right\rangle_F = \frac{1}{2} \frac{d \left\| V \right\|_F^2}{dt}$

We note $l = \min_{i=1,\dots,p} l_i$

$$f(t,V_t) \le -l \|V(t)\|_F^2 + (\sqrt{p|\Omega|} \|J\|_F + I) \|V(t)\|_F$$

Thus, if

$$\left\|V(t)\right\|_{F} \geq 2\frac{\sqrt{p\left|\Omega\right|}.\left\|J\right\|_{F}+I}{l} \stackrel{def}{=} R, f(t, V_{t}) \leq -\frac{lR^{2}}{2} \stackrel{def}{=} -\delta < 0$$

Let us show that the open route of F of center 0 and radius R, B_R , is stable under the dynamics of equation. We know that V(t) is defined for all $t \ge 0s$ and that f < 0 on ∂B_R , the boundary of B_R . We consider three cases for the initial condition V_0 . If $||V_0||_C < R$ and set $T = \sup\{t \mid \forall s \in [0, t], V(s) \in B_{\mathbb{R}}\}.$ Suppose that $T \in \mathbb{R}$, then V(T) is defined and belongs to $\overline{B_R}$, the closure of B_R , because $\overline{B_R}$ is closed, in $\partial B_{\scriptscriptstyle P}$, we $\frac{d}{dt} \|V\|_F^2 \Big|_{t=T} = f(T, V_T) \le -\delta < 0$ $V(T) \in \partial B_R$. Thus we deduce that for $\varepsilon > 0$ and small enough, $V(T+\varepsilon) \in B_R$ which contradicts the definition of T. Thus $T \notin R$ and $\overline{B_R}$ is stable.

Because f<0 on $\partial B_R, V(0) \in \partial B_R$ implies that $\forall t>0, V(t) \in B_R$. Finally we consider the case $V(0) \in C\overline{B_R}$. Suppose that $\forall t>0, V(t) \not\in \overline{B_R}$, then

$$\forall t>0, \frac{d}{dt}\|V\|_F^2 \leq -2\delta, \quad \text{thus} \quad \|V(t)\|_F \quad \text{is}$$
 monotonically decreasing and reaches the value of R in finite time when $V(t)$ reaches ∂B_R . This contradicts our assumption. Thus $\exists T>0 \,|\, V(T)\in B_R$.

Proposition 1.1: Let s and t be measured simple functions on X. for $E \in M$, define

$$\phi(E) = \int_{E} s \, d\mu \qquad (1)$$
Then ϕ is a measure on M .
$$\int_{X} (s+t) d\mu = \int_{X} s \, d\mu + \int_{X} t d\mu \qquad (2)$$

Proof: If s and if E_1, E_2, \ldots are disjoint members of M whose union is E, the countable additivity of μ shows that



$$\phi(E) = \sum_{i=1}^{n} \alpha_i \mu(A_i \cap E) = \sum_{i=1}^{n} \alpha_i \sum_{r=1}^{\infty} \mu(A_i \cap E_r)$$
$$= \sum_{r=1}^{\infty} \sum_{i=1}^{n} \alpha_i \mu(A_i \cap E_r) = \sum_{r=1}^{\infty} \phi(E_r)$$

Also, $\varphi(\phi) = 0$, so that φ is not identically ∞ . Next, let s be as before, let $\beta_1, ..., \beta_m$ be the distinct values of t,and let $B_j = \{x : t(x) = \beta_j\}$ If $E_{ij} = A_i \cap B_j$, the $\int_{E_{ij}} (s+t) d\mu = (\alpha_i + \beta_j) \mu(E_{ij})$ and $\int_{E_i} s d\mu + \int_{E_i} t d\mu = \alpha_i \mu(E_{ij}) + \beta_j \mu(E_{ij})$

Thus (2) holds with E_{ij} in place of X. Since X is the disjoint union of the sets E_{ij} $(1 \le i \le n, 1 \le j \le m)$, the first half of our proposition implies that (2) holds.

Theorem 1.1: If K is a compact set in the plane whose complement is connected, if f is a continuous complex function on K which is holomorphic in the interior of, and if $\varepsilon > 0$, then there exists a polynomial P such that $|f(z) = P(z)| < \varepsilon$ for all $z \varepsilon K$. If the interior of K is empty, then part of the hypothesis is vacuously satisfied, and the conclusion holds for every $f \varepsilon C(K)$. Note that K need to be connected.

Proof: By Tietze's theorem, f can be extended to a continuous function in the plane, with compact support. We fix one such extension and denote it again by f. For any $\delta > 0$, let $\omega(\delta)$ be the supremum of the numbers $\left| f(z_2) - f(z_1) \right|$ Where z_1 and z_2 are subject to the condition $\left| z_2 - z_1 \right| \leq \delta$. Since f is uniformly continous, we have $\lim_{\delta \to 0} \omega(\delta) = 0$ (1) From now on, δ will be fixed. We shall prove that there is a polynomial P such that

$$|f(z)-P(z)| < 10,000 \omega(\delta) \quad (z \in K)$$
 (2)

By (1), this proves the theorem. Our first objective is the construction of a function $\Phi \mathcal{E} C_c(R^2)$, such that for all z

$$|f(z) - \Phi(z)| \le \omega(\delta),$$
 (3)

$$\left| (\partial \Phi)(z) \right| < \frac{2\omega(\delta)}{\delta},$$
 (4)

And

$$\Phi(z) = -\frac{1}{\pi} \iint_{\mathcal{X}} \frac{(\partial \Phi)(\zeta)}{\zeta - z} d\zeta d\eta \qquad (\zeta = \xi + i\eta), \tag{5}$$

Where X is the set of all points in the support of Φ whose distance from the complement of K does not δ . (Thus X contains no point which is "far within" K.) We construct Φ as the convolution of f with a smoothing function A. Put a(r) = 0 if $r > \delta$, put

$$a(r) = \frac{3}{\pi \delta^2} (1 - \frac{r^2}{\delta^2})^2$$
 $(0 \le r \le \delta),$ (6)

And define

$$A(z) = a(|z|) \tag{7}$$

For all complex z . It is clear that $A \varepsilon C_c(R^2)$. We claim that

$$\iint_{\mathbb{R}^s} A = 1,\tag{8}$$

$$\iint_{\mathbb{R}^2} \partial A = 0, \tag{9}$$

$$\iint_{\rho^3} \left| \partial A \right| = \frac{24}{15\delta} < \frac{2}{\delta},\tag{10}$$

The constants are so adjusted in (6) that (8) holds. (Compute the integral in polar coordinates), (9) holds simply because A has compact support. To compute (10), express ∂A in polar coordinates, and note that $\partial A/\partial \theta = 0$,

$$\partial A/\partial r = -a'$$
,

Now define

$$\Phi(z) = \iint_{\mathbb{R}^2} f(z - \zeta) A d\xi d\eta = \iint_{\mathbb{R}^2} A(z - \zeta) f(\zeta) d\xi d\eta$$
 (11)

Since f and A have compact support, so does Φ . Since

$$\Phi(z) - f(z)$$

$$= \iint_{\mathbb{R}^2} [f(z - \zeta) - f(z)] A(\xi) d\xi d\eta \quad (12)$$

And $A(\zeta) = 0$ if $|\zeta| > \delta$, (3) follows from (8). The difference quotients of A converge boundedly to the corresponding partial derivatives, since

 $A\varepsilon C_c(R^2)$. Hence the last expression in (11) may



be differentiated under the integral sign, and we obtain

$$(\partial \Phi)(z) = \iint_{R^2} (\overline{\partial A})(z - \zeta) f(\zeta) d\xi d\eta$$

$$= \iint_{R^2} f(z - \zeta)(\partial A)(\zeta) d\xi d\eta$$

$$= \iint_{R^2} [f(z - \zeta) - f(z)](\partial A)(\zeta) d\xi d\eta \qquad (13)$$

The last equality depends on (9). Now (10) and (13) give (4). If we write (13) with Φ_x and Φ_y in place of $\partial \Phi$, we see that Φ has continuous partial derivatives, if we can show that $\partial \Phi = 0$ in G, where G is the set of all $z \in K$ whose distance from the complement of K exceeds δ . We shall do this by showing that

$$\Phi(z) = f(z) \quad (z \in G); \quad (14)$$

Note that $\partial f = 0$ in G, since f is holomorphic there. Now if $z \in G$, then $z - \zeta$ is in the interior of K for all ζ with $|\zeta| < \delta$. The mean value property for harmonic functions therefore gives, by the first equation in (11),

$$\Phi(z) = \int_0^\delta a(r)rdr \int_0^{2\pi} f(z - re^{i\theta})d\theta$$
$$= 2\pi f(z) \int_0^\delta a(r)rdr = f(z) \iint_{\mathbb{R}^2} A = f(z)$$
(15)

For all $z \in G$, we have now proved (3), (4), and (5) The definition of X shows that X is compact and that X can be covered by finitely many open discs $D_1,...,D_n$, of radius 2δ , whose centers are not in K. Since S^2-K is connected, the center of each D_j can be joined to ∞ by a polygonal path in S^2-K . It follows that each D_j contains a compact connected set E_j , of diameter at least 2δ , so that S^2-E_j is connected and so that $K\cap E_j=\phi$. with $r=2\delta$. There are functions $g_j\mathcal{E}H(S^2-E_j)$ and constants b_j so that the inequalities.

$$\left| Q_{j}(\zeta, z) \right| < \frac{50}{\delta}, \qquad (16)$$

$$\left| Q_{j}(\zeta, z) - \frac{1}{z - \zeta} \right| < \frac{4,000\delta^{2}}{\left| z - \zeta \right|^{2}} \qquad (17)$$

Hold for $z \notin E_i$ and $\zeta \in D_i$, if

$$Q_{i}(\zeta, z) = g_{i}(z) + (\zeta - b_{i})g_{i}^{2}(z)$$
 (18)

Let Ω be the complement of $E_1 \cup ... \cup E_n$. Then Ω is an open set which contains K. Put $X_1 = X \cap D_1$ and $X_2 = (X_1 \cap D_2) \cap (X_1 \cap A_2 \cap A_3)$

$$X_{j} = (X \cap D_{j}) - (X_{1} \cup ... \cup X_{j-1}),$$
 for

 $2 \le j \le n$,

Define

$$R(\zeta, z) = Q_i(\zeta, z)$$
 $(\zeta \varepsilon X_i, z \varepsilon \Omega)$ (19)

And

$$F(z) = \frac{1}{\pi} \iint_{X} (\partial \Phi)(\zeta) R(\zeta, z) d\zeta d\eta \qquad (20)$$
$$(z \in \Omega)$$

Since.

$$F(z) = \sum_{j=1}^{\infty} \frac{1}{\pi} \iint_{X_{j}} (\partial \Phi)(\zeta) Q_{j}(\zeta, z) d\xi d\eta, \qquad (21)$$

(18) shows that F is a finite linear combination of the functions g_j and g_j^2 . Hence $F \varepsilon H(\Omega)$. By (20), (4), and (5) we have

$$|F(z) - \Phi(z)| < \frac{2\omega(\delta)}{\pi \delta} \iint_{\mathbf{v}} |R(\zeta, z)|$$

$$-\frac{1}{z-\zeta}|d\xi d\eta \quad (z \in \Omega) \quad (22)$$

Observe that the inequalities (16) and (17) are valid with R in place of Q_j if $\zeta \in X$ and $z \in \Omega$. Now fix $z \in \Omega$., put $\zeta = z + \rho e^{i\theta}$, and estimate the integrand in (22) by (16) if $\rho < 4\delta$, by (17) if $4\delta \le \rho$. The integral in (22) is then seen to be less

$$2\pi \int_0^{4\delta} \left(\frac{50}{\delta} + \frac{1}{\rho} \right) \rho d\rho = 808\pi\delta \tag{23}$$

And

$$2\pi \int_{4\delta}^{\infty} \frac{4,000\delta^2}{\rho^2} \rho d\rho = 2,000\pi\delta.$$
 (24)

Hence (22) yields

than the sum of

$$|F(z) - \Phi(z)| < 6{,}000\omega(\delta)$$
 $(z \in \Omega)$ (25)

Since $F \in H(\Omega)$, $K \subset \Omega$, and $S^2 - K$ is connected, Runge's theorem shows that F can be uniformly approximated on K by polynomials. Hence (3) and (25) show that (2) can be satisfied. This completes the proof.



Lemma 1.0: Suppose $f \in C_c(\mathbb{R}^2)$, the space of all continuously differentiable functions in the plane, with compact support. Put

$$\partial = \frac{1}{2} \left(\frac{\partial}{\partial x} + i \frac{\partial}{\partial y} \right) \tag{1}$$

Then the following "Cauchy formula" holds:

$$f(z) = -\frac{1}{\pi} \iint_{R^2} \frac{(\partial f)(\zeta)}{\zeta - z} d\xi d\eta$$
$$(\zeta = \xi + i\eta) \tag{2}$$

Proof: This may be deduced from Green's theorem. However, here is a simple direct proof:

Put
$$\varphi(r,\theta) = f(z + re^{i\theta}), r > 0, \theta$$
 real

If $\zeta = z + re^{i\theta}$, the chain rule gives

$$(\partial f)(\zeta) = \frac{1}{2}e^{i\theta} \left[\frac{\partial}{\partial r} + \frac{i}{r} \frac{\partial}{\partial \theta} \right] \varphi(r,\theta) \tag{3}$$

The right side of (2) is therefore equal to the limit, as $\varepsilon \to 0$, of

$$-\frac{1}{2}\int_{\varepsilon}^{\infty}\int_{0}^{2\pi} \left(\frac{\partial\varphi}{\partial r} + \frac{i}{r}\frac{\partial\varphi}{\partial\theta}\right) d\theta dr \tag{4}$$

For each $r>0, \varphi$ is periodic in θ , with period 2π . The integral of $\partial\varphi/\partial\theta$ is therefore 0, and (4) becomes

$$-\frac{1}{2\pi} \int_0^{2\pi} d\theta \int_{\varepsilon}^{\infty} \frac{\partial \varphi}{\partial r} dr = \frac{1}{2\pi} \int_0^{2\pi} \varphi(\varepsilon, \theta) d\theta \tag{5}$$

As $\varepsilon \to 0$, $\varphi(\varepsilon, \theta) \to f(z)$ uniformly. This gives (2)

If $X^{\alpha}\in a$ and $X^{\beta}\in k\big[X_1,...X_n\big]$, then $X^{\alpha}X^{\beta}=X^{\alpha+\beta}\in a$, and so A satisfies the condition (*). Conversely,

$$(\sum_{\alpha \in A} c_{\alpha} X^{\alpha})(\sum_{\beta \in \mathbb{I}^{n}} d_{\beta} X^{\beta}) = \sum_{\alpha, \beta} c_{\alpha} d_{\beta} X^{\alpha + \beta} \qquad (finite sums),$$

and so if A satisfies (*), then the subspace generated by the monomials $X^{\alpha}, \alpha \in a$, is an ideal. The proposition gives a classification of the monomial ideals in $k \big[X_1, ... X_n \big]$: they are in one to one correspondence with the subsets A of \square^n satisfying (*). For example, the monomial ideals in $k \big[X \big]$ are exactly the ideals $(X^n), n \geq 1$, and the zero ideal (corresponding to the empty set A). We write $\left\langle X^{\alpha} \mid \alpha \in A \right\rangle$ for the ideal corresponding to

A (subspace generated by the X^{α} , $\alpha \in a$).

LEMMA 1.1. Let S be a subset of \square ⁿ. The the ideal α generated by $X^{\alpha}, \alpha \in S$ is the monomial ideal corresponding to

$$A = \{ \beta \in \square^n \mid \beta - \alpha \in \square^n, \quad some \ \alpha \in S \}$$

Thus, a monomial is in a if and only if it is divisible by one of the $X^{\alpha}, \alpha \in S$

PROOF. Clearly A satisfies (*), and $a \subset \langle X^{\beta} \mid \beta \in A \rangle$. Conversely, if $\beta \in A$, then $\beta - \alpha \in \Box^n$ for some $\alpha \in S$, and $X^{\beta} = X^{\alpha}X^{\beta-\alpha} \in a$. The last statement follows from the fact that $X^{\alpha} \mid X^{\beta} \Leftrightarrow \beta - \alpha \in \Box^n$. Let $A \subset \Box^n$ satisfy (*). From the geometry of A, it is clear that there is a finite set of elements $S = \{\alpha_1, ... \alpha_s\}$ of A such that $A = \{\beta \in \Box^n \mid \beta - \alpha_i \in \Box^2, some \alpha_i \in S\}$ (The α_i 's are the corners of A) Moreover, $a = \langle X^{\alpha} \mid \alpha \in A \rangle$ is generated by the monomials $X^{\alpha_i}, \alpha_i \in S$.

DEFINITION 1.0. For a nonzero ideal a in $k[X_1,...,X_n]$, we let (LT(a)) be the ideal generated by $\{LT(f) | f \in a\}$

LEMMA 1.2 Let a be a nonzero ideal in $k[X_1,...,X_n]$; then (LT(a)) is a monomial ideal, and it equals $(LT(g_1),...,LT(g_n))$ for some $g_1,...,g_n \in a$.

PROOF. Since (LT(a)) can also be described as the ideal generated by the leading monomials (rather than the leading terms) of elements of a.

THEOREM 1.2. Every *ideal* a in $k[X_1,...,X_n]$ is finitely generated; more precisely, $a=(g_1,...,g_s)$ where $g_1,...,g_s$ are any elements of a whose leading terms generate LT(a)

PROOF. Let $f \in a$. On applying the division algorithm, we find $f = a_1g_1 + ... + a_sg_s + r, \qquad a_i, r \in k\big[X_1,...,X_n\big]$



, where either r=0 or no monomial occurring in it is divisible by any $LT(g_i)$. But $r=f-\sum a_ig_i\in a$, and therefore $LT(r)\in LT(a)=(LT(g_1),...,LT(g_s))$, implies that every monomial occurring in r is divisible by one in $LT(g_i)$. Thus r=0, and $g\in (g_1,...,g_s)$.

DEFINITION 1.1. A finite subset $S = \{g_1, | ..., g_s\}$ of an ideal a is a standard ((Grobner) bases for a if $(LT(g_1), ..., LT(g_s)) = LT(a)$. In other words, S is a standard basis if the leading term of every element of a is divisible by at least one of the leading terms of the g_i .

THEOREM 1.3 The ring $k[X_1,...,X_n]$ is Noetherian i.e., every ideal is finitely generated.

PROOF. For n=1, k[X] is a principal ideal domain, which means that every ideal is generated by single element. We shall prove the theorem by induction on n. Note that the obvious map $k[X_1,...X_{n-1}][X_n] \rightarrow k[X_1,...X_n]$ is an isomorphism – this simply says that every polynomial f in n variables $X_1,...X_n$ can be expressed uniquely as a polynomial in X_n with coefficients in $k[X_1,...,X_n]$:

$$f(X_1,...X_n) = a_0(X_1,...X_{n-1})X_n^r + ... + a_r(X_1,...X_{n-1})$$

Thus the next lemma will complete the proof

LEMMA 1.3. If A is Noetherian, then so also is A[X]

PROOF. For a polynomial

$$f(X) = a_0 X^r + a_1 X^{r-1} + \ldots + a_r, \quad a_i \in A, \quad a_0 \neq 0,$$
 r is called the degree of f , and a_0 is its leading coefficient. We call 0 the leading coefficient of the polynomial 0. Let a be an ideal in $A[X]$. The leading coefficients of the polynomials in a form an ideal a in A , and since A is Noetherian, a will be finitely generated. Let g_1, \ldots, g_m be elements of a whose leading coefficients generate a , and let a be the maximum degree of a . Now

let $f \in a$, and suppose f has degree s > r, say, $f = aX^s + ...$ Then $a \in a$, and so we can write $a = \sum b_i a_i$, $b_i \in A$, $a_i = leading \ coefficient \ of \ g_i$

 $f - \sum b_i g_i X^{s-r_i}$, $r_i = \deg(g_i)$, has degree $< \deg(f)$. By continuing in this way, we find that $mod(g_1,...g_m)$ With $f \equiv f_{t}$ polynomial of degree t < r For each d < r, let a_d be the subset of A consisting of 0 and the leading coefficients of all polynomials in a of degree d; it is again an ideal in A. Let $g_{d,1},...,g_{d,m_d}$ be polynomials of degree d whose leading coefficients generate $\boldsymbol{a}_{\boldsymbol{d}}$. Then the same argument as above shows that any polynomial f_d in degree can of d $mod(g_{d,1},...g_{d,m_d})$ With f_{d-1} $f_d \equiv f_{d-1}$ of degree $\leq d-1$. On applying this remark we repeatedly $f_t \in (g_{r-1,1}, ..., g_{r-1,m_{s-1}}, ..., g_{0,1}, ..., g_{0,m_0})$ Hence

$$f_t \in (g_1, ... g_m g_{r-1,1}, ... g_{r-1,m_{r-1}}, ..., g_{0,1}, ..., g_{0,m_0})$$

and so the polynomials $g_1, ..., g_{0,m_0}$ generate a

Definition 2 (**RCPI**). An instance of a role containment problem (RCPI) is given by a triple hP,R,X.u w A.ri. An RCPI is said to be satisfied if and only if [[X.u]]P0 _ [[A.r]]P0 for each P0 such that P _ 7!R P0. In this case we also say that P satisfies X.u w A.r under R.

III. COMPLEXITY REDUCTION

This section describes several reductions that transform one RCPI into another that is typically less expensive to evaluate. Our findings in Section 5 indicate that, when using our model checking technique and our platform configuation, these reductions often make the difference between being unable to evaluate an RCPI and being able to do so. We conjecture that they may also reduce the cost of applying other approaches to solve RCPI problems, such as one based one on the proof method of Sistla [20, 21].

One of the great successes of category theory in computer science has been the development of a "unified theory" of the constructions underlying denotational semantics. In the untyped λ -calculus, any term may appear in the function position of an application. This means



that a model D of the λ -calculus must have the property that given a term t whose interpretation is $d \in D$, Also, the interpretation of a functional abstraction like $\lambda x \cdot x$ is most conveniently defined as a function from $D \, to \, D$, which must then be regarded an element of $\psi: [D \to D] \to D$ be the function that picks out elements of D to represent elements of $[D \rightarrow D]$ and $\phi: D \to [D \to D]$ be the function that maps elements of D to functions of D. Since $\psi(f)$ is intended to represent the function f as an element of D, it makes sense to require that $\phi(\psi(f)) = f$, that is, $\psi \, o \psi \! = \! i d_{[D \to D]}$ Furthermore, we often want to view every element of D as representing some function from D to D and require that elements representing the same function be equal -

$$\psi(\varphi(d)) = d$$

or

$$\psi \circ \phi = id_D$$

The latter condition is called extensionality. These conditions together imply that ϕ and ψ are inverses--- that is, D is isomorphic to the space of functions from D to D that can be the functional abstractions: interpretations of $D \cong [D \to D]$.Let us suppose we are working with the untyped λ -calculus, we need a solution of the equation $D \cong A + [D \rightarrow D]$, where A is some predetermined domain containing interpretations for elements of C. Each element of D corresponds to either an element of A or an element of $[D \rightarrow D]$, with a tag. This equation can be solved by finding least fixed points of the function $F(X) = A + |X \rightarrow X|$ from domains to domains --- that is, finding domains X such that $X \cong A + [X \longrightarrow X]$, and such that for any domain Y also satisfying this equation, there is an embedding of X to Y --- a pair of maps

$$X igcup_{f^R}^f Y$$
Such that
$$f^R \circ f = id_X$$

$$f \circ f^R \subseteq id_Y$$
Where $f \subseteq g$ means that
$$f \ approximates \ g \ \text{in some ordering representing}$$

their information content. The key shift of perspective from the domain-theoretic to the more general category-theoretic approach lies in considering F not as a function on domains, but as a functor on a category of domains. Instead of a least fixed point of the function, F.

Definition 1.3: Let K be a category and $F: K \to K$ as a functor. A fixed point of F is a pair (A,a), where A is a K-object and $a: F(A) \to A$ is an isomorphism. A prefixed point of F is a pair (A,a), where A is a K-object and a is any arrow from F(A) to A

Definition 1.4: An ω -chain in a category K is a diagram of the following form:

$$\Delta = D_o \xrightarrow{f_o} D_1 \xrightarrow{f_1} D_2 \xrightarrow{f_2} \dots$$

Recall that a cocone μ of an ω -chain Δ is a K-object X and a collection of K -arrows $\{\mu_i:D_i\to X\,|\,i\ge 0\}$ such that $\mu_i=\mu_{i+1}o\ f_i$ for all $i\ge 0$. We sometimes write $\mu:\Delta\to X$ as a reminder of the arrangement of μ 's components Similarly, a colimit $\mu:\Delta\to X$ is a cocone with the property that if $\nu:\Delta\to X'$ is also a cocone then there exists a unique mediating arrow $k:X\to X'$ such that for all $i\ge 0$,, $v_i=k\ o\ \mu_i$. Colimits of ω -chains are sometimes referred to as ω -co limits. Dually, an ω^{op} -chain in K is a diagram of the following form:

 f_o f_1 $\Delta = D_o \longleftarrow D_1 \longleftarrow D_2 \longleftarrow \dots$ $\mu: X \to \Delta$ of an ω^{op} - chain Δ is a **K**-object X and a collection of **K**-arrows $\{\mu_i : D_i \mid i \geq 0\}$ such that for all $i \ge 0$, $\mu_i = f_i \circ \mu_{i+1}$. An ω^{op} limit of an ω^{op} – chain Δ is a cone $\mu: X \to \Delta$ with the property that if $\nu: X^{'} \to \Delta$ is also a cone, then there exists a unique mediating arrow $k: X' \to X$ such that for all $i \ge 0$, $\mu_i \circ k = \nu_i$. We write \perp_k (or just \perp) for the distinguish initial object of K, when it has one, and $\bot \to A$ for the unique arrow from \perp to each K-object A. It is also convenient to write $\Delta^- = D_1 \longrightarrow D_2 \longrightarrow \dots$ to denote all of Δ except D_o and f_0 . By analogy, μ^- is $\{\mu_i \mid i \geq 1\}$. For the images of Δ and μ



under
$$F$$
 we write $F(\Delta) = F(D_o) \xrightarrow{F(f_o)} F(D_1) \xrightarrow{F(f_1)} F(D_2) \xrightarrow{F(f_2)} \dots$ and $F(\mu) = \{F(\mu_i) | i \ge 0\}$

We write F^i for the **i**-fold iterated composition of F — that is, $F^o(f) = f$, $F^1(f) = F(f)$, $F^2(f) = F(F(f))$, etc. With these definitions we can state that every monitonic function on a complete lattice has a least fixed point:

Lemma 1.4. Let K be a category with initial object \bot and let $F: K \to K$ be a functor. Define the $\omega - chain \Delta$ by

$$\Delta = \perp \xrightarrow{f(\bot)} F(\bot) \xrightarrow{F(\bot \to F(\bot))} F^2(\bot) \xrightarrow{F^2(\bot \to F(\bot))} \cdots \cdots$$
If both $\mu : \Delta \to D$ and $F(\mu) : F(\Delta) \to F(D)$ are colimits, then (D,d) is an intial F-algebra, where $d : F(D) \to D$ is the mediating arrow from $F(\mu)$ to the cocone μ^-

Theorem 1.4 Let a DAG G given in which each node is a random variable, and let a discrete conditional probability distribution of each node given values of its parents in G be specified. Then the product of these conditional distributions yields a joint probability distribution P of the variables, and (G,P) satisfies the Markov condition.

Proof. Order the nodes according to an ancestral ordering. Let X_1, X_2, \ldots, X_n be the resultant ordering. Next define.

$$P(x_1, x_2,...x_n) = P(x_n | pa_n) P(x_{n-1} | Pa_{n-1})...$$

.. $P(x_2 | pa_2) P(x_1 | pa_1),$

Where PA_i is the set of parents of X_i of in G and $P(x_i \mid pa_i)$ is the specified conditional probability distribution. First we show this does indeed yield a joint probability distribution. Clearly, $0 \le P(x_1, x_2, ...x_n) \le 1$ for all values of the variables. Therefore, to show we have a joint distribution, as the variables range through all their possible values, is equal to one. To that end, Specified conditional distributions are the conditional distributions they notationally represent in the joint distribution. Finally, we show the Markov condition is satisfied. To do this, we need show for $1 \le k \le n$ that

whenever

$$P(pa_k) \neq 0, if \ P(nd_k \mid pa_k) \neq 0$$

 $and \ P(x_k \mid pa_k) \neq 0$
 $then \ P(x_k \mid nd_k, pa_k) = P(x_k \mid pa_k),$

Where ND_k is the set of nondescendents of X_k of in G. Since $PA_k \subseteq ND_k$, we need only show $P(x_k \mid nd_k) = P(x_k \mid pa_k)$. First for a given k, order the nodes so that all and only nondescendents of X_k precede X_k in the ordering. Note that this ordering depends on k, whereas the ordering in the first part of the proof does not. Clearly then

$$\begin{aligned} ND_k &= \left\{X_1, X_2, X_{k-1}\right\} \\ Let \\ D_k &= \left\{X_{k+1}, X_{k+2}, X_n\right\} \\ \text{follows } \sum_{k,l} d_{k+1} \end{aligned}$$

We define the m^{th} cyclotomic field to be the field $Q[x]/(\Phi_m(x))$ Where $\Phi_m(x)$ is the m^{th} cyclotomic polynomial. $Q[x]/(\Phi_m(x))$ $\Phi_m(x)$ has degree $\varphi(m)$ over Q since $\Phi_m(x)$ has degree $\varphi(m)$. The roots of $\Phi_m(x)$ are just the primitive m^{th} roots of unity, so the complex embeddings of $Q[x]/(\Phi_m(x))$ are simply the $\varphi(m)$ maps

$$\sigma_{k}: Q[x]/(\Phi_{m}(x)) \mapsto C,$$

$$1 \le k < m, (k, m) = 1, \quad where$$

$$\sigma_{k}(x) = \xi_{m}^{k},$$

 ξ_m being our fixed choice of primitive m^{th} root of unity. Note that $\xi_m^k \in Q(\xi_m)$ for every k; it follows that $Q(\xi_m) = Q(\xi_m^k)$ for all k relatively prime to m. In particular, the images of the σ_i coincide, so $Q[x]/(\Phi_m(x))$ is Galois over Q. This means that we can write $Q(\xi_m)$ for $Q[x]/(\Phi_m(x))$ without much fear of ambiguity; we will do so from now on, the identification being $\xi_m \mapsto x$. One advantage of this is that one can easily talk about cyclotomic fields being extensions of one another, or intersections or compositums; all of these things take place considering them as



subfield of C. We now investigate some basic properties of cyclotomic fields. The first issue is whether or not they are all distinct; to determine this, we need to know which roots of unity lie in $Q(\xi_m)$. Note, for example, that if m is odd, then $-\xi_m$ is a $2m^{th}$ root of unity. We will show that this is the only way in which one can obtain any non- m^{th} roots of unity.

LEMMA 1.5 If m divides n, then $Q(\xi_m)$ is contained in $Q(\xi_m)$

PROOF. Since $\xi^{n/m} = \xi_m$, we have $\xi_m \in Q(\xi_n)$, so the result is clear

LEMMA 1.6 If m and n are relatively prime, then

$$Q(\xi_m, \xi_n) = Q(\xi_{nm})$$

and

$$Q(\xi_m) \cap Q(\xi_n) = Q$$

(Recall the $Q(\xi_m,\xi_n)$ is the compositum of $Q(\xi_m)$ and $Q(\xi_n)$)

PROOF. One checks easily that $\xi_m \xi_n$ is a primitive mn^{th} root of unity, so that

$$Q(\xi_{mn}) \subseteq Q(\xi_{m}, \xi_{n})$$

$$[Q(\xi_m,\xi_n):Q] \leq [Q(\xi_m):Q][Q(\xi_n:Q]$$

 $= \varphi(m)\varphi(n) = \varphi(mn);$

Since $Q(\xi_{mn}):Q=\varphi(mn)$; this implies that $Q(\xi_m,\xi_n)=Q(\xi_{nm})$ We know that $Q(\xi_m,\xi_n)$ has degree $\varphi(mn)$ over Q, so we must have

$$[Q(\xi_m,\xi_n):Q(\xi_m)]=\varphi(n)$$

and

$$[Q(\xi_m, \xi_n): Q(\xi_m)] = \varphi(m)$$

$$[Q(\xi_m):Q(\xi_m)\cap Q(\xi_n)]\geq \varphi(m)$$

And thus that $Q(\xi_m) \cap Q(\xi_n) = Q$

PROPOSITION 1.2 For any m and n

$$Q(\xi_m, \xi_n) = Q(\xi_{[m,n]})$$
And
$$Q(\xi_m) \cap Q(\xi_n) = Q(\xi_{(m,n)});$$

here [m, n] and (m, n) denote the least common multiple and the greatest common divisor of m and n, respectively.

PROOF. Write $m = p_1^{e_1} \dots p_k^{e_k}$ and $p_1^{f_1} \dots p_k^{f_k}$ where the p_i are

distinct primes. (We allow e_i or f_i to be zero)

$$Q(\xi_m) = Q(\xi_{n^{e_1}})Q(\xi_{n^{e_2}})...Q(\xi_{n^{e_k}})$$

and

$$Q(\xi_n) = Q(\xi_{p_1,f_1})Q(\xi_{p_2,f_2})...Q(\xi_{p_k,f_k})$$

Thus

$$\begin{split} Q(\xi_{m},\xi_{n}) &= Q(\xi_{p_{1}^{e_{1}}})......Q(\xi_{p_{2}^{e_{k}}})Q(\xi_{p_{1}^{f_{1}}})...Q(\xi_{p_{k}^{f_{k}}}) \\ &= Q(\xi_{p_{1}^{e_{1}}})Q(\xi_{p_{1}^{f_{1}}})...Q(\xi_{p_{k}^{e_{k}}})Q(\xi_{p_{k}^{f_{k}}}) \\ &= Q(\xi_{p_{1}^{\max(e_{1},f_{1})}})......Q(\xi_{p_{n}^{\max(e_{k},f_{k})}}) \\ &= Q(\xi_{p_{1}^{\max(e_{1},f_{1})}}......p_{1}^{\max(e_{k},f_{k})}) \\ &= Q(\xi_{[m,n]}); \end{split}$$

An entirely similar computation shows that $Q(\xi_m) \cap Q(\xi_n) = Q(\xi_{(m,n)})$

Mutual information measures the information transferred when X_i is sent and Y_i is received, and is defined as

$$I(x_i, y_i) = \log_2 \frac{P(x_i/y_i)}{P(x_i)} bits$$
 (1)

In a noise-free channel, **each** y_i is uniquely connected to the corresponding x_i , and so they constitute an input –output pair (x_i, y_i) for which

$$P(\frac{x_i}{y_j}) = 1 \text{ and } I(x_i, y_j) = \log_2 \frac{1}{P(x_i)}$$
 bits;

that is, the transferred information is equal to the self-information that corresponds to the input x_i In a very noisy channel, the output y_i and input x_i would be completely uncorrelated, and so

$$P(\frac{x_i}{y_j}) = P(x_i)$$
 and also $I(x_i, y_j) = 0$; that is,

there is no transference of information. In general, a given channel will operate between these two extremes. The mutual information is defined between the input and the output of a given channel. An average of the calculation of the mutual information for all input-output pairs of a given channel is the average mutual information:



$$I(X,Y) = \sum_{i,j} P(x_i, y_j) I(x_i, y_j) = \sum_{i,j} P(x_i, y_j) \log_2 \left[\frac{P(x_i/y_j)}{P(x_i)} \right]$$

bits per symbol. This calculation is done over the input and output alphabets. The average mutual information. The following expressions are useful for modifying the mutual information expression:

$$P(x_i, y_j) = P(\frac{x_i}{y_j})P(y_j) = P(\frac{y_j}{x_i})P(x_i)$$

$$P(y_j) = \sum_i P(\frac{y_j}{x_i})P(x_i)$$

$$P(x_i) = \sum_i P(\frac{x_i}{y_j})P(y_j)$$

Ther

$$I(X,Y) = \sum_{i,j} P(x_i, y_j)$$

$$= \sum_{i,j} P(x_i, y_j) \log_2 \left[\frac{1}{P(x_i)} \right]$$

$$-\sum_{i,j} P(x_i, y_j) \log_2 \left[\frac{1}{P(x_i/y_j)} \right]$$

$$\sum_{i,j} P(x_i, y_j) \log_2 \left[\frac{1}{P(x_i)} \right]$$

$$= \sum_{i} \left[P(x_i / y_j) P(y_j) \right] \log_2 \frac{1}{P(x_i)}$$

$$\sum_{i} P(x_i) \log_2 \frac{1}{P(x_i)} = H(X)$$

$$I(X,Y) = H(X) - H(X/Y)$$

Where

$$H(X/Y) = \sum_{i,j} P(x_i, y_j) \log_2 \frac{1}{P(X_i/Y_i)}$$
 is

usually called the equivocation. In a sense, the equivocation can be seen as the information lost in the noisy channel, and is a function of the backward conditional probability. The observation of an output symbol y_j provides H(X) - H(X/Y) bits of information. This difference is the mutual information of the channel. *Mutual Information: Properties* Since

$$P(\frac{x_i}{y_j})P(y_j) = P(\frac{y_j}{x_i})P(x_i)$$

The mutual information fits the condition

$$I(X,Y) = I(Y,X)$$

And by interchanging input and output it is also true that

$$I(X,Y) = H(Y) - H(\frac{Y}{X})$$

Where

$$H(Y) = \sum_{j} P(y_j) \log_2 \frac{1}{P(y_j)}$$

This last entropy is usually called the noise entropy. Thus, the information transferred through the channel is the difference between the output entropy and the noise entropy. Alternatively, it can be said that the channel mutual information is the difference between the number of bits needed for determining a given input symbol before knowing the corresponding output symbol, and the number of bits needed for determining a given input symbol after knowing the corresponding output symbol

$$I(X,Y) = H(X) - H(X/Y)$$

As the channel mutual information expression is a difference between two quantities, it seems that this parameter can adopt negative values. However, and is spite of the fact that for some y_j , $H(X \mid y_j)$ can be larger than H(X), this is not possible for the average value calculated over all the outputs:

$$\sum_{i,j} P(x_i, y_j) \log_2 \frac{P(x_i/y_j)}{P(x_i)} = \sum_{i,j} P(x_i, y_j) \log_2 \frac{P(x_i, y_j)}{P(x_i)P(y_j)}$$

Ther

$$-I(X,Y) = \sum_{i,j} P(x_i, y_j) \frac{P(x_i)P(y_j)}{P(x_i, y_i)} \le 0$$

Because this expression is of the form

$$\sum_{i=1}^{M} P_i \log_2(\frac{Q_i}{P_i}) \le 0$$

The above expression can be applied due to the factor $P(x_i)P(y_j)$, which is the product of two probabilities, so that it behaves as the quantity Q_i , which in this expression is a dummy variable that fits the condition $\sum_i Q_i \leq 1$. It can be concluded that the average mutual information is a nonnegative number. It can also be equal to zero, when the input and the output are independent of each other. A related entropy called the joint entropy is defined as

$$H(X,Y) = \sum_{i,j} P(x_i, y_j) \log_2 \frac{1}{P(x_i, y_j)}$$
$$= \sum_{i,j} P(x_i, y_j) \log_2 \frac{P(x_i)P(y_j)}{P(x_i, y_j)}$$
$$+ \sum_{i,j} P(x_i, y_j) \log_2 \frac{1}{P(x_i)P(y_j)}$$



Theorem 1.5: Entropies of the binary erasure channel (BEC) The BEC is defined with an alphabet of two inputs and three outputs, with symbol probabilities.

 $P(x_1) = \alpha$ and $P(x_2) = 1 - \alpha$, and transition probabilities

$$P(\frac{y_3}{x_2}) = 1 - p \text{ and } P(\frac{y_2}{x_1}) = 0,$$
and $P(\frac{y_3}{x_1}) = 0$
and $P(\frac{y_1}{x_2}) = p$
and $P(\frac{y_3}{x_2}) = 1 - p$

Lemma 1.7. Given an arbitrary restricted time-discrete, amplitude-continuous channel whose restrictions are determined by sets F_n and whose density functions exhibit no dependence on the state s, let n be a fixed positive integer, and p(x) an arbitrary probability density function on Euclidean n-space. $p(y \mid x)$ for the density $p_n(y_1,...,y_n \mid x_1,...x_n)$ and F for F_n . For any real number a, let

$$A = \left\{ (x, y) : \log \frac{p(y \mid x)}{p(y)} > a \right\} \tag{1}$$

Then for each positive integer u , there is a code (u,n,λ) such that

$$\lambda \le ue^{-a} + P\{(X,Y) \notin A\} + P\{X \notin F\}$$

Where

 $P\{(X,Y) \in A\} = \int_{A} \dots \int p(x,y) dx dy, \qquad p(x,y) = p(x)p(y \mid x)$

$$P\{X \in F\} = \int_{F} \dots \int p(x) dx$$

Proof: A sequence $x^{(1)} \in F$ such that

$$P\left\{Y \in A_{x^{1}} \mid X = x^{(1)}\right\} \ge 1 - \varepsilon$$

where
$$A_x = \{y: (x, y) \in A\};$$

Choose the decoding set B_1 to be $A_{x^{(1)}}$. Having chosen $x^{(1)},\ldots,x^{(k-1)}$ and B_1,\ldots,B_{k-1} , select $x^k\in F$ such that

$$P\left\{Y \in A_{x^{(k)}} - \bigcup_{i=1}^{k-1} B_i \mid X = x^{(k)}\right\} \ge 1 - \varepsilon;$$

Set $B_k = A_{x^{(k)}} - \bigcup_{i=1}^{k-1} B_i$, If the process does not terminate in a finite number of steps, then the sequences $\mathcal{X}^{(i)}$ and decoding sets

 $B_i, i=1,2,...,u$, form the desired code. Thus assume that the process terminates after t steps. (Conceivably t=0). We will show $t \ge u$ by showing

 $\varepsilon \le te^{-a} + P\{(X,Y) \notin A\} + P\{X \notin F\}$. We proceed as follows.

Let

$$B = \bigcup_{j=1}^{t} B_{j}. \quad (If \quad t = 0, \text{ take } B = \emptyset). \text{ Then}$$

$$P\{(X,Y) \in A\} = \int_{(x,y)\in A} p(x,y)dxdy$$

$$= \int_{x} p(x) \int_{y \in A_{x}} p(y \mid x) dy dx$$

$$= \int_{x} p(x) \int_{y \in B \cap A_{x}} p(y \mid x) dy dx + \int_{x} p(x)$$

IV. EXPERIMENTAL DESIGN

We evaluate the performance of our scheme and study various "what-if" scenarios through detailed simulation experiments. We compare our scheme (2) against existing alternatives of using a least recently used (LRU) or a least frequently used (LFU) cache replacement strategy.

A. Algorithms

Ideals. Let A be a ring. Recall that an *ideal a* in A is a subset such that a is subgroup of A regarded as a group under addition;

$$a \in a, r \in A \Rightarrow ra \in A$$

The ideal generated by a subset S of A is the intersection of all ideals A containing a ---- it is easy to verify that this is in fact an ideal, and that it consist of all finite sums of the form $\sum r_i s_i$ with

$$r_i \in A, s_i \in S$$
. When $S = \{s_1,, s_m\}$, we shall write $(s_1,, s_m)$ for the ideal it generates.

Let a and b be ideals in A. The set $\left\{a+b \,|\, a\in a, b\in b\right\} \text{ is an ideal, denoted by } a+b \quad . \text{ The ideal generated by } \left\{ab \,|\, a\in a, b\in b\right\} \text{ is denoted by } ab \text{ . Note that } ab \subset a \cap b \text{ . Clearly } ab \text{ consists of all finite sums } \sum a_i b_i \quad \text{with } a_i \in a \quad \text{and} \quad b_i \in b \quad , \quad \text{and if } a \in a \quad \text{and} \quad b_i \in b \quad , \quad \text{and if } a \in a \quad \text{and} \quad b \in b \quad .$



 $a = (a_1, ..., a_m)$ and $b = (b_1, ..., b_n)$, then $ab = (a_1b_1, ..., a_ib_i, ..., a_mb_n)$. Let a be an ideal of A. The set of cosets of a in A forms a ring A/aand $a \mapsto a + a$ is a homomorphism $\phi: A \mapsto A/a$. The map $b \mapsto \phi^{-1}(b)$ is a one to one correspondence between the ideals of A/aand the ideals of A containing a An ideal p if prime if $p \neq A$ and $ab \in p \Rightarrow a \in p$ or $b \in p$. Thus p is prime if and only if A/p is nonzero has the property $b \neq 0 \Rightarrow a = 0$, i.e., A/p is an ab = 0, integral domain. An ideal m is maximal if $m \neq A$ and there does not exist an ideal n contained strictly between m and A. Thus m is maximal if and only if A/m has no proper nonzero ideals, and so is a field. Note that m maximal $\Rightarrow m$ prime. The ideals of $A \times B$ are all of the form $a \times b$, with a and b ideals in A and B. To see this, note that if c is an ideal in $A \times B$ and $(a,b) \in c$, then $(a,0) = (a,b)(1,0) \in c$ $(0,b) = (a,b)(0,1) \in c$ This shows that $c = a \times b$ with $a = \{a \mid (a,b) \in c \text{ some } b \in b\}$

B together with a homomorphism $i_B:A\to B$. A homomorphism of A -algebra $B\to C$ is a homomorphism of rings $\varphi:B\to C$ such that $\varphi(i_B(a))=i_C(a)$ for all $a\in A$. An A-algebra B is said to be finitely generated (or of finite-type over A) if there exist elements $x_1,...,x_n\in B$ such that every element of B can be expressed as a polynomial in the x_i with coefficients in i(A), i.e., such that the homomorphism $A\big[X_1,...,X_n\big]\to B$ sending X_i to x_i is surjective. A ring homomorphism $A\to B$ is finite, and B is finitely generated as an A-module. Let B be a field, and let A be a B-algebra. If B in B is injective, we can identify B with its

Let A be a ring. An A-algebra is a ring

 $b = \{b \mid (a,b) \in c \text{ some } a \in a\}$

. **Polynomial rings.** Let k be a field. A *monomial* in $X_1, ..., X_n$ is an expression of the form

image, i.e., we can regard k as a subring of A. If

1=0 in a ring R, the R is the zero ring, i.e., $R = \{0\}$

 $X_1^{a_1}...X_n^{a_n}, \qquad a_i \in N$. The *total degree* of the monomial is $\sum a_i$. We sometimes abbreviate it by X^{α} , $\alpha = (a_1, ..., a_n) \in \square^n$ The elements of the polynomial ring $k[X_1,...,X_n]$ are finite sums $\sum c_{a_1...a_n} X_1^{a_1} ... X_n^{a_n}, \qquad c_{a_1...a_n} \in k, \quad a_j \in \square$ With the obvious notions of equality, addition and multiplication. Thus the monomials from basis for $k[X_1,...,X_n]$ as a k-vector space. The ring $k[X_1,...,X_n]$ is an integral domain, and the only units in it are the nonzero constant polynomials. A polynomial $f(X_1,...,X_n)$ is irreducible if it is nonconstant and has only the obvious factorizations, i.e., $f = gh \Rightarrow g$ or h is constant. **Division in** k[X]. The division algorithm allows us to divide a nonzero polynomial into another: let f and g be polynomials in k[X] with $g \neq 0$; then there exist unique polynomials $q, r \in k[X]$ such that f = qg + r with either r = 0 or $\deg r < \deg g$. Moreover, there is an algorithm for deciding whether $f \in (g)$, namely, find r and check whether it is zero. Moreover, the Euclidean algorithm allows to pass from finite set of generators for an ideal in $k \mid X \mid$ to a single generator by successively replacing each pair of generators with their greatest common divisor. (Pure) lexicographic ordering (lex). monomials are ordered by lexicographic(dictionary) order. More precisely, let $\alpha = (a_1, ... a_n)$ and $\beta = (b_1, ..., b_n)$ be two elements of \square^n ; then $\alpha > \beta$ and $X^{\alpha} > X^{\beta}$ (lexicographic ordering) if, in the vector difference $\alpha - \beta \in \square$, the left most nonzero entry is positive. For example,

 $XY^2 > Y^3Z^4$; $X^3Y^2Z^4 > X^3Y^2Z$. Note that this isn't quite how the dictionary would order them: it would put XXXYYZZZZ after XXXYYZ. Graded reverse lexicographic order (grevlex). Here monomials are ordered by total degree, with ties broken by reverse lexicographic ordering. Thus, $\alpha > \beta$ if $\sum a_i > \sum b_i$, or $\sum a_i = \sum b_i$ and in $\alpha - \beta$ the right most nonzero entry is negative. For example:

$$X^4Y^4Z^7 > X^5Y^5Z^4$$
 (total degree greater)
 $XY^5Z^2 > X^4YZ^3$, $X^5YZ > X^4YZ^2$



Orderings on $k[X_1,...X_n]$. Fix an ordering on the monomials in $k[X_1,...X_n]$. Then we can write an element f of $k[X_1,...X_n]$ in a canonical fashion, by re-ordering its elements in decreasing order. For example, we would write

$$f = 4XY^{2}Z + 4Z^{2} - 5X^{3} + 7X^{2}Z^{2}$$
as
$$f = -5X^{3} + 7X^{2}Z^{2} + 4XY^{2}Z + 4Z^{2} \quad (lex)$$
or
$$f = 4XY^{2}Z + 7X^{2}Z^{2} - 5X^{3} + 4Z^{2} \quad (grevlex)$$

Let $\sum a_{\alpha}X^{\alpha}\in k\big[X_1,...,X_n\big]$, in decreasing order:

$$f = a_{\alpha_0} X^{\alpha_0} +_{\alpha_1} X^{\alpha_1} + ..., \qquad \alpha_0 > \alpha_1 > ..., \quad \alpha_0$$

- The *multidegree* of f to be multdeg(f)=
- The leading coefficient of f to be LC(f)
- The leading monomial of f to be LM(f)
- The leading term of f to be LT(f) =

For the polynomial $f = 4XY^2Z + ...$, the multidegree is (1,2,1), the leading coefficient is 4, the leading monomial is XY^2Z , and the leading term is $4XY^2Z$. The division algorithm in $k[X_1,...X_n]$. Fix a monomial ordering in \square^2 . Suppose given a polynomial f and an ordered set $(g_1,...g_s)$ of polynomials; the division algorithm then constructs polynomials $a_1,...a_s$ and r such that $f = a_1 g_1 + ... + a_s g_s + r$ Where either r = 0 or no monomial in r is divisible by any of $LT(g_1),...,LT(g_s)$ $LT(g_1) | LT(f)$, divide g_1 into f to get $f = a_1 g_1 + h, \qquad a_1 = \frac{LT(f)}{LT(g_1)} \in k[X_1, ..., X_n]$ If $LT(g_1)|LT(h)$, repeat the process until $f = a_1 g_1 + f_1$ (different a_1) with $LT(f_1)$ not

divisible by $LT(g_1)$. Now divide g_2 into f_1 , and

so on, until $f = a_1 g_1 + ... + a_s g_s + r_1$

 $LT(r_1)$ not divisible by any $LT(g_1),...LT(g_s)$ **Step 2:** Rewrite $r_1 = LT(r_1) + r_2$, and repeat Step r_2 for $f = a_1 g_1 + ... + a_s g_s + LT(r_1) + r_3$ (different a_i 's) Monomial ideals. In general, an ideal awill contain a polynomial without containing the individual terms of the polynomial; for example, the ideal $a = (Y^2 - X^3)$ contains $Y^2 - X^3$ but not Y^2 or X^3

DEFINITION 1.5. An ideal a is monomial if $\sum c_{\alpha}X^{\alpha} \in a \Rightarrow X^{\alpha} \in a$ all α with $c_{\alpha} \neq 0$.

 $f = a_{\alpha_0} X^{\alpha_0} +_{\alpha_1} X^{\alpha_1} + ...,$ $\alpha_0 > \alpha_1 > ...,$ $\alpha_0 \neq 0$ **PROPOSITION 1.3.** Let a be a monomial ideal, and let $A = \{ \alpha \mid X^{\alpha} \in a \}$. Then A satisfies the condition $\alpha \in A$, $\beta \in \square$ $^n \Rightarrow \alpha + \beta \in$ And a is the k -subspace of $k[X_1,...,X_n]$ generated by the X^{α} , $\alpha \in A$. Conversely, of A is a subset of \square ⁿ satisfying (*), then the k-subspace a of $k[X_1,...,X_n]$ generated by $\{X^{\alpha} \mid \alpha \in A\}$ is a monomial ideal.

> PROOF. It is clear from its definition that a monomial ideal a is the k -subspace of $k[X_1,...,X_n]$ generated by the set of monomials it contains. If

> $X^{\alpha} \in a$ and $X^{\beta} \in k[X_1, ..., X_n]$

If a permutation is chosen uniformly and at random from the n! possible permutations in S_n , then the counts $C_i^{(n)}$ of cycles of length j are dependent random variables. The joint distribution of $C^{(n)} = (C_1^{(n)}, ..., C_n^{(n)})$ follows from Cauchy's formula, and is given by

$$P[C^{(n)} = c] = \frac{1}{n!}N(n,c) = 1\left\{\sum_{j=1}^{n} jc_{j} = n\right\} \prod_{j=1}^{n} \left(\frac{1}{j}\right)^{c_{j}} \frac{1}{c_{j}!}, \quad (1.1)$$

for $c \in \square_+^n$.

Lemma_{1.7} For nonnegative integers m_1 m_n ,

$$E\left(\prod_{j=1}^{n} (C_{j}^{(n)})^{[m_{j}]}\right) = \left(\prod_{j=1}^{n} \left(\frac{1}{j}\right)^{m_{j}}\right) 1 \left\{\sum_{j=1}^{n} j m_{j} \le n\right\}$$
(1.4)

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Proof. This can be established directly by exploiting cancellation of the form $c_j^{[m_j]}/c_j^!=1/(c_j-m_j)!$ when $c_j\geq m_j$, which occurs between the ingredients in Cauchy's formula and the falling factorials in the moments. Write $m=\sum jm_j$. Then, with the first sum indexed by $c=(c_1,...c_n)\in \square_+^n$ and the last sum indexed by $d=(d_1,...,d_n)\in \square_+^n$ via the correspondence $d_j=c_j-m_j$, we have

$$\begin{split} E\Bigg(\prod_{j=1}^{n}(C_{j}^{(n)})^{[m_{j}]}\Bigg) &= \sum_{c}P[C^{(n)} = c] \prod_{j=1}^{n}(c_{j})^{[m_{j}]} \\ &= \sum_{c:c_{j} \geq m_{j} \ for \ all \ j} \mathbf{1} \bigg\{\sum_{j=1}^{n} jc_{j} = n\bigg\} \prod_{j=1}^{n} \frac{(c_{j})^{[m_{j}]}}{j^{c_{j}}c_{j}!} \\ &= \prod_{j=1}^{n} \frac{1}{j^{m_{j}}} \sum_{d} \mathbf{1} \bigg\{\sum_{j=1}^{n} jd_{j} = n - m\bigg\} \prod_{j=1}^{n} \frac{1}{j^{d_{j}}(d_{j})!} \end{split}$$

This last sum simplifies to the indicator $1(m \le n)$, corresponding to the fact that if $n-m \ge 0$, then $d_j = 0$ for j > n-m, and a random permutation in S_{n-m} must have some cycle structure $(d_1, ..., d_{n-m})$. The moments of $C_j^{(n)}$ follow immediately as

$$E(C_i^{(n)})^{[r]} = j^{-r} 1 \{ jr \le n \}$$
 (1.2)

We note for future reference that (1.4) can also be written in the form

$$E\left(\prod_{j=1}^{n} (C_{j}^{(n)})^{[m_{j}]}\right) = E\left(\prod_{j=1}^{n} Z_{j}^{[m_{j}]}\right) 1\left\{\sum_{j=1}^{n} j m_{j} \le n\right\},\tag{1.3}$$

Where the Z_j are independent Poisson-distribution random variables that satisfy $E(Z_j) = 1/j$

The marginal distribution of cycle counts provides a formula for the joint distribution of the cycle counts C_j^n , we find the distribution of C_j^n using a combinatorial approach combined with the inclusion-exclusion formula.

Lemma 1.8. For $1 \le j \le n$,

$$P[C_j^{(n)} = k] = \frac{j^{-k}}{k!} \sum_{l=0}^{\lfloor n/j \rfloor - k} (-1)^l \frac{j^{-l}}{l!}$$
 (1.1)

Proof. Consider the set I of all possible cycles of length j, formed with elements chosen from $\{1,2,...n\}$, so that $|I|=n^{[j]/j}$. For each $\alpha \in I$, consider the "property" G_{α} of having α ; that is,

 G_{α} is the set of permutations $\pi \in S_n$ such that α is one of the cycles of π . We then have $|G_{\alpha}| = (n-j)!$, since the elements of $\{1, 2, ..., n\}$ not in α must be permuted among themselves. To use the inclusion-exclusion formula we need to calculate the term S_r , which is the sum of the probabilities of the r-fold intersection of properties, summing over all sets of r distinct properties. There are two cases to consider. If the r properties are indexed by r cycles having no elements in common, then the intersection specifies how rj elements are moved by the permutation, and there $(n-rj)!1(rj \leq n)$ permutations the There are $n^{[rj]}/(j^r r!)$ intersection. such intersections. For the other case, some two distinct properties name some element in common, so no permutation can have both these properties, and the r -fold intersection is empty. Thus

$$S_r = (n - rj)!1(rj \le n)$$

$$\times \frac{n^{[rj]}}{j^r r!} \frac{1}{n!} = 1 (rj \le n) \frac{1}{j^r r!}$$

Finally, the inclusion-exclusion series for the number of permutations having exactly k properties is

$$\sum_{l\geq 0} (-1)^l \binom{k+l}{l} S_{k+l,}$$

Which simplifies to (1.1) Returning to the original hat-check problem, we substitute j=1 in (1.1) to obtain the distribution of the number of fixed points of a random permutation. For k = 0, 1, ..., n,

$$P[C_1^{(n)} = k] = \frac{1}{k!} \sum_{l=0}^{n-k} (-1)^l \frac{1}{l!},$$
(1.2)

and the moments of $C_1^{(n)}$ follow from (1.2) with j=1. In particular, for $n \ge 2$, the mean and variance of $C_1^{(n)}$ are both equal to 1. The joint distribution of $(C_1^{(n)},...,C_b^{(n)})$ for any $1 \le b \le n$ has an expression similar to (1.7); this too can be derived by inclusion-exclusion. For any $c=(c_1,...,c_b) \in \square$ with $m=\sum ic_i$,

$$P[(C_1^{(n)},...,C_b^{(n)})=c]$$

$$= \left\{ \prod_{i=1}^{b} \left(\frac{1}{i} \right)^{c_i} \frac{1}{c_i!} \right\} \sum_{\substack{l \ge 0 \text{ with} \\ \sum |l_i \le n-m}} (-1)^{l_1 + \dots + l_b} \prod_{i=1}^{b} \left(\frac{1}{i} \right)^{l_i} \frac{1}{l_i!}$$
(1.3)

The joint moments of the first b counts $C_1^{(n)},...,C_b^{(n)}$ can be obtained directly from (1.2) and (1.3) by setting $m_{b+1}=...=m_n=0$



The limit distribution of cycle counts

It follows immediately from Lemma 1.2 that for each fixed j, as $n \rightarrow \infty$,

$$P[C_j^{(n)} = k] \rightarrow \frac{j^{-k}}{k!} e^{-1/j}, \quad k = 0, 1, 2, ...,$$

So that $C_j^{(n)}$ converges in distribution to a random variable Z_j having a Poisson distribution with mean 1/j; we use the notation $C_j^{(n)} \to_d Z_j$ where $Z_j \square P_o(1/j)$ to describe this. Infact, the limit random variables are independent.

Theorem 1.6 The process of cycle counts converges in distribution to a Poisson process of \square with intensity j^{-1} . That is, as $n \to \infty$,

$$(C_1^{(n)}, C_2^{(n)}, ...) \rightarrow_d (Z_1, Z_2, ...)$$
 (1.1)

Where the Z_j , j=1,2,..., are independent Poisson-distributed random variables with $E(Z_j) = \frac{1}{i}$

Proof. To establish the converges in distribution one shows that for each fixed $b \ge 1$, as $n \to \infty$,

$$P[(C_1^{(n)},...,C_b^{(n)})=c] \to P[(Z_1,...,Z_b)=c]$$

Error rates

The proof of Theorem says nothing about the rate of convergence. Elementary analysis can be used to estimate this rate when b=1. Using properties of alternating series with decreasing terms, for k=0,1,...,n,

$$\frac{1}{k!} \left(\frac{1}{(n-k+1)!} - \frac{1}{(n-k+2)!} \right) \le \left| P[C_1^{(n)} = k] - P[Z_1 = k] \right|$$

$$\le \frac{1}{k!(n-k+1)!}$$

It follows that

$$\frac{2^{n+1}}{(n+1)!} \frac{n}{n+2} \le \sum_{k=0}^{n} \left| P[C_1^{(n)} = k] - P[Z_1 = k] \right| \le \frac{2^{n+1} - 1}{(n+1)!}$$
 (1.11)

Since

$$P[Z_1 > n] = \frac{e^{-1}}{(n+1)!} (1 + \frac{1}{n+2} + \frac{1}{(n+2)(n+3)} + \dots) < \frac{1}{(n+1)!},$$

We see from (1.11) that the total variation distance between the distribution $L(C_1^{(n)})$ of $C_1^{(n)}$ and the distribution $L(Z_1)$ of Z_1

B. Infinite State Space Reduction

Any analysis technique that operates by exhaustively examining reachable states must address the fact that in our analysis problem the size of reachable policy states is unbounded, and hence the state space is infinite. In this section we present a subspace of bounded size that was previously identified [12] and that has the property that, given any query, any restriction rule, and any initial state, there exists a reachable state in which the query is violated if and only if there exists a reachable state within the bounded state space in which the query is violated. Given a policy state P, a restriction rule R, and a query O = X.u w A.r. the state space that must be considered consists of all states that are reachable from P under R and that are composed of statements in P and in N, in which N is constructed as follows: $N = \{A.r -D \mid r \text{ 2 Names}(P) \land A,D \text{ 2 Principals}(P) [$ NewPrinc(P,Q)}, NewPrinc(P,Q) is a set of new principals of size 2K, K = |SigRoles(P,Q)|, and SigRoles(P,Q) is the set $\{X.u\}$ [$\{A.r1 \mid A.r - A.r1.r2\}$ 2 P} [$\{B1.r1,B2.r2 \mid A.r - B1.r1 \setminus B2.r2 \mid 2P\}$. It is shown by Li et al. [12] that this set of reachable states has the desired property. We consider the number of new principals used here to be conservative in the sense that any fewer number of principals does not guarantee this desired property.

Establish the asymptotics of $Pigl[A_n(C^{(n)})igr]$ under

conditions (A_0) and (B_{01}) , where

$$A_n(C^{(n)}) = \bigcap_{1 \le i \le n} \bigcap_{r_i + 1 \le j \le r_i} \{C_{ij}^{(n)} = 0\},$$

and
$$\zeta_i = (r_i / r_{id}) - 1 = O(i^{-g})$$
 as $i \to \infty$, for

some g' > 0. We start with the expression

$$P[A_n(C^{(n)})] = \frac{P[T_{0m}(Z) = n]}{P[T_{0m}(Z) = n]}$$

$$\prod_{\substack{1 \le i \le n \\ r_i^{'} + 1 \le j \le r_i}} \left\{ 1 - \frac{\theta}{ir_i} (1 + E_{i0}) \right\}$$
 (1.1)

$$P[T_{0n}(Z') = n]$$

$$= \frac{\theta d}{n} \exp \left\{ \sum_{i>1} [\log(1+i^{-1}\theta d) - i^{-1}\theta d] \right\}$$

$$\left\{1 + O(n^{-1}\varphi_{\{1,2,7\}}(n))\right\}$$
 (1.2)

and

$$P[T_{\alpha}(Z')=n]$$

$$= \frac{\theta d}{n} \exp \left\{ \sum_{i \ge 1} [\log(1 + i^{-1}\theta d) - i^{-1}\theta d] \right\}$$

$$\left\{1 + O(n^{-1}\varphi_{\{1,2,7\}}(n))\right\}$$
 (1.3)



Where $\varphi_{\{1,2,7\}}(n)$ refers to the quantity derived from Z. It thus follows that $P[A_n(C^{(n)})] \square Kn^{-\theta(1-d)}$ for a constant K, depending on Z and the r_i and computable explicitly from (1.1)-(1.3), if Conditions (A_0) and (B_{01}) are satisfied and if $\zeta_i^*=O(i^{-g})$ from some g>0, since, under these circumstances, both $n^{-1}\varphi_{\{1,2,7\}}(n)$ and $n^{-1}\varphi_{\{1,2,7\}}(n)$ tend to zero as $n\to\infty$. In particular, for polynomials and square free polynomials, the relative error in this asymptotic approximation is of order n^{-1} if g>1.

For
$$0 \le b \le n/8$$
 and $n \ge n_0$, with n_0 $d_{TV}(L(C[1,b]), L(Z[1,b]))$ $\le d_{TV}(L(C[1,b]), L(Z[1,b]))$

 $\leq \varepsilon_{(7,7)}(n,b),$

Where $\mathcal{E}_{\{7,7\}}(n,b) = O(b/n)$ under Conditions $(A_0),(D_1)$ and (B_{11}) Since, by the Conditioning Relation,

$$L(C[1,b]|T_{0b}(C)=l)=L(Z[1,b]|T_{0b}(Z)=l),$$

It follows by direct calculation that

$$d_{TV}(L(C[1,b]), L(Z[1,b]))$$

$$= d_{TV}(L(T_{0b}(C)), L(T_{0b}(Z)))$$

$$= \max_{A} \sum_{r \in A} P[T_{0b}(Z) = r]$$

$$\left\{1 - \frac{P[T_{bn}(Z) = n - r]}{P[T_{0n}(Z) = n]}\right\}$$
(1.4)

Suppressing the argument Z from now on, we thus obtain

$$\begin{split} d_{TV}(L(C[1,b]),L(Z[1,b])) \\ &= \sum_{r \geq 0} P[T_{0b} = r] \left\{ 1 - \frac{P[T_{bn} = n - r]}{P[T_{0n} = n]} \right\}_{+} \\ &\leq \sum_{r > n/2} P[T_{0b} = r] + \sum_{r = 0}^{[n/2]} \frac{P[T_{0b} = r]}{P[T_{0b} = n]} \\ &\times \left\{ \sum_{s = 0}^{n} P[T_{0b} = s](P[T_{bn} = n - s] - P[T_{bn} = n - r] \right\}_{+} \\ &\leq \sum_{r > n/2} P[T_{0b} = r] + \sum_{r = 0}^{[n/2]} P[T_{0b} = r] \end{split}$$

$$\times \sum_{s=0}^{\lfloor n/2 \rfloor} P[T_{0b} = s] \frac{\left\{ P[T_{bn} = n - s] - P[T_{bn} = n - r] \right\}}{P[T_{0n} = n]}$$

$$+ \sum_{s=0}^{\lfloor n/2 \rfloor} P[T_{0b} = r] \sum_{s=\lfloor n/2 \rfloor + 1}^{n} P[T = s] P[T_{bn} = n - s] / P[T_{0n} = n]$$

The first sum is at most $2n^{-1}ET_{0b}$; the third is bound by

$$\begin{split} &(\max_{n/2 < s \le n} P[T_{0b} = s]) / P[T_{0n} = n] \\ & \le \frac{2\varepsilon_{\{10.5(1)\}}(n/2, b)}{n} \frac{3n}{\theta P_{\theta}[0, 1]}, \\ & \frac{3n}{\theta P_{\theta}[0, 1]} 4n^{-2} \phi_{\{10.8\}}^*(n) \sum_{r=0}^{[n/2]} P[T_{0b} = r] \sum_{s=0}^{[n/2]} P[T_{0b} = s] \frac{1}{2} |r - s| \\ & \le \frac{12\phi_{\{10.8\}}^*(n)}{\theta P_{\theta}[0, 1]} \frac{ET_{0b}}{n} \end{split}$$

Hence we may take

$$\varepsilon_{\{7,7\}}(n,b) = 2n^{-1}ET_{0b}(Z) \left\{ 1 + \frac{6\phi_{\{10.8\}}^*(n)}{\theta P_{\theta}[0,1]} \right\} P$$

$$+ \frac{6}{\theta P_{\theta}[0,1]} \varepsilon_{\{10.5(1)\}}(n/2,b)$$
 (1.5)

Required order under Conditions $(A_0), (D_1)$ and $(B_{\!\scriptscriptstyle 11}), \ {
m if} \ S(\infty) < \infty. \ {
m If not}, \ \phi_{\scriptscriptstyle \{10.8\}}^*\left(n
ight) \ {
m can be}$ replaced by $\phi_{\{10.11\}}^*(n)$ in the above, which has the required order, without the restriction on the r_i implied by $S(\infty) < \infty$. Examining the Conditions $(A_0),(D_1)$ and (B_{11}) , it is perhaps surprising to find that (B_{11}) is required instead of just (B_{01}) ; that is, that we should need $\sum_{i>2} l \varepsilon_{il} = O(i^{-a_1})$ to hold for some $a_1 > 1$. A first observation is that a similar problem arises with the rate of decay of \mathcal{E}_{i1} as well. For this reason, n_1 is replaced by n_1 . This makes it possible to replace condition (A_1) by the weaker pair of conditions (A_0) and (D_1) in the eventual assumptions needed for $\mathcal{E}_{\{7,7\}}(n,b)$ to be of order O(b/n); the decay rate requirement of order $i^{-1-\gamma}$ is shifted from \mathcal{E}_{i1} itself to its first difference. This is needed to obtain the right approximation error for the random mappings since example. However, all the applications make far more stringent assumptions



about the $\varepsilon_{i1}, l \geq 2$, than are made in (B_{11}) . The critical point of the proof is seen where the initial difference $P[T_{bn}^{(m)} = s] - P[T_{bn}^{(m)} = s + 1]$ The $\mathcal{E}_{f_{10,10}}(n)$, which should be small, contains a far tail element from n_1 of the form $\phi_1^{\theta}(n) + u_1^*(n)$, which is only small if $a_1 > 1$, being otherwise of order $O(n^{1-a_1+\delta})$ for any $\delta > 0$, since $a_2 > 1$ is in any case assumed. For $s \ge n/2$, this gives rise to a contribution of order $O(n^{-1-a_1+\delta})$ in the estimate of the $P[T_{bn} = s] - P[T_{bn} = s + 1],$ which, in remainder of the proof, is translated into a contribution of order $O(tn^{-1-a_1+\delta})$ for differences of the form $P[T_{bn} = s] - P[T_{bn} = s+1]$, finally leading to a contribution of order $bn^{-a_1+\delta}$ for any $\delta > 0$ in $\mathcal{E}_{\{7,7\}}(n,b)$. Some improvement would seem to be possible, defining the function g by $g(w) = 1_{\{w=s\}} - 1_{\{w=s+t\}}$, differences that are of the form $P[T_{bn} = s] - P[T_{bn} = s + t]$ can be directly estimated, at a cost of only a single contribution of the form $\phi_1^{\theta}(n) + u_1^*(n)$. Then, iterating the cycle, in which one estimate of a difference in point probabilities is improved to an estimate of smaller order, a bound of the form

$$\begin{aligned} \left| P[T_{bn} = s] - P[T_{bn} = s + t] \right| &= O(n^{-2}t + n^{-1 - a_1 + \delta}) \\ \text{for any } \delta > 0 \text{ could perhaps be attained, leading to} \\ \text{a final error estimate in order } O(bn^{-1} + n^{-a_1 + \delta}) \\ \text{for any } \delta > 0 \text{, to replace } \varepsilon_{\{7.7\}}(n,b). \text{ This would} \\ \text{be of the ideal order } O(b/n) \text{ for large enough } b, \\ \text{but would still be coarser for small } b. \end{aligned}$$

With b and n as in the previous section, we wish to show that

$$\left| d_{TV}(L(C[1,b]), L(Z[1,b])) - \frac{1}{2}(n+1)^{-1} \left| 1 - \theta \right| E \left| T_{0b} - ET_{0b} \right| \right| \le \varepsilon_{\{7,8\}}(n,b),$$

Where

$$\begin{split} & \mathcal{E}_{\{7.8\}}(n,b) = O(n^{-1}b[n^{-1}b + n^{-\beta_{12} + \delta}]) \quad \text{for any} \\ & \mathcal{S} > 0 \quad \text{under Conditions} \quad (A_0), (D_1) \quad \text{and} \quad (B_{12}), \\ & \text{with} \quad \beta_{12} \quad \text{. The proof uses sharper estimates. As} \\ & \text{before, we begin with the formula} \end{split}$$

$$\begin{aligned} d_{TV}(L(C[1,b]), L(Z[1,b])) \\ &= \sum_{r \ge 0} P[T_{0b} = r] \left\{ 1 - \frac{P[T_{bn} = n - r]}{P[T_{0n} = n]} \right\}_{+} \end{aligned}$$

Now we observe that

$$\left| \sum_{r\geq 0} P[T_{0b} = r] \left\{ 1 - \frac{P[T_{bn} = n - r]}{P[T_{0n} = n]} \right\}_{+} - \sum_{r=0}^{[n/2]} \frac{P[T_{0b} = r]}{P[T_{0n} = n]} \right| \times \left| \sum_{s=[n/2]+1}^{n} P[T_{0b} = s] (P[T_{bn} = n - s] - P[T_{bn} = n - r]) \right| \le 4n^{-2} E T_{0b}^{2} + \left(\max_{n/2 < s \le n} P[T_{0b} = s] \right) / P[T_{0n} = n] + P[T_{0b} > n / 2] \right| \le 8n^{-2} E T_{0b}^{2} + \frac{3\varepsilon_{\{10.5(2)\}}(n / 2, b)}{\theta P_{0}[0.1]},$$

$$(1.1)$$

We have

$$\Big| \sum_{r=0}^{[n/2]} \frac{P[T_{0b} = r]}{P[T_{0n} = n]}$$

$$\times \left\{ \sum_{s=0}^{\lfloor n/2 \rfloor} P[T_{0b} = s] (P[T_{bn} = n - s] - P[T_{bn} = n - r] \right\}_{+}$$

$$- \left\{ \sum_{s=0}^{\lfloor n/2 \rfloor} P[T_{0b} = s] \frac{(s - r)(1 - \theta)}{n + 1} P[T_{0n} = n] \right\}_{+}$$

$$\leq \frac{1}{n^{2}P[T_{0n} = n]} \sum_{r \geq 0} P[T_{0b} = r] \sum_{s \geq 0} P[T_{0b} = s] |s - r|
\times \left\{ \varepsilon_{\{10.14\}}(n,b) + 2(r \vee s) |1 - \theta| n^{-1} \left\{ K_{0}\theta + 4\phi_{\{10.8\}}^{*}(n) \right\} \right\}
\leq \frac{6}{\theta n P_{\theta}[0,1]} E T_{0b} \varepsilon_{\{10.14\}}(n,b)
+ 4 |1 - \theta| n^{-2} E T_{0b}^{2} \left\{ K_{0}\theta + 4\phi_{\{10.8\}}^{*}(n) \right\}
\left(\frac{3}{\theta n P_{\theta}[0,1]}\right) \right\},$$
(1.2)

The approximation in (1.2) is further simplified by noting that

$$\sum_{r=0}^{[n/2]} P[T_{0b} = r] \left\{ \sum_{s=0}^{[n/2]} P[T_{0b} = s] \frac{(s-r)(1-\theta)}{n+1} \right\}_{+}$$

$$-\left\{\sum_{s=0} P[T_{0b} = s] \frac{(s-r)(1-\theta)}{n+1}\right\}_{\perp} = \left\{\frac{(s-r)(1-\theta)}{n+1}\right\}_{\perp}$$

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$$\leq \sum_{r=0}^{\lfloor n/2 \rfloor} P[T_{0b} = r] \sum_{s > \lfloor n/2 \rfloor} P[T_{0b} = s] \frac{(s-r)|1-\theta|}{n+1} \\
\leq |1-\theta| n^{-1} E(T_{0b} 1\{T_{0b} > n/2\}) \leq 2|1-\theta| n^{-2} ET_{0b}^2, \quad (1.3)$$
and then by observing that

$$\sum_{r>\{n/2\}} P[T_{0b} = r] \left\{ \sum_{s\geq 0} P[T_{0b} = s] \frac{(s-r)(1-\theta)}{n+1} \right\}$$

$$\leq n^{-1} \left| 1 - \theta \right| (ET_{0b}P[T_{0b} > n/2] + E(T_{0b}1\{T_{0b} > n/2\}))$$

$$\leq 4 \left| 1 - \theta \right| n^{-2}ET_{0b}^{2}$$
(1.4)

Combining the contributions of (1.2) – (1.3), we thus find tha $d_{TV}(L(C[1,b]), L(Z[1,b]))$

$$-(n+1)^{-1} \sum_{r\geq 0} P[T_{0b} = r] \left\{ \sum_{s\geq 0} P[T_{0b} = s](s-r)(1-\theta) \right\}_{+}$$

$$\leq \mathcal{E}_{(n+1)}(n,h)$$

$$= \frac{3}{\theta P_{c}[0.1]} \Big\{ \varepsilon_{\{10.5(2)\}}(n/2,b) + 2n^{-1}ET_{0b}\varepsilon_{\{10.14\}}(n,b) \Big\}$$

$$+2n^{-2}ET_{0b}^{2}\left\{4+3\left|1-\theta\right|+\frac{24\left|1-\theta\right|\phi_{\{10.8\}}^{*}(n)}{\theta P_{\theta}[0,1]}\right\} \tag{1.5}$$

The quantity $\mathcal{E}_{\{7.8\}}(n,b)$ is seen to be of the order claimed under Conditions $(A_0),(D_1)$ and (B_{12}) , provided that $S(\infty)<\infty$; this supplementary condition can be removed if $\phi_{\{10.8\}}^*(n)$ is replaced by $\phi_{\{10.11\}}^*(n)$ in the definition of $\mathcal{E}_{\{7.8\}}(n,b)$, has the required order without the restriction on the r_i implied by assuming that $S(\infty)<\infty$. Finally, a direct calculation now shows that

$$\sum_{r\geq 0} P[T_{0b} = r] \left\{ \sum_{s\geq 0} P[T_{0b} = s](s-r)(1-\theta) \right\}_{+}$$
$$= \frac{1}{2} |1-\theta| E |T_{0b} - ET_{0b}|$$

C. Unrestricted Role Reduction

We now present a result that in many cases enables us to further reduce the size of the state space that must be explored. The idea is that given an initial state P, we can often perform the analysis using a smaller initial state and

obtain identical results. It is important to note that this is significantly different than the Lower Bound LB(P) and Upper

Bound UB(P) programs described in [12]. The LB(P) program removes all statements defining a role C.r 62 SR from P for the purpose of evaluating a query of the form $X.u\ w\ \{B\mid B\ 2\ Principals\}$, whereas the UB(P) program adds a special principal

_ representing all principals to each growth unrestricted role for the purposes of evaluating a query of the form {B | B 2 Principals} w A.r. Our reduction program removes all statements defining a role C.r 62 SR ^ C.r 62 GR for the purpose of evaluating a query of the form X.u w A.r. We define _R, a binary relation over policy states, by

P1_R P2 if and only if P1_7!R P2 and P2_7!R P1. Note that _R is an equivalence relation: (1) it is reflexive, as P_7!R P for all P; (2) it is symmetric by construction; (3) it is transitive because _7!R is transitive. We also define the core of P, coreR(P), to be the subset of P consisting of those statements that define growthrestricted or shrink-restricted roles.

Theorem 1. Given any policy states P1 and P2 and any restriction rule R, P1 _R P2 if and only if coreR(P1) = coreR(P2).

Proof. The "if" direction is straightforward: starting from any state P, it is clear that any other state with the same core is reachable from P. For the "only if" direction, suppose coreR(P1) 6= coreR(P2). There are two cases, depending on whether the cores differ in statements that define roles that are growthrestricted or roles that are shrink restricted. If coreR(P1)\coreR(P2) contains a statement defining a growth restricted role, then P2 _ 7!R P1 does not hold. If coreR(P1)\coreR(P2) contains a statement defining a shrink restricted role, then P1 7!R P2 does not hold. Given this result, it is clear that the least state (under the subset ordering) that is equivalent to a given P is coreR(P). It follows from the definition of R and the transitivity of 7!R that for all P0, P 7!R P0 if and only if coreR(P) 7!R P0. Moreover, coreR(P) is the least set for which this is true. Thus, given an initial state P, it is both correct and more efficient to perform our analysis using coreR (P) as the initial state instead. Definition 3. Given a policy P and restriction rule R, URR(P, R) = coreR(P)

Example 1.0. Consider the $O = (0, ..., 0) \in \square^n$. For an arbitrary vector r, the coordinates of the point x = O + r are equal to the respective coordinates of $r: x = (x^1, ..., x^n)$ and $r = (x^1, ..., x^n)$. The vector r such as in the example is called the position vector or the radius vector of the point X. (Or, in greater detail: r is the radius-vector of x w.r.t an origin O). Points are frequently specified by their radiusvectors. This presupposes the choice of O as the "standard origin". Let us summarize. We have considered \square ⁿ and interpreted its elements in two ways: as points and as vectors. Hence we may say that we leading with the two copies of \square^n : $\square^n =$ $\square^n = \{\text{vectors}\}\$ Operations with vectors: multiplication by a number, addition. Operations with points and vectors: adding a vector



to a point (giving a point), subtracting two points (giving a vector). \square * treated in this way is called an n-dimensional affine space. (An "abstract" affine space is a pair of sets, the set of points and the set of vectors so that the operations as above are defined axiomatically). Notice that vectors in an affine space are also known as "free vectors". Intuitively, they are not fixed at points and "float freely" in space. From \square " considered as an affine space we can precede in two opposite directions: \square as an Euclidean space \Leftarrow \square as an affine space $\Rightarrow \Box^n$ as a manifold. Going to the left means introducing some extra structure which will make the geometry richer. Going to the right means forgetting about part of the affine structure; going further in this direction will lead us to the so-called "smooth (or differentiable) manifolds". The theory of differential forms does not require any extra geometry. So our natural direction is to the right. The Euclidean structure, however, is useful for examples and applications. So let us say a few words about it:

Remark 1.0. Euclidean geometry. In \Box ⁿ considered as an affine space we can already do a good deal of geometry. For example, we can consider lines and planes, and quadric surfaces like an ellipsoid. However, we cannot discuss such things as "lengths", "angles" or "areas" and "volumes". To be able to do so, we have to introduce some more definitions, making \Box ⁿ a Euclidean space. Namely, we define the length of a vector $a = (a^1, ..., a^n)$ to be

$$|a| := \sqrt{(a^1)^2 + \dots + (a^n)^2}$$
 (1)

After that we can also define distances between points as follows:

$$d(A,B) := \left| \overrightarrow{AB} \right| \tag{2}$$

One can check that the distance so defined possesses natural properties that we expect: is it always non-negative and equals zero only for coinciding points; the distance from A to B is the same as that from B to A (symmetry); also, for three points, A, B and C, we have $d(A,B) \le d(A,C) + d(C,B)$ (the "triangle inequality"). To define angles, we first introduce the scalar product of two vectors

$$(a,b) := a^1 b^1 + \dots + a^n b^n$$
 (3)

Thus $|a| = \sqrt{(a,a)}$. The scalar product is

also denote by dot: a.b = (a,b), and hence is often referred to as the "dot product". Now, for nonzero vectors, we define the angle between them by the equality

$$\cos \alpha := \frac{(a,b)}{|a||b|} \tag{4}$$

The angle itself is defined up to an integral multiple of 2π . For this definition to be consistent we have to ensure that the r.h.s. of (4) does not exceed 1 by the absolute value. This follows from the inequality

$$(a,b)^2 \le |a|^2 |b|^2$$
 (5)

known as the Cauchy-Bunyakovsky-Schwarz inequality (various combinations of these three names are applied in different books). One of the ways of proving (5) is to consider the scalar square of the linear combination a+tb, where $t \in R$. As $(a+tb,a+tb) \geq 0$ is a quadratic polynomial in t which is never negative, its discriminant must be less or equal zero. Writing this explicitly yields (5). The triangle inequality for distances also follows from the inequality (5).

Example 1.1. Consider the function $f(x) = x^i$ (the i-th coordinate). The linear function dx^i (the differential of x^i) applied to an arbitrary vector h is simply h^i . From these examples follows that we can rewrite df as

$$df = \frac{\partial f}{\partial x^1} dx^1 + \dots + \frac{\partial f}{\partial x^n} dx^n, \tag{1}$$

which is the standard form. Once again: the partial derivatives in (1) are just the coefficients (depending on x); $dx^1, dx^2, ...$ are linear functions giving on an arbitrary vector h its coordinates $h^1, h^2, ...$, respectively. Hence

$$df(x)(h) = \partial_{hf(x)} = \frac{\partial f}{\partial x^1} h^1 + \frac{\partial f}{\partial x^2} h^2 + \frac$$

$$... + \frac{\partial f}{\partial x^n} h^n, \qquad (2)$$

Theorem 1.7. Suppose we have a parametrized curve $t \mapsto x(t)$ passing through $x_0 \in \square^n$ at $t = t_0$ and with the velocity vector $x(t_0) = \upsilon$ Then $\frac{df(x(t))}{dt}(t_0) = \partial_{\upsilon} f(x_0) = df(x_0)(\upsilon) \tag{1}$

Proof. Indeed, consider a small increment of the parameter $t:t_0\mapsto t_0+\Delta t$, Where $\Delta t\mapsto 0$. On the other hand, we have $f(x_0+h)-f(x_0)=df(x_0)(h)+\beta(h)|h|$ for an arbitrary vector h, where $\beta(h)\to 0$ when



 $h \rightarrow 0$. Combining it together, for the increment of f(x(t)) we obtain

$$f(x(t_0 + \Delta t) - f(x_0)$$

$$= df(x_0)(\upsilon.\Delta t + \alpha(\Delta t)\Delta t)$$

$$+\beta(\upsilon.\Delta t + \alpha(\Delta t)\Delta t).$$
 $\upsilon\Delta t + \alpha(\Delta t)\Delta t$

$$= df(x_0)(\upsilon).\Delta t + \gamma(\Delta t)\Delta t$$

For a certain $\gamma(\Delta t)$ such that $\gamma(\Delta t) \to 0$ when $\Delta t \to 0$ (we used the linearity of $df(x_0)$). By the definition, this means that the derivative of f(x(t)) at $t=t_0$ is exactly $df(x_0)(\upsilon)$. The statement of the theorem can be expressed by a simple formula:

$$\frac{df(x(t))}{dt} = \frac{\partial f}{\partial x^1} x^1 + \dots + \frac{\partial f}{\partial x^n} x^n$$
 (2)

To calculate the value Of df at a point x_0 on a given vector v one can take an arbitrary curve passing Through x_0 at t_0 with v as the velocity vector at t_0 and calculate the usual derivative of f(x(t)) at $t=t_0$.

Theorem 1.8. For functions $f, g: U \to \square$, $U \subset \square^n$.

$$d(f+g) = df + dg \tag{1}$$

$$d(fg) = df \cdot g + f \cdot dg \tag{2}$$

Proof. Consider an arbitrary point x_0 and an arbitrary vector v0 stretching from it. Let a curve v(t)1 be such that v(t)2 = v3 and v(t)3 = v4. Hence

$$d(f+g)(x_0)(v) = \frac{d}{dt}(f(x(t)) + g(x(t)))$$

at $t = t_0$ and

$$d(fg)(x_0)(v) = \frac{d}{dt}(f(x(t))g(x(t)))$$

at $t=t_0$ Formulae (1) and (2) then immediately follow from the corresponding formulae for the usual derivative Now, almost without change the theory generalizes to functions taking values in \square m instead of \square . The only difference is that now the differential of a map $F:U \to \square$ m at a point x will be a linear function taking vectors in \square n to vectors in \square m (instead of \square). For an arbitrary vector $h \in \square$ n ,

$$F(x+h) = F(x) + dF(x)(h) + \beta(h)|h|$$
(3)

Where $\beta(h) \to 0$ when $h \to 0$. We have $dF = (dF^1, ..., dF^m)$ and

$$dF = \frac{\partial F}{\partial x^{1}} dx^{1} + \dots + \frac{\partial F}{\partial x^{n}} dx^{n}$$

$$= \begin{pmatrix} \frac{\partial F^{1}}{\partial x^{1}} \dots \frac{\partial F^{1}}{\partial x^{n}} \\ \dots & \dots & \dots \\ \frac{\partial F^{m}}{\partial x^{1}} \dots \frac{\partial F^{m}}{\partial x^{n}} \end{pmatrix} \begin{pmatrix} dx^{1} \\ \dots \\ dx^{n} \end{pmatrix}$$

$$(4)$$

In this matrix notation we have to write vectors as vector-columns.

Theorem 1.9. For an arbitrary parametrized curve x(t) in \square^n , the differential of a map $F: U \to \square^m$ (where $U \subset \square^n$) maps the velocity vector x(t) to the velocity vector of the curve F(x(t)) in \square^m :

$$\frac{dF(x(t))}{dt} = dF(x(t))(x(t)) \tag{1}$$

Proof. By the definition of the velocity vector,

$$x(t + \Delta t) = x(t) + \dot{x}(t) \cdot \Delta t + \alpha(\Delta t) \Delta t \tag{2}$$

Where $\alpha(\Delta t) \to 0$ when $\Delta t \to 0$. By the definition of the differential,

$$F(x+h) = F(x) + dF(x)(h) + \beta(h)|h$$
 (3)

Where $\beta(h) \rightarrow 0$ when $h \rightarrow 0$, we obtain

$$F(x(t + \Delta t)) = F(x + \underbrace{x(t).\Delta t + \alpha(\Delta t)\Delta t}_{h})$$

$$= F(x) + dF(x)(x(t)\Delta t + \alpha(\Delta t)\Delta t) +$$

$$\beta(x(t)\Delta t + \alpha(\Delta t)\Delta t). \left| x(t)\Delta t + \alpha(\Delta t)\Delta t \right|$$

$$= F(x) + dF(x)(x(t)\Delta t + \gamma(\Delta t)\Delta t$$

For some $\gamma(\Delta t) \rightarrow 0$ when $\Delta t \rightarrow 0$. This

precisely means that dF(x)x(t) is the velocity vector of F(x). As every vector attached to a point can be viewed as the velocity vector of some curve passing through this point, this theorem gives a clear geometric picture of dF as a linear map on vectors.



Theorem 1.10 Suppose we have two maps $F: U \to V$ and $G: V \to W$, where $U \subset \square^n, V \subset \square^m, W \subset \square^p$ (open domains). Let $F: x \mapsto y = F(x)$. Then the differential of the composite map $GoF: U \to W$ is the composition of the differentials of F and G:

$$d(GoF)(x) = dG(y)odF(x)$$
 (4)

Proof. We can use the description of the differential .Consider a curve x(t) in \square^n with the

velocity vector x. Basically, we need to know to which vector in \Box p it is taken by d(GoF). the curve (GoF)(x(t) = G(F(x(t))). By the same theorem, it equals the image under dG of the Anycast Flow vector to the curve F(x(t)) in \Box m . Applying the theorem once again, we see that the velocity vector to the curve F(x(t)) is the image

under dF of the vector x(t) . Hence $d(GoF)(x) = dG(dF(x)) \qquad \text{for an arbitrary}$ vector x .

Corollary 1.0. If we denote coordinates in \square by $(x^1,...,x^n)$ and in \square by $(y^1,...,y^m)$, and write

$$dF = \frac{\partial F}{\partial x^{1}} dx^{1} + \dots + \frac{\partial F}{\partial x^{n}} dx^{n}$$
 (1)

$$dG = \frac{\partial G}{\partial y^{1}} dy^{1} + \dots + \frac{\partial G}{\partial y^{n}} dy^{n}, \qquad (2)$$

Then the chain rule can be expressed as follows:

$$d(GoF) = \frac{\partial G}{\partial y^{1}} dF^{1} + \dots + \frac{\partial G}{\partial y^{m}} dF^{m}, \qquad (3)$$

Where dF^i are taken from (1). In other words, to get d(GoF) we have to substitute into (2) the expression for $dy^i = dF^i$ from (3). This can also be expressed by the following matrix formula:

$$d(GoF) = \begin{pmatrix} \frac{\partial G^{1}}{\partial y^{1}} & \dots & \frac{\partial G^{1}}{\partial y^{m}} \\ \dots & \dots & \dots \\ \frac{\partial G^{p}}{\partial y^{1}} & \dots & \frac{\partial G^{p}}{\partial y^{m}} \end{pmatrix} \begin{pmatrix} \frac{\partial F^{1}}{\partial x^{1}} & \dots & \frac{\partial F^{1}}{\partial x^{n}} \\ \dots & \dots & \dots & \dots \\ \frac{\partial F^{m}}{\partial x^{1}} & \dots & \frac{\partial F^{m}}{\partial x^{n}} \end{pmatrix} \begin{pmatrix} dx^{1} & \dots & dx^{n} \\ \dots & \dots & \dots & \dots \\ dx^{n} & \dots & \dots \end{pmatrix}$$
(4)

i.e., if dG and dF are expressed by matrices of partial derivatives, then d(GoF) is expressed by the product of these matrices. This is often written

$$\left(\frac{\partial z^{1}}{\partial x^{1}} \dots \frac{\partial z^{1}}{\partial x^{n}}\right) = \left(\frac{\partial z^{1}}{\partial y^{1}} \dots \frac{\partial z^{1}}{\partial y^{m}}\right) \\
\dots \dots \dots \\
\left(\frac{\partial z^{p}}{\partial x^{1}} \dots \frac{\partial z^{p}}{\partial x^{n}}\right) = \left(\frac{\partial z^{1}}{\partial y^{1}} \dots \frac{\partial z^{1}}{\partial y^{m}} \dots \frac{\partial z^{p}}{\partial y^{m}}\right)$$

$$\begin{pmatrix}
\frac{\partial y^{1}}{\partial x^{1}} & \dots & \frac{\partial y^{1}}{\partial x^{n}} \\
\dots & \dots & \dots \\
\frac{\partial y^{m}}{\partial x^{1}} & \dots & \frac{\partial y^{m}}{\partial x^{n}}
\end{pmatrix}, (5)$$

Or

$$\frac{\partial z^{\mu}}{\partial x^{a}} = \sum_{i=1}^{m} \frac{\partial z^{\mu}}{\partial y^{i}} \frac{\partial y^{i}}{\partial x^{a}},$$
 (6)

Where it is assumed that the dependence of $y \in \square^m$ on $x \in \square^n$ is given by the map F, the dependence of $z \in \square^p$ on $y \in \square^m$ is given by the map G, and the dependence of $z \in \square^p$ on $x \in \square^n$ is given by the composition GoF.

Definition 1.6. Consider an open domain $U \subset \square^n$. Consider also another copy of \square^n , denoted for distinction \square_y^n , with the standard coordinates $(y^1...y^n)$. A system of coordinates in the open domain U is given by a map $F:V \to U$, where $V \subset \square_y^n$ is an open domain of \square_y^n , such that the following three conditions are satisfied:

- (1) F is smooth;
- (2) F is invertible;
- (3) $F^{-1}: U \to V$ is also smooth

The coordinates of a point $x \in U$ in this system are the standard coordinates of $F^{-1}(x) \in \square_{y}^{n}$

In other words,

$$F:(y^1...,y^n) \mapsto x = x(y^1...,y^n)$$
 (1)

Here the variables $(y^1,...,y^n)$ are the "new" coordinates of the point x



Example 1.2. Consider a curve in \square ² specified in polar coordinates as

$$x(t): r = r(t), \varphi = \varphi(t) \tag{1}$$

We can simply use the chain rule. The map $t\mapsto x(t)$ can be considered as the composition of the maps $t\mapsto (r(t),\varphi(t)),(r,\varphi)\mapsto x(r,\varphi)$. Then, by the chain rule, we have

$$x = \frac{dx}{dt} = \frac{\partial x}{\partial r}\frac{dr}{dt} + \frac{\partial x}{\partial \varphi}\frac{d\varphi}{dt} = \frac{\partial x}{\partial r}r + \frac{\partial x}{\partial \varphi}\varphi$$
 (2)

Here r and φ are scalar coefficients depending on t, whence the partial derivatives $\frac{\partial x}{\partial r}, \frac{\partial x}{\partial \varphi}$ are vectors depending on point in \Box ². We can compare this with the formula in the "standard" coordinates:

 $x = e_1 x + e_2 y$. Consider the vectors $\frac{\partial x}{\partial r}, \frac{\partial x}{\partial \varphi}$.

Explicitly we have

$$\frac{\partial x}{\partial r} = (\cos \varphi, \sin \varphi)$$

$$\frac{\partial x}{\partial \varphi} = (-r \sin \varphi, r \cos \varphi)$$
(3)

From where it follows that these vectors make a basis at all points except for the origin (where r=0). It is instructive to sketch a picture, drawing vectors corresponding to a point as starting from that point. Notice that $\frac{\partial x}{\partial r}, \frac{\partial x}{\partial \varphi}$ are, respectively, the velocity vectors for the curves $r\mapsto x(r,\varphi)$ $(\varphi=\varphi_0\ fixed)$ and $\varphi\mapsto x(r,\varphi)$ $(r=r_0\ fixed)$. We can conclude that for an arbitrary curve given in polar coordinates the velocity vector will have components (r,φ) if as a basis we take $e_r \coloneqq \frac{\partial x}{\partial r}, e_{\varphi} \coloneqq \frac{\partial x}{\partial \varphi}$:

$$x = e_r r + e_{\varphi} \varphi \tag{5}$$

A characteristic feature of the basis e_r, e_{φ} is that it is not "constant" but depends on point. Vectors "stuck to points" when we consider curvilinear coordinates.

Proposition 1.3. The velocity vector has the same appearance in all coordinate systems.

Proof. Follows directly from the chain rule and the transformation law for the basis e_i . In particular, the elements of the basis $e_i = \frac{\partial x}{\partial x^i}$ (originally, a formal notation) can be understood directly as the velocity vectors of the coordinate lines

 $x^i \mapsto x(x^1,...,x^n)$ (all coordinates but x^i are fixed). Since we now know how to handle velocities in arbitrary coordinates, the best way to treat the differential of a map $F: \Box^n \to \Box^m$ is by its action on the velocity vectors. By definition, we set

$$dF(x_0): \frac{dx(t)}{dt}(t_0) \mapsto \frac{dF(x(t))}{dt}(t_0) \tag{1}$$

Now $dF(x_0)$ is a linear map that takes vectors attached to a point $x_0 \in \square^n$ to vectors attached to the point $F(x) \in \square^m$

$$dF = \frac{\partial F}{\partial x^{1}} dx^{1} + \dots + \frac{\partial F}{\partial x^{n}} dx^{n}$$

$$(e_{1}, \dots, e_{m}) \begin{pmatrix} \frac{\partial F^{1}}{\partial x^{1}} \dots \frac{\partial F^{1}}{\partial x^{n}} \\ \dots \dots \dots \\ \frac{\partial F^{m}}{\partial x^{1}} \dots \frac{\partial F^{m}}{\partial x^{n}} \end{pmatrix} \begin{pmatrix} dx^{1} \\ \dots \\ dx^{n} \end{pmatrix}, \tag{2}$$

In particular, for the differential of a function we always have

$$df = \frac{\partial f}{\partial x^1} dx^1 + \dots + \frac{\partial f}{\partial x^n} dx^n, \tag{3}$$

Where x^i are arbitrary coordinates. The form of the differential does not change when we perform a change of coordinates.

Example 1.3 Consider a 1-form in \Box ² given in the standard coordinates:

A = -ydx + xdy In the polar coordinates we will have $x = r\cos\varphi$, $y = r\sin\varphi$, hence

$$dx = \cos\varphi dr - r\sin\varphi d\varphi$$

$$dy = \sin \varphi dr + r \cos \varphi d\varphi$$

Substituting into A, we get

$$A = -r \sin \varphi (\cos \varphi dr - r \sin \varphi d\varphi)$$

$$+r\cos\varphi(\sin\varphi dr+r\cos\varphi d\varphi)$$

$$= r^2 (\sin^2 \varphi + \cos^2 \varphi) d\varphi = r^2 d\varphi$$

Hence $A=r^2d\varphi$ is the formula for A in the polar coordinates. In particular, we see that this is again a 1-form, a linear combination of the differentials of coordinates with functions as coefficients. Secondly, in a more conceptual way, we can define a 1-form in a domain U as a linear function on vectors at every point of U:

$$\omega(\upsilon) = \omega_1 \upsilon^1 + \dots + \omega_n \upsilon^n, \tag{1}$$



If $\upsilon = \sum e_i \upsilon^i$, where $e_i = \frac{\partial x}{\partial x^i}$. Recall that the differentials of functions were defined as linear functions on vectors (at every point), and $dx^{i}(e_{j}) = dx^{i} \left(\frac{\partial x}{\partial x^{j}} \right) = \delta_{j}^{i}$ at

every point X.

Theorem 1.9. For arbitrary 1-form ω and path γ , the integral ω does not change if we change parametrization of γ provide the orientation remains the same.

Proof: Consider
$$\left\langle \omega(x(t)), \frac{dx}{dt} \right\rangle$$
 and $\left\langle \omega(x(t(t'))), \frac{dx}{dt'} \right\rangle$ As $\left\langle \omega(x(t(t'))), \frac{dx}{dt'} \right\rangle = \left| \left\langle \omega(x(t(t'))), \frac{dx}{dt'} \right\rangle \cdot \frac{dt}{dt'} \right\rangle$

D. Cone of Influence Reduction

A given security policy P may contain statements that do not affect the membership of queried roles. Such extraneous statements can safely be filtered from the policy model in order to reduce the size of the problem. This reduction removes statements that are said to be outside of a role's cone of influence. This reduction is particularly significant because it has the potential to remove linked or intersection inclusion type statements that contribute to a larger number of principals in the model. Given a set of roles, _, the following definition constructs a set of roles in P. A role, _, is in the constructed set if the membership of some role in _ depends in some way on the membership of _. This dependency is reflective of relationship established between roles via RT statements. We call this set of roles DefRoles.

Definition 4. Let _ and M be sets of roles, and P be a policy. We define DefRoles(P,_,M) to be the least

roles O satisfying the following conditions:

- __ O (_ 2 O ^_ B.r1 2 P ^ B.r1 62M)) B.r1 2 O (_ 2 O ^_ B.r1.r2 2 P ^ D 2 Principals)) ((B.r1 62M) B.r1 2 O) ^ (D.r2 62M) D.r2 2 O))
- (_ 2 O ^ _ B.r1 \ C.r2 2 P)) ((B.r1 62 M) B.r1 2 O) ^ (C.r2 62M) C.r2 2 O))

Using DefRoles, we define the Cone of Influence (COI) as a policy constructed from those statements that define roles in DefRoles. In other words, we retain only those statements that influence the answer to an RCPI.

Definition 5. Given an RCPI hP,R,X.u w A.ri, we define

COI(hP,R,X.u w A.ri)

 $= P_DefRoles(P, \{X.u\}, SR)[DefRoles(P, \{A.r\}, GR)]$

Theorem 2. Given any RCPI hP,R,X.u w A.ri, let P0 = COI(hP,R,X.u w A.ri). Then hP,R,X.u w A.riis satisfied if and only if hP0,R,X.u w A.ri is satisfied. Proof for this and subsequent reductions are given in [17].

Let p be a rational prime and let $K = \square (\zeta_p)$. We write ζ for ζ_p or this section. Recall that Khas degree $\varphi(p) = p-1$ over \square . We wish to show that $O_K = \square[\zeta]$. Note that ζ is a root of $x^p - 1$, and thus is an algebraic integer; since O_K is a ring we have that $\Box [\zeta] \subseteq O_{\kappa}$. We give a proof without assuming unique factorization of ideals. We begin with some norm and trace computations. Let j be an integer. If j is not divisible by p, then ζ^{j} is a primitive p^{th} root of unity, and thus its conjugates are $\zeta, \zeta^2, ..., \zeta^{p-1}$. Therefore

$$Tr_{K/\Box}(\zeta^{j}) = \zeta + \zeta^{2} + ... + \zeta^{p-1} = \Phi_{p}(\zeta) - 1 = -1$$

If p does divide j, then $\zeta^{j} = 1$, so it has only the one conjugate 1, and $Tr_{K/\square}(\zeta^j) = p-1$ By linearity of the trace, we find that

$$Tr_{K/\square} (1-\zeta) = Tr_{K/\square} (1-\zeta^2) = \dots$$

$$=Tr_{K/\square}(1-\zeta^{p-1})=p$$

We also need to compute the norm of $1-\zeta$. For this, we use the factorization

$$x^{p-1} + x^{p-2} + \dots + 1 = \Phi_p(x)$$

$$=(x-\zeta)(x-\zeta^2)...(x-\zeta^{p-1});$$

Plugging in x = 1 shows that

$$p = (1 - \zeta)(1 - \zeta^{2})...(1 - \zeta^{p-1})$$

Since the $(1-\zeta^j)$ are the conjugates of $(1-\zeta)$, this shows that $N_{K/\square}(1-\zeta)=p$ The key result for determining the ring of integers O_{K} is the following.

LEMMA 1.9

$$(1-\zeta)O_{\kappa}\cap\Box=p\Box$$



Proof. We saw above that p is a multiple of $(1-\zeta)$ in O_K , so the inclusion $(1-\zeta)O_K \cap \square \supseteq p\square$ is immediate. Suppose now that the inclusion is strict. Since $(1-\zeta)O_K \cap \square$ is an ideal of \square containing $p\square$ and $p\square$ is a maximal ideal of \square , we must have $(1-\zeta)O_K \cap \square = \square$ Thus we can write $1=\alpha(1-\zeta)$

For some $\alpha \in O_K$. That is, $1-\zeta$ is a unit in O_K .

COROLLARY 1.1 For any $\alpha \in O_K$, $Tr_{K/\square} ((1-\zeta)\alpha) \in p\square$ PROOF. We have

$$Tr_{K/\square} ((1-\zeta)\alpha) = \sigma_{1}((1-\zeta)\alpha) + ... + \sigma_{p-1}((1-\zeta)\alpha)$$

$$= \sigma_{1}(1-\zeta)\sigma_{1}(\alpha) + ... + \sigma_{p-1}(1-\zeta)\sigma_{p-1}(\alpha)$$

$$= (1-\zeta)\sigma_{1}(\alpha) + ... + (1-\zeta^{p-1})\sigma_{p-1}(\alpha)$$

Where the σ_i are the complex embeddings of K (which we are really viewing as automorphisms of K) with the usual ordering. Furthermore, $1-\zeta^j$ is a multiple of $1-\zeta$ in O_K for every $j\neq 0$. Thus

 $Tr_{K/\square}(\alpha(1-\zeta)) \in (1-\zeta)O_K$ Since the trace is also a rational integer.

PROPOSITION 1.4 Let p be a prime number and let $K = |\Box(\zeta_p)|$ be the p^{th} cyclotomic field. Then

$$O_K = \square[\zeta_p] \cong \square[x]/(\Phi_p(x));$$
 Thus $1, \zeta_p, ..., \zeta_p^{p-2}$ is an integral basis for O_K .

PROOF. Let $\alpha \in O_K$ and write

$$\begin{split} \alpha &= a_0 + a_1 \zeta + \ldots + a_{p-2} \zeta^{p-2} & \text{ With } \quad a_i \in \Box \ . \\ \alpha &(1-\zeta) = a_0 (1-\zeta) + a_1 (\zeta - \zeta^2) + \ldots \\ &+ a_{p-2} (\zeta^{p-2} - \zeta^{p-1}) \end{split}$$

By the linearity of the trace and our above calculations we find that $Tr_{K/\square}\left(\alpha(1-\zeta)\right)=pa_0$ We also have

 $Tr_{K/\square}\left(\alpha(1-\zeta)\right)\in p\square$, so $a_0\in\square$ Next consider the algebraic integer

$$(\alpha - a_0)\zeta^{-1} = a_1 + a_2\zeta + \dots + a_{p-2}\zeta^{p-3}$$
; This is an algebraic integer since $\zeta^{-1} = \zeta^{p-1}$ is. The same

argument as above shows that $a_1 \in \square$, and continuing in this way we find that all of the a_i are in \square . This completes the proof.

Let $K = \square$, then the local ring $\square_{(p)}$ is simply the subring of \square of rational numbers with denominator relatively prime to p. Note that this ring $\square_{(p)}$ is not the ring \square_p of padic integers; to get \square none must complete \square (n). The usefulness of $O_{K,p}$ comes from the fact that it has a particularly simple ideal structure. Let a be any proper ideal of $O_{K,p}$ and consider the ideal $a \cap O_{\kappa}$ of O_{κ} . We $a = (a \cap O_K)O_{K,n}$; That is, that a is generated by the elements of a in $a \cap O_K$. It is clear from the definition of an ideal that $a \supseteq (a \cap O_K)O_{K,n}$. To prove the other inclusion, let α be any element of a. Then we can write $\alpha = \beta / \gamma$ where $\beta \in O_K$ and $\gamma \notin p$. In particular, $\beta \in a$ (since $\beta / \gamma \in a$ and a is an ideal), so $\beta \in O_K$ and $\gamma \notin p$, so $\beta \in a \cap O_K$. Since $1/\gamma \in O_{K_n}$, this implies that $\alpha = \beta / \gamma \in (a \cap O_K)O_{K,p}$, claimed.We can use this fact to determine all of the ideals of $O_{K,p}$. Let a be any ideal of $O_{K,p}$ and consider the ideal factorization of $a \cap O_K$ in O_K . write it as $a \cap O_K = p^n b$ For some n and some ideal b, relatively prime to p, we claim first that $bO_{K,p} = O_{K,p}$. We now find that

$$a = (a \cap O_K)O_{K,p} = p^n bO_{K,p} = p^n O_{K,p}$$

Since $bO_{K,p}$. Thus every ideal of $O_{K,p}$ has the form $p^nO_{K,p}$ for some n; it follows immediately that $O_{K,p}$ is noetherian. It is also now clear that $p^nO_{K,p}$ is the unique non-zero prime ideal in $O_{K,p}$. Furthermore, the inclusion $O_K \mapsto O_{K,p} / pO_{K,p}$ Since $pO_{K,p} \cap O_K = p$, this map is also surjection, since the residue class of $\alpha / \beta \in O_{K,p}$ (with $\alpha \in O_K$ and $\beta \notin p$) is the image of $\alpha\beta^{-1}$ in $O_{K/p}$, which makes sense since β is invertible



in $O_{K/p}$. Thus the map is an isomorphism. In particular, it is now abundantly clear that every non-zero prime ideal of $O_{K,p}$ is maximal.

show that $O_{K,p}$ is a Dedekind domain, it remains to show that it is integrally closed in K. So let $\gamma \in K$ be a root of a polynomial with coefficients in $O_{K,p}$; write this polynomial as $x^m + \frac{\alpha_{m-1}}{\beta_{m-1}} x^{m-1} + \ldots + \frac{\alpha_0}{\beta_0}$ With $\alpha_i \in O_K$ and $\beta_i \in O_{K-p}$. Set $\beta = \beta_0 \beta_1 \ldots \beta_{m-1}$. Multiplying by β^m we find that $\beta \gamma$ is the root of a monic polynomial with coefficients in O_K . Thus $\beta \gamma \in O_K$; since $\beta \notin p$, we have $\beta \gamma / \beta = \gamma \in O_{K,p}$. Thus $O_{K,p}$ is integrally close in K.

COROLLARY 1.2. Let K be a number field of degree n and let α be in O_K then $N_{K/\square}$ $(\alpha O_K) = \left| N_{K/\square} (\alpha) \right|$

PROOF. We assume a bit more Galois theory than usual for this proof. Assume first that K/\square is Galois. Let σ be an element of $Gal(K/\square)$. It is clear that $\sigma(O_K)/\sigma(\alpha)\cong O_{K/\alpha}$; since $\sigma(O_K)=O_K$, this shows that $N_{K/\square}^{'}(\sigma(\alpha)O_K)=N_{K/\square}^{'}(\alpha O_K)$. Taking the product over all $\sigma\in Gal(K/\square)$, we have $N_{K/\square}^{'}(N_{K/\square}(\alpha)O_K)=N_{K/\square}^{'}(\alpha O_K)^n$ Since $N_{K/\square}(\alpha)$ is a rational integer and O_K is a free \square -module of rank n.

 $O_{\scriptscriptstyle{K}}/N_{\scriptscriptstyle{K/\square}}(\alpha)O_{\scriptscriptstyle{K}}$ Will have order $N_{\scriptscriptstyle{K/\square}}(\alpha)^n$; therefore

$$N_{K/\square}^{'}(N_{K/\square}(\alpha)O_{K}) = N_{K/\square}(\alpha O_{K})^{n}$$

This completes the proof. In the general case, let L be the Galois closure of K and set [L:K] = m.

E. Encryption Keys

We use symmetric keys to encrypt and decrypt attribute values. These keys are distributed only to the brokers that are trusted with the attribute values. The system will never deliver these keys to clients. This reduces the number of nodes that are trusted with sensitive keys, and that take part in key management protocols. Note that this does not affect security since local brokers encrypt and decrypt attribute values on behalf of connected clients, and

deliver events to clients over secure links. To support cryptographic properties such as key freshness, and forward and backward secrecy [22], the system requires key management service(s). The most suitable key management strategy depends on the broker-network architecture. For EDSAC21 we assume a stable configuration with static, multi-hop, interbroker connections and are investigating a treebased approach [22]. However, the dynamic nature of a peer-to-peer routing layer presents special problems, and we are also evaluating an alternative, ad-hoc network based approach [23]. Efficient group key management [24] is not the focus of this paper. Overall, the efficiency of key distribution will have little impact on performance, since symmetric keys are distributed only to brokers, as opposed to publishers and subscribers. Relatively few entities are involved in key dissemination, and changes will be infrequent. However, correct key management is essential for the security of the system.

V. CASE STUDY

We now illustrate our architecture for a city in which the publish/subscribe systems of different emergency services interoperate securely and efficiently. We use a break-in to a university building as an example. Fig. 4 shows the principals, brokers and messages discussed below. We assume that equipment failure has left the police network partititioned, and that broker b1 is connected only through the fire network. 1) We focus on two police officers on night shift; part of their duty is to respond to notifications of burglaries. We assume that the event-type BurglaryEvent is already advertised when the officers come on duty. This means that a rendezvous node b5 is assigned for the type and subscriptions can be made. We shall see further advertisements, and subsequent publications, can be made as burglaries are detected in different areas. We assume that both officers authenticate with their local OASIS service on coming on duty and, assuming that their credentials are valid, acquire the role with associated privilege to send subscription messages: s1 and s2 respectively.

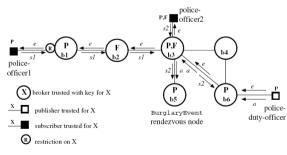


Fig. 4. Notifying two police officers of a BurglaryEvent

Officer 1 is a probationary officer, who moves between different parts of the city. Officer 2 is located in West Cambridge. Suppose that at the start



2 of her shift officer subscribes BurglaryEvent(location = 'West Cambridge'). Since this subscription requires filtering on the location attribute, and this attribute is encrypted with the police key (recall the event type hierarchy shown in Fig. 2), the officer knows that her local broker must be trusted with the police key, i.e. a P broker. Officer 1 tries to subscribe to all burglary events police with code less than BurglaryEvent(polCode < 4), but the request is only partially granted. Instead, the subscription is restricted, as described in Section 4, to deliver only those events that occur in the officer's current location. This restriction, which is based on a dynamically checked environmental constraint, is shown in Fig. 4, attached to his broker connection.

- 2) Any broker through which s1 and/or s2 travel (towards their rendezvous node and then along the reverse path of advertisements) will update its internal routing state appropriately. Note that our security architecture augments standard Hermes subscription setup behaviour when we reach broker b2. Whilst s1 travels through this broker, the broker is not part of the police network, and thus will not have access to the police key. Therefore this broker will be forced to degrade routing efficiency by ignoring police officer 1's filter on the polCode attribute, which it cannot decrypt, and routing all events forward.
- 3) We show a duty-officer at a police station who must notify police officers of reported burglaries. Like officers 1 and 2, the duty-officer authenticates himself with his local OASIS service, and acquires privileges to advertise BurglaryEvents. Again, his local broker needs access to the police key. The consequent advertisement message is shown as a in Fig. 4. This step could occur in parallel with a subscription, see Step 1. If a broker notices that an existing subscription matches a new advertisement, it will resend the subscription message along the reverse path of the new advertisement towards the publisher. All this occurs at the start of the officers' sessions, a long time (in publish/ subscribe terms) before the actual burglary occurs.

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