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Automatic Malware Signature Generation

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Thanks...

Summary

In most recent years the proliferation of malicious software, namely Malware, has had a massive increase: according to AV Atlas Dashboard [1] the new malware samples (and PUA - Potentially Unwanted Application) currently detected every day are about 440.000 (at the time of writing), and this number is predicted to only keep growing. The total number of known Microsoft Windows malicious software (and PUA) passed from about 55 million in 2011, to about 400 million in 2016, and finally to nearly 830 million now. The huge number of malware samples out there in the wild renders the detection through manually generated signatures (patterns that identify malicious code) infeasible and consequently imposes the urgent need of tools able to automatically detect malware and possibly describe it in an human-interpretable way. Several methodologies have been proposed through the years, ranging from signature-based detection (especially with Yara Rules) to various Machine Learning approaches like Decision Trees, Naive Bayes models and Neural Networks.

This thesis presents a novel model built upon previous works in **ML-based** (Machine Learning) automatic PE (Microsoft Windows Portable Executable) malware detection and description and introduces a new evaluation procedure on the learnt implicit representation/signature of malware samples that may prove the applicability of its usage in the Malicious family prediction and ranking tasks. The model is trained on an open source large scale dataset of malware and benignware samples with the aim of creating high quality implicit signatures capable of correctly detecting (and describing) unseen malware samples as well as obfuscated malware and new variants, with high True Positive Rate (TPR) and high Recall at low False Positive Rates (FPRs).

The **Proposed Model**'s results in the different tasks (both the ones it was trained on - Malicious/Benign label and descriptive tags prediction - and the additional malware family prediction and ranking tasks) were compared to the previous models' ones. In particular, the **ALOHA** model proposed in [2] and the **Joint Embedding** model described in [3] were selected as reference models.

The results show that the **Proposed Model** generates implicit signatures (samples embeddings) which provide comparable TPRs, Accuracies, Recalls, Precisions and F1 Scores at low false positive rates with respect to the ones produced by the mentioned previous methods. When tested on the Malware Family prediction and ranking tasks, however, the results were less promising. After more detailed analyses, it was apparent that malicious samples belonging to different families but with similar characteristics and behaviour were mapped close to each other in the model's learnt embedding space, resulting in the same family label being predicted for all those samples. This implies that unfortunately, for now, the proposed model lacks the capability to generalize over the more interesting task of malware family categorization.

Nevertheless, future works capable of overcoming this limitation may be very useful to the IT-Security field in the current scenario and could even enable the generation of explicit (and thus more interpretable) signatures derived from the learnt implicit ones.

Acknowledgements

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Chapter 1

Introduction

Malware first made its appearance in the 1960s. Then, hackers used to design computer viruses mainly for fun, as an exciting prank/experiment; their creations would generally display harmless messages and then spread to other computers [4]. There are numerous examples of malware created at that time within a laboratory setting: for example the **Darwing game** in 1962 and **Creeper** in 1971.

In the early 1980s, the concept of malware caught on in the technology industry, and numerous examples of viruses and worms appeared both on Apple and IBM personal computers. With the introduction of the World Wide Web and the commercial internet in the 1990s it eventually became widely popularized, so much that Yisreal Radai coined the term **malware** in 1990.

```
Elk Cloner: The program with a personality

It will get on all your disks
It will infiltrate your chips
Yes it's Cloner!

It will stick to you like glue
It will modify ram too
Send in the Cloner!
```

Figure 1.1: Poem displayed by Elk Cloner virus

The previously mentioned 1960s and 1970s malware were all kept within a laboratory environment and never managed to escape to the wild. **Elk Cloner** (1981) was the first known virus to have been able to escape its creation environment. It spread through infected floppy disks but did not cause deliberate harm simply displaying the 'poem' shown in figure 1.1 on the user's screen. Following the success of that prank gone wild, the first Microsoft PC virus, called **Brain**, was then created in 1986. Again, like *Elk Cloner*, Brain was mostly annoying rather than harmful, but it was also the first known virus capable of concealing its presence on the disk thus evading detection. In 1988 the first worm, called **Morris** worm, an experimental, self-propagating, self-replicating program was released on the internet [5]. In 1988 made its appearance also the first example of intentionally harmful virus, the **Vienna** virus, which encrypted data and destroyed files. This led to the creation of the first antivirus tool ever [4].

In the following decades malware has evolved both regarding its complexity and malware sample numbers. In particular, according to AV Atlas [1], in 2019 and 2020 (and until mid 2021 - that is until the time of writing) the number of newly generated malware blew with respect to previous years, as shown in graph 1.2, to the point that currently approximately 5.1 new Microsoft Windows malware (and PUA) are generated per second, $\sim 18,500$ per hour. As reported by AV Atlas dashboard and as shown in graph 1.3, in recent years the total amount of unique malware (and PUA) variants reached impressive numbers, to the point that they are now more than

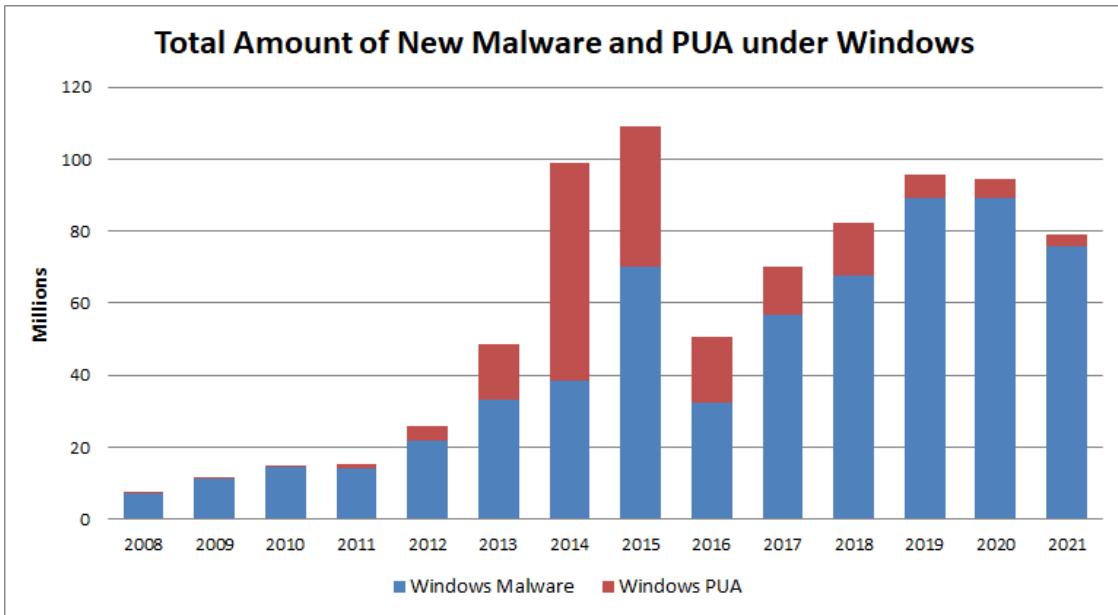


Figure 1.2: AV Atlas - Total Amount of New Malware and PUA

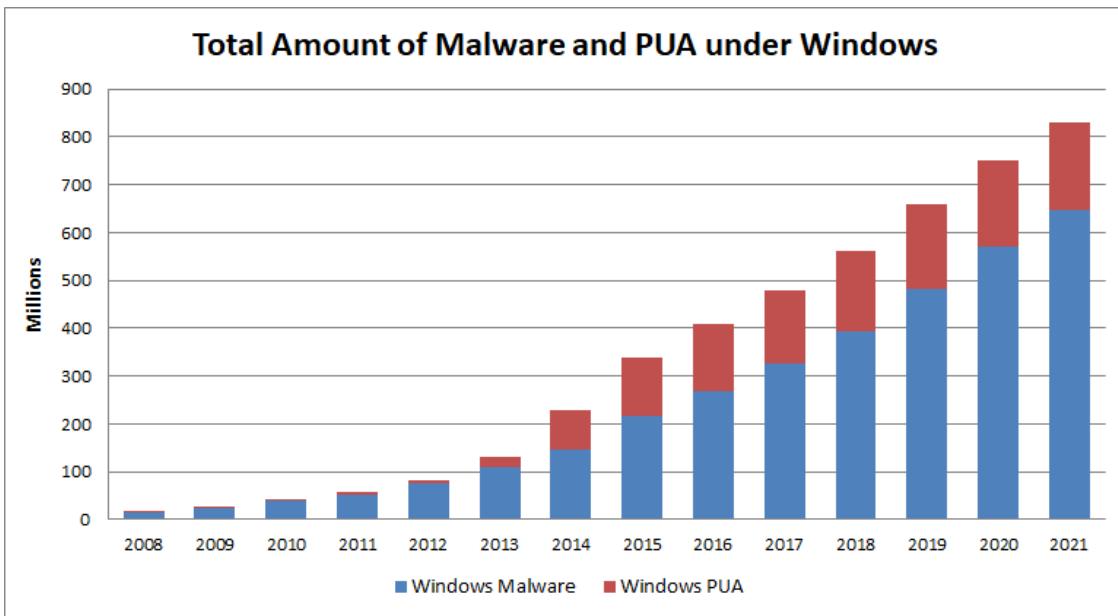


Figure 1.3: AV Atlas - Total Amount of Malware and PUA

830 millions. Moreover, nowadays malware commonly use obfuscation and other sophisticated techniques, such as Polymorphism and Metamorphism, to evolve their structure thus evading detection.

For all these reasons, signature-based detection techniques, such as manually generated Yara Rules, which are typically used by most commercial anti-virus solutions, are becoming inefficient in the present scenario. In fact, it is now straight up impossible for analysts to manually analyse each malware variant that is found in the wild. Moreover, even when a new malware family is identified and a appropriate amount of its samples are analysed, the generated signature may not be capable of detecting new variations or may even be rendered useless through the use of obfuscation and/or polymorphic mechanisms. There is therefore the need for automated malware analysis solutions capable of automatically generating (implicit or explicit) signatures effective at distinguishing malicious from benign code while being less susceptible to code modifications and

obfuscation attempts.

This thesis presents a research aimed at satisfying this need for automated malware detection solutions. In particular, it presents a novel model built upon previous works on ML-based (Machine Learning) automatic malware detection and description designed for PE (Microsoft Windows Portable Executable) files. Moreover, it introduces a new evaluation procedure that may prove the applicability of the model learnt implicit representation/signature of malware samples in the Malware family prediction and ranking tasks. These tasks are particularly interesting for malware analysts since they enable them to quickly categorize malware samples as being part of specific sets (families) of samples with common behavioural and structural characteristics.

The proposed framework life cycle can be conceptually divided in three phases: **model architecture definition**, **model training and evaluation** and finally **model deployment**. In particular, in the first phase the FNN (Feedforward Neural Network) model architecture is defined and implemented taking inspiration from previous works such as the **ALOHA** and the **Joint Embedding** models presented in papers [2] and [3], respectively. In the second phase, instead, the proposed model is trained on an open source large scale dataset of malware and benignware samples (Sorel20M [6]) with the aim of creating high quality implicit signatures capable of detecting (and describing) unseen malware samples, as well as obfuscated malware and new variants, with high True Positive Rate (TPR) and high Recall at low False Positive Rates (FPRs). Moreover, in this phase the model is also tested on the Malware family prediction and ranking tasks using a novel dataset, containing 1000 samples belonging to 10 of the most widespread malware families in Italy at the time of writing, specifically created for that purpose. In both datasets the samples are directly represented by their numerical feature representation extracted statically from specific fields of each sample's Windows Portable Executable (PE) file header. The proposed model thus relies exclusively on static analysis features which are generally simpler, less computationally intensive and thus faster than dynamic analysis ones (executable's behaviour characteristics). Finally, in the last phase the final proposed model architecture is deployed in the wild. In particular, it can be used as an automatic Malware detection tool that provides additional description tags useful for remediation. Moreover, potentially, if the corresponding evaluation results allow it, it could also be used to provide information about the malware family each analysed sample most probably belongs to, among the set of families of interest.

This thesis focuses on the first two phases previously mentioned. In particular, it concentrates on defining, training and evaluating the best model architecture possible for the tasks at hand. However, some code optimization challenges resulting from the slowness of the code in the instance used for the experiments meant that it could be possible to train the model only with the first half of the samples of the Sorel20M dataset in a reasonable time, with some approximations on the samples dispersion when random sampling them from the dataset. This led to slightly worse performance than what is theoretically possible with the current architecture using the whole dataset. Nevertheless, the deployment of the proposed model on a real-world scenario is theoretically possible with the current final architecture, although it would be better to train the proposed model on the whole Sorel20M dataset on a better instance first in order to see its true potential.

Chapter 2

Background

2.1 Malware

Malware, short for *malicious software*, is a general term for all types of programs designed to perform harmful or undesirable actions on a system. In fact in the context of IT security the term *malicious software* commonly means [7]:

Software which is used with the aim of attempting to breach a computer system's security policy with respect to Confidentiality, Integrity and/or Availability.

Malware consists of programming artefacts (code, scripts, active content, and other software) designed to disrupt or deny operation, gain unauthorized access to system resources, gather information that leads to loss of privacy or exploitation, and other abusive behaviour. Malware is not (and should not be confused with) defective software - software that has a legitimate purpose but contains harmful bugs (programming errors).

Different companies, organizations and people describe malware in various ways. For example **Microsoft** defines it in a generic way:

Malware is a catch-all term to refer to any software designed to cause damage to a single computer, server, or computer network [8].

The **National Institute of Standards and Technology (NIST)**, on the other hand, presents multiple definitions for malware, describing it as "hardware, firmware, or software that is intentionally included or inserted in a system for a harmful purpose" [9].

In another more specific definition **NIST** affirms that Malware is:

A program that is inserted into a system, usually covertly, with the intent of compromising the confidentiality, integrity, or availability of the victim's data, applications, or operating system or of otherwise annoying or disrupting the victim [9].

The computer system whose security policy is attempted to be breached is usually known as the **target** for the malware. The cybercriminal who originally launched the malware with the purpose of attacking one or more targets, on the other hand, is generally referred to as the "initiator of the malware". Furthermore, depending on the malware type, the initiator may or may not exactly know what the set of targets is [7].

According to the above definitions software is defined as malicious in relation to an attempted breach of the target's **security policy**. In other words, software is often identified as malware based on its *intended use*, rather than the particular technique or technology used to build it.

2.1.1 Why is Malware used

Generally, cybercriminals use malware to access targets' sensitive data, extort ransoms, or simply cause as much damage as possible to the affected systems.

More generally malware serves a variety of purposes. For example, the most common cybercriminals' uses of malware are: [10]

- **To profit financially (either directly or through the sale of their products or services).** For example, attackers may use malware to infect targets' devices with the purpose of stealing their credit account information or cryptocurrency. Alternatively, they may sell their malware to other cybercriminals or as a service offering (*malware-as-a-service*).
- **As a means of revenge or to carry out a personal agenda.** For example, Brian Krebs of Krebs on Security was struck by a big DDoS attack in 2016 after having talked about a DDoS attacker on his blog.
- **To carry out a political or social agenda.** For example, there exist many Nation-state actors (such as state-run hacker groups in China and North Korea) and hacker groups such as Anonymous.
- **As a way to entertain themselves.** Some cybercriminals perpetrate attacks on victims just for fun.

Obviously there are also reasons for **non-malicious** actors to create and/or deploy some types of malware too - for example they can be used to test a system's security, to spy on someone as part of a legal act/police operation, etc.

2.1.2 Common Malware types

There are different ways of categorizing malware; one way is by *how* the malicious software spreads. Another one is by *what it does* once it has successfully infected its victim's computers (i.e. what is its payload, how it exploits the system or makes it vulnerable).

By how they spread

Terms like *trojan*, *virus* and *worm* are commonly used interchangeably to indicate generic malware, but they actually describe three subtly different ways malicious software can infect target computers [11]:

- **Trojan horse.** Generally speaking, a *Trojan Horse*, commonly referred to as a "Trojan", is any program that disguises itself as legitimate and invites the user to download and run it, concealing a malicious payload. When executed, the payload - malicious routines - may immediately take effect and cause many undesirable effects, such as deleting the user files or installing additional malware or PUAs (Potentially Unwanted Apps), depending on the payload attached to them [5].

Trojans may hide in games, apps, or even software patches, or they may rely on social engineering and be embedded in attachments included in phishing emails. However, they cannot self-replicate but rely on the system operators to activate.

- **Virus.** The term "computer virus" is used for describing a **passive** self-replicating malicious program. Usually spread via infected websites, file sharing, or email attachment downloads, it lies dormant until the infected host file or program is activated. At that point it spreads to other executables (and/or boot sectors) by embedding copies of itself into those files. A virus, in fact, in order to spread from one computer to another, usually relies on the infected files possibly ending up, by some means or another, in the target system. The mean of transport (file, media file, network file, etc.) is often referred to as the virus

vector. Depending on how complex the virus code is, it may be able to modify its copies upon replication. For the transport of the infected files to the target system(s), the virus may rely on an unsuspecting human user (who for example uses a USB drive containing the infected file) or may initiate the transfer itself (for example, it may send the infected files as an e-mail attachment) [7].

Viruses may also perform other harmful actions other than just replicating, such as creating a backdoor for later use, damaging files, stealing information, creating botnets, render advertisements or even damaging equipment.

- **Worm.** A worm is a self-replicating, **active** malicious program that exploits various system vulnerabilities to spread over the network. Particularly, it relies on vulnerabilities present in the target's operating system or installed software. Worms usually consume a lot of bandwidth and processing resources due to continuous scanning and may render the host unstable, sometimes even causing crashes. Apart from replicating, computer worms also contain "payloads" which are pieces of code written to perform various nefarious actions on the affected hosts, for example stealing data, deleting files or creating bots - which can lead the infected systems to become part of a botnet [5].

Moreover, attackers can also install malware "manually" on a computer, either by gaining physical access to the target system or by using privilege escalation methods to obtain remote administrator access [12].

By what they do

There is a wide range of potential attack techniques used by malware, here are briefly described some of them:

- **Adware.** *Adware* is any software package which automatically plays, displays, or downloads advertisements to a computer. These can be in the form of pop-up ads, ad banners in websites, or advertisements displayed by software, that lure the user into making a purchase. The goal of Adware is, in fact, to generate revenue for its author.

Often times adware comes even bundled with "free", or discounted, versions of non-malicious software and/or applications since it is usually seen by developers as a way to recover development costs [13].

Adware, by itself, is annoying but somewhat harmless, since it is solely designed to deliver ads; however, it often comes bundled with spyware (such as keyloggers), and/or other privacy-invasive software that is capable of tracking user activity and steal information [14].

- **Backdoor.** A *backdoor*, also called Remote Access Trojan (RAT), is a vulnerability deliberately buried into software's code allowing to bypass typical protection mechanisms, like credentials-based login authentication. Once a system has been compromised (by others types of malware or other methods), one or more backdoors may be installed. This is done with the purpose of allowing the attacker easier access in the future without alerting the user or the system's security programs. Moreover, backdoors may also be installed before other malicious software, to allow attackers entry [13].

On the other hand, it is not uncommon to see many (non-malicious) device or software manufacturers ship their products with intentionally hidden backdoors to allow company personnel or law enforcement access to the system when needed [15]. Alternatively, backdoors are sometimes hidden in programs also by intelligence services.

- **Browser Hijacker.** A *Browser Hijacker*, also called "hijackware", is a type of malicious program which considerably modifies the behaviour of the victim's web browser. It can be used to make money off unwanted ads, to steal information from users, or to infect the systems with other malware by redirecting users to malicious websites [15].

- **Bots / Botnet.** In general, *bots* (short for 'robots') are software programs designed to automatically perform specific operations. Some bots are used for legitimate and harmless purposes such as video programming, video gaming, internet auctions and online contest, among other functions. It is however becoming increasingly common to see bots being used maliciously. Malicious bots can be (and usually are) used to form botnets. A botnet is defined as a network of host computers (zombies/bots) that is controlled by an attacker - the *bot-master* [5]. Botnets are frequently used for DDoS (Distributed Denial of Service) attacks, but there are other ways that botnets can be useful to cybercriminals: for example they can be employed to carry out brute force attacks on websites or to distribute malware [10].
- **Crypto-miner.** Crypto-miners are a relatively new malware family. Cybercriminals employ this type of malicious tools to mine Bitcoin and/or other similar digital currencies on the target machine by exploiting the victim system's computing power, without the owner realising it. Obviously, the mined coins end up in the attackers' digital crypto wallets. Recently, a more modern method of crypto-mining that works within browsers (also called crypto-jacking), has become quite popular [16].
- **File-less malware.** As the term suggests, File-less malware is a type of malware that operates from a victim's computer memory, not from files on the hard drive, taking advantage of legitimate code and tools (known as "LOLBins" [10]) already existing within the system. File-less malware leaves no malware files to scan and no malicious processes to detect, and is therefore harder to detect and remove than traditional malware: it is, in fact, up to ten times more successful [17]. Furthermore, it also renders forensic analysis more difficult because it disappears from the victim's computer upon rebooting.
- **Keylogger.** Keyloggers, often considered as being a sub-category of spyware, are malicious programs which secretly track keystrokes on a keyboard, without the system's owner consent - action that is usually referred to as *keylogging* or *Keystroke logging*. The collected data is stored and sent to the attacker who uses it to figure out highly sensitive information such as passwords, usernames and payment details. Keylogging can be performed in various ways, ranging from hardware and software-based approaches to the more sophisticated electromagnetic and acoustic analysis [13].
- **RAM Scraper.** *RAM scraper* malware, also known as *Point-of-Sale (POS)* malware, targets POS systems like cash registers or vendor portals, harvesting data temporarily stored in RAM (Random Access Memory). Doing so the attacker can easily access unencrypted credit card numbers [15].
- **Ransomware.** *Ransomware*, also known as "encryption" or "crypto" Trojan, is a malicious program that, after having infected a host or network, holds the system captive and displays a message requesting a ransom to the host/network users. In particular, it encrypts data on the infected system (or anyway it locks down the system preventing user access) and only unblocks it when the correct password - decryption key - is entered. Without the latter, is practically impossible to regain access to the system. Digital currencies such as Bitcoin and Ether are the most common means of payment, making it difficult to track cybercriminals. Moreover, paying the ransom does not guarantee the user to receive the necessary decryption key. Additionally, some forms of ransomware threaten victims to publicize sensitive information within the encrypted data in order to convince them to pay the ransom.
- **Rogue Security Software.** *Rogue Security Software* can be considered as a form of scareware. This type of malware programs presents itself as a security tool to remove risks from the user's system. In reality, this fake security software installs more malware onto the system [15].
- **Rootkit.** A *rootkit* is generally thought as a type of malicious software, or a collection of software tools, designed to remotely access or control a computer without being detected by users or security programs. An attacker who has installed a rootkit on a system is able to perform a wide range of malicious activities such as remotely executing files, logging user

activities, installing hidden malware, etc. Since a rootkit operates stealthily and continually hides its presence, its prevention, detection and removal can be difficult and often relies on manual methods [14].

More recently, the term "rootkit" has often been used to refer also to concealment routines in a malicious program. These highly advanced and complex routines are written to hide malware within legitimate processes on the infected computer. In fact, once a malicious program has been installed on a system, it is essential that it remains hidden, to avoid detection and disinfection [13].

- **Scareware.** Scareware is a generic term for malware that uses social engineering to frighten and manipulate users, inducing them into thinking their system is vulnerable or has been compromised. However, in reality no danger has actually been detected: it is a scam. The attack succeeds when the user purchases unwanted - and potentially dangerous - software in an attempt to eliminate the "threat". Generally, the suggested software is additional malware or allegedly protective software with no value whatsoever [16].

Some versions of scareware act as a sort of shadow version of ransomware; they claim to have taken control of the victim's system and demand a ransom. However they are actually just trying to fool the victim [12].

- **Spyware.** *Spyware*, another name for *privacy-invasive software*, is a type of malicious software that spies on the infected system's user activity. Specifically it can collect various types of personal information about users, such as Internet browsing habits, credit card details and passwords, without their knowledge. The information gathered is then sent back to the responsible cybercriminal(s).

However, the functions of spyware often go far beyond simple activity monitoring and information gathering. In fact, they may also interfere with the user's control of the computer in other ways, such as installing additional software, redirecting web browser activity and changing computer settings [13].

Law enforcement, government agencies and information security organizations often use spyware to monitor communications in a sensitive environment or during an investigation [18].

Other cyber-threats

Other cyber threats which are not strictly, and should not be confused with, malware are, for example:

- **Software Bug.** A software bug is an error, or flaw, in a computer program code or system that causes it to produce an incorrect or unexpected result, or to behave in unintended ways. Usually, most of these defects arise from human errors made in the program's source code. Minor bugs only slightly affect the behaviour of a program and can therefore remain undiscovered for quite a long time. On the other hand, more significant bugs can cause crashes or freezes [14].
- **Software Vulnerability.** In computer security, a vulnerability is a hole or a weakness in an application, which can be a design flaw or an implementation bug, that can be exploited by a threat actor, such as an attacker to bypass access controls such as user authentication, override access privileges, steal data or perform other unauthorized actions within a computer system potentially causing harm to the stakeholders of an application. Stakeholders include the application owner, application users, and other entities that rely on the application [19].
- **Malvertising.** Malvertising is the use of legitimate ads or ad networks to covertly deliver malware to unsuspecting users' computers. For example, a cybercriminal might pay to place a malicious ad on a legitimate website. When clicked, it either redirects the victims to a malicious website or installs malware on their computer. In some cases, the code embedded in an ad might even execute automatically without any action from the user, a technique referred to as a "drive-by-download".

- **Phishing.** Phishing is a type of social engineering attack in which the attacker attempts, through email messages (or other means), to trick users into divulging passwords or other personal and financial information, downloading a malicious attachment or visiting a website that installs malware on their systems.

Some phishing emails are highly sophisticated and can deceive even experienced users, especially if the attacker has successfully compromised a known contact's email account and uses it to spread phishing attacks or malware such as worms. Others are less sophisticated and simply spam as many emails as possible with messages such as "Check your bank account details" [20].

- **Spam.** In cybersecurity, unsolicited emails are generally referred to as *spam*. Typically, spam includes emails carrying unwanted advertisements, fraud attempts, links to malicious websites or malicious attachments. Most spam emails contain one or more of the following characteristic traits [15]:

- Poor spelling and grammar
- Unusual sender address
- Unrealistic claims
- Suspicious links

General considerations on malware types

Malware samples are usually categorised both by their means of infection and their behavioural category: for instance, *WannaCry* is a ransomware worm. Moreover, a particular piece of malware may have various forms with different attack vectors: e.g., the banking malware called *Emotet* has been spotted in the wild as both a trojan and a worm [12]. Finally, many instances of malware fit into multiple categories: for example *Stuxnet* is both a worm, a virus and a rootkit.

2.2 Detection evasion

From the creation of the first malwares in 1960s, when hackers used to design computer viruses mainly for fun, a strong competition between attackers and defenders has risen. To defend from malware attacks, anti-malware groups have been developing increasingly complex (and clever) new techniques. On the other hand, malware developers have conceived and adopted new tactics/methods to avoid the malware detectors [21].

The first type of anti-malware tools were mostly based on the assumption that malware structures do not change appreciably during time. In fact, initially, the malware machine code was completely unprotected. This allowed analysts to exploit opcode sequences to recognise specific malware families. Recently, however, a big advancement led to the so-called "second generation" malware [22] which, to evade such opcode signatures, employs several obfuscation techniques and can create variants of itself. This posed a challenge to anti-malware developers.

The first time a malware has been recognised to exhibit detection avoidance behaviour was in 1986 with the *Brain* virus [23]. In fact, such malware managed to conceal the infected disk section whenever the user attempted to read it, forcing the computer to display clean data instead of the infected part. From that moment on, the ever increasing popularity of detection evasion techniques among malware writers has shown that malware survival has become the number one priority: the longer the malware remains undetected, the more harm it can do and the more profitable it is to its writer [5].

2.2.1 Reverse-Engineering

Reverse engineering, in broad terms, indicates the process of extracting knowledge, ideas, design philosophy etc. from anything man-made [24].

Software reverse engineering is, of course, the application of reversing methodologies and techniques to extract knowledge from a software product to better understand its inner workings.

Reversing is used extensively by both malicious actors and investigators but with opposing purposes. Malware developers often use it to discover vulnerabilities in systems or programs, while analysts and antivirus software developers use it to analyse malicious programs to understand how they work, what damages they can cause, how they infect the system and reproduce, how they can be removed, detected and avoided.

2.2.2 Malware analysis

Malware analysis is the process of extracting as much information as possible from malicious samples discovered in the wild, which usually are in the form of machine code executables (compiled executables), in order to determine their purpose and functionality (and threats associated). This process allows security teams to develop effective detection techniques against the analysed malicious code, contain its damage, reverse its effects on the system, develop removal tools that can delete it from infected machines (to cleanly remove a piece of malware from an infected machine it is usually not enough to delete the binary itself) and design methods to guard systems against future infections [25].

Initially malware analysts/researchers had to manually analyse each malware sample. This process is however complex, requires high expertise, and is time-intensive. Moreover, the number of malware samples that need to be analysed on a daily basis is nowadays of the order of hundreds. This implies that the analysis of malware samples can no longer be done exclusively manually. Several analysis tools have been developed in recent years to facilitate analysts in analysing malware samples.

Traditionally, there are two main types of analysis: *static* and *dynamic*. Moreover, these two types can be, and frequently are, combined together (*hybrid* analysis) in various stages of malware analysis to optimize results [5].

Static analysis

Static analysis consists of examining an executable file's code without actually executing it. Static analysis techniques usually extract peculiar features from malicious samples in order to be able to recognise them and distinguish them from benign ones. The features usually extracted are, for example, string signatures, byte-sequence n-grams, library or API calls, opcode frequency distributions, peculiar attributes found in the executable header etc. However, this approach, being based on signatures/features extracted from already analysed samples, is not much effective on zero-day and evolutionary malware.

A malware analyst performing manual static analysis usually disassembles the binary first, meaning that he 'translates' the program's machine code instructions back into assembly language ones generating a more human-interpretable code listing. On the latter, control flow, data flow analysis, and many others static techniques can be employed to try understating the program functionality and inner workings, among other useful information [25].

Static analysis advantages are, among others, that it takes into account the entire program code and it is also usually faster (and safer) than the dynamic one. However, a general disadvantage of static analysis is that many times the information collected during this type of analysis is very simple and not always sufficient for a conclusive decision on the malicious intent of a file. It is, however, good practice to start the analysis of a suspicious executable file extracting as much information as possible through various static techniques before passing to the dynamic counterpart. The information statically extracted may in fact provide useful knowledge to better apply dynamic techniques and enhance the final results.

Additionally, another common problem to deal with when using static analysis is that, since malicious code is written directly by the adversary, it can be purposefully designed to be hard to analyse statically. For example, analysis evasion techniques like packing, encryption and obfuscation can be exploited by malware authors to hinder both disassembly and code analysis steps typical of static analysis approaches, ultimately leading to incorrect or useless information [5].

Dynamic analysis

Contrary to static analysis, *dynamic techniques* analyse the program's code while or after execution in a controlled environment. These techniques, while being non-exhaustive, they have the significant advantage that they analyse only those instructions that are actually executed by the running process. This implies that dynamic analysis is less susceptible to anti-analysis attempts like code obfuscation or anti-disassembly [25]. Moreover, dynamic analysis is also more effective in terms of malicious behaviour detection, since it doesn't look at the disassembled code but, through the use of monitoring tools, it tracks the operations that the code performs on the file system, registry, network etc. It is however, computationally more expensive and time consuming.

Basic dynamic analysis consists of observing the sample under analysis interacting with the system. For example, this can be done taking a snapshot of the original system state, introducing the malware into the system, executing it and finally comparing the new system state with the original one. The changes detected can then be used for infection removal on infected systems and/or for modelling effective signatures/features.

Advanced dynamic analysis, on the other hand, consists of directly examining the executed malware internal state while it is being run. This is done typically by monitoring the APIs and OS function calls invoked, the files created and/or deleted, the registry changes and the data processed by the program under analysis during its interaction with the system. The information extracted in this way can be used to understand the malware behaviour and functionality [5].

When using dynamic techniques, however, malware analysts don't simply run malware executables on their own computer, which most probably is even connected to Internet, as they could easily escape the analysis environment and infect other hosts/networks. It is, in fact, advised to deploy dynamic techniques on "safe" and controlled (isolated) environments such as dedicated stand-alone (and isolated) hosts, virtual machines or emulators.

The use of clean dedicated hosts, reinstalled after each dynamic analysis run, is however not the most efficient solution due to the environment re-installation process overheads. On the other hand, using virtual machines (for example VMware) to perform dynamic analysis is more efficient. In fact, in this case, since the malware only affects the virtual machine environment, it is enough, after a dynamic analysis run, to simply discard the infected hard disk image and replace it with a clean one. Unfortunately, a significant drawback is that the malware being analysed may determine it is running in a virtualized environment and, as a result, modify its behaviour. To counter this last problem one could make use of emulators, which are theoretically undetectable by analysed malware. These tools, however, run the code under analysis significantly slower and are therefore sometimes detectable using specially crafted time-related code.

Hybrid analysis

Hybrid Analysis is the combination of static and dynamic analysis. It is a technique that integrates run-time information extracted through dynamic analysis with information extracted through static analysis in order to have a complete view of the malware's behaviour while avoiding the problems posed by anti-analysis techniques as much as possible.

2.2.3 Anti-reversing

Anti-reversing techniques are techniques originally meant to make the reverse engineering process difficult for a hacker or any malicious user. The main objective of various anti-reverse engineering techniques is simply to complicate the process of reversing as much as possible. For example an attacker could use the disassembly of a binary in order to get an insight of the logic of the code as well as getting hidden information.

Recently anti-reversing techniques are, however, extensively used also by malware authors in order to make their creations difficult to analyse in an attempt to postpone detection as much as possible.

There exist several anti-reversing approaches, each with its own advantages and disadvantages. However it is common practice to use a combination of more than one of them. In the next sections some of the more common anti-reversing techniques are discussed.

2.2.4 Anti-disassembly

This part needs to be summarized since it is too long and too specific.

Anti-disassembly techniques use specially crafted code and/or data in a program to cause disassembly analysis tools to generate an incorrect program listing [26]. The attackers' usage of these techniques thus implies a time-consuming analysis for malware analysts, ultimately preventing the retrieval of the source code in a reasonable time.

Any executable code can be reverse engineered, but by armouring their code with anti-disassembly and anti-debugging techniques, attackers increase the skill level required by analysts. Furthermore, anti-disassembly techniques may also inhibit various automated analysis tools and heuristic-based engines which take advantage of disassembly analysis to identify or classify malware.

These techniques exploit the inherent weaknesses present in disassembler algorithms. Moreover, disassemblers, in order to work properly, make certain assumptions on the code being analysed. However, when these assumptions are not met, there is an opportunity for malware authors to deceive the analyst.

For example, while disassembling a program, sequences of executable code can have multiple disassembly representations, some of which may be invalid and obscure the real purpose of the program. Thus, the malware authors, in order to add anti-disassembly functionality to their creations, can produce sequences of code that deceive the disassembler into outputting a list of instructions that differs from those that would be executed [26].

There are two types of disassembler algorithms: linear and flow-oriented (recursive). The linear one is easier to implement, but it is also more simplistic and error-prone.

Linear Disassemblers

The *linear* disassembly strategy is based upon the basic assumption that the program's instructions are organized one after the other, linearly. In fact, this type of disassemblers iterates over a block of code, disassembling one instruction at a time, sequentially, without deviating. More specifically, the tool uses the size of the currently disassembled instruction to figure out what bytes to disassemble next, without accounting for control-flow instructions [26].

Linear disassemblers are easy to implement and work reasonably well when working with small sections of code. They introduce, however, occasional errors even with non-malicious binaries. The main drawback of this technique is that it blindly disassembles code until the end of the section, assuming the data is nothing but instructions packed together, without being able to distinguish between code, data and pointers.

In a PE-formatted executable file, for example, the executable code is typically contained inside a single ".text" section. However, for almost all binaries, this code section contains also data, such as pointer values. These pointers will be blindly disassembled and interpreted by the linear disassembler as instructions.

Malware authors can exploit this weakness of linear-disassembly algorithms implanting data bytes that form the opcodes of multi-byte instructions in the code section.

Flow-Oriented Disassemblers

The *flow-oriented* (or *recursive*) disassembly strategy is more advanced than the previous one and is, in fact, the one used by most commercial disassemblers like *IDA Pro* [26].

Differently from the linear strategy, the flow oriented one examines each instruction, builds a list of locations to disassemble (the ones reached by code) and keeps track of the code flow.

This implies that, if disassembling a code section we find a JMP instruction, this type of disassembler will not blindly parse the bytes immediately following the JMP instruction's ones, but it will disassemble the bytes at the jump destination address.

This behaviour is more resilient and generally provides better results, but also implies a greater complexity.

In fact, while a linear disassembler has no choices to make about which instructions to disassemble at any given time, flow-oriented disassemblers have to make choices and assumptions, in particular when dealing with conditional branches and call instructions.

Particularly, in the case of conditional branches, the disassembler needs to follow both the false branch (most flow-oriented disassemblers will process the false branch of any conditional jump first) and the true one. In typical compiler-generated code there would be no difference in output if the disassembler processes first one branch or the other. However, in handwritten assembly code and anti-disassembly code, taking first one branch or the other can often produce different disassembly for the same block of code, leading to problems in analysis.

Anti-Disassembly Techniques

Jump Instructions with the Same Target One of the most used anti-disassembly techniques consists of two consecutive conditional *jump* instructions both pointing to the same target [26].

Here is an example:

```

1  74 03  jz  loc
2  75 01  jnz loc
3
4  loc:

```

Listing 2.1: Jump Instructions with the Same Target

In this case, the conditional jump '**jz loc**' is immediately followed by a jump to the same target but with opposite condition: '**jnz loc**'. This implies that the location **loc** will always be jumped to. Consequently, the combination of **jz** with **jnz** acts, in this case, like an unconditional **jmp** instruction. A disassembler, however, since it disassembles just one instruction at a time, won't recognize this combination as being an unconditional branch. During the disassembly process, in fact, if a **jnz** instruction is encountered, the disassembler will take the false branch of the instruction and will continue disassembling, even though this branch will never be executed in practice.

Jump Instructions with a Constant Condition Another common anti-disassembly technique is composed of a single conditional *jump* instruction with an always true (or false) condition [26].

Example:

```

1  33 C0  xor eax, eax
2  74 01  jz  loc
3
4  loc:

```

Listing 2.2: Jump Instructions with a Constant Condition example

The first instruction in the example code, **xor eax, eax**, sets the **EAX** register to zero and, consequently, it sets the zero flag. The next instruction, **jz** (jump if zero flag is set), appears to be a conditional jump but in reality is not conditional at all. In fact the zero flag will always be set at this point in the program execution. The disassembler, however, will process the false branch first, even if in reality it would never trigger.

Impossible Disassembly The simple anti-disassembly techniques mentioned above are frequently coupled with the use of a, so called, *rogue byte*. A *rogue byte* is a data byte strategically placed after a conditional *jump* instruction in order to trick the disassembler. The byte inserted usually is the opcode for a multi-byte instruction, therefore disassembling it prevents the real following instruction from being properly disassembled. This byte is called *rogue byte* because it is not part of the program logic flow and it is inserted in the code with the only purpose of fooling the disassembler [26].

In all these cases, however, a reverse engineer is able to properly disassemble the code with the use of interactive disassemblers like IDA Pro, ignoring the *rogue bytes*.

However, there are some conditions in which no traditional assembly listing can accurately represent the instructions that are executed. Exploiting these conditions we obtain what are called *impossible disassembly* techniques. The code produced using these techniques can however be disassembled, but only using a vastly different representation of the code than what is provided by currently available disassemblers.

The core idea behind these techniques is to make the *rogue byte* part of a legitimate instruction that is executed at runtime. This way the *rogue byte* becomes not ignorable during disassembly. In this scenario any given byte may be a part of multiple instructions that are executed. This is done using *jmp* instructions. The processor, while running the code, will interpret and execute the bytes following the logical flow of the program, so there is no limitation on the number of instructions the same byte can be part of; a disassembler, however, has such limitations since it will usually represent a single byte as being part of a single instruction.

Example:

```

1  EB
2      JMP -1
3  FF
4      INC EAX
5  C0
6  48    DEC EAX

```

Listing 2.3: Impossible Disassembly example

In this simple example the first instruction is a 2-byte *jmp -1* instruction (**EB FF**). Its target is its own second byte. At run time this causes no errors because the **FF** byte is the first byte of the next instruction *inc eax* (**FF C0**).

However, when disassembling, if the disassembler interprets the **FF** byte as part of the *jmp* instruction, it won't be able to interpret it also as the beginning of the *inc eax* instruction. While the **FF** byte is in reality part of both instructions that actually execute, the disassembler is not able to recognise this.

The 4-byte example code increments the **EAX** register, and then decrements it, therefore it is essentially a complex **NOP** sequence. Being a simple, and small, sequence it could be inserted at any location in a program code in order to fool disassemblers. However this sequence it is also easily recognisable by reverse engineers and substituted with **NOP** instructions using IDA Pro or other instruments and/or scripts. Another alternative is to interpret this sequence as data bytes forcing the disassembler to skip it.

However this was only a simple example sequence. More complex and ingenious sequences can be made to fool disassemblers while being harder to detect.

Obscuring Flow Control

Control-flow analysis (CFA) is a static-code-analysis technique for determining the control flow of a program. Modern disassemblers like IDA Pro are able to correlate function calls and extract high-level information about the program knowing how functions are related to each other [26].

Control-flow analysis can however be easily defeated by malware authors.

The Function Pointer Problem Function pointers are a common programming idiom present in programming languages such as **C**, while being extensively used in the background in object oriented languages like **C++** and **Java** [26].

As opposed to referencing a data value, a function pointer points to executable code within memory. Dereferencing the function pointer yields the referenced function, which can be invoked and passed arguments to as in a normal function call. Since, doing so, the function is being invoked indirectly through a variable instead of directly through a fixed identifier or address, such invocation is also known as an "indirect" call. In assembly code this corresponds to a *call* instruction with a function pointer as argument.

Function pointers, however, greatly reduce the information that can be automatically extracted by the disassembler about the program control flow. Moreover, if function pointers are used in specially crafted, or non-standard code, the resulting code can be difficult to reverse-engineer without the use of dynamic analysis techniques.

As a result, function pointers, in combination with other anti-disassembly techniques, can greatly increase the complexity and difficulty of reverse-engineering.

Return Pointer Abuse Among the instructions capable of transferring control within a program we already mentioned the *call* and *jmp* instructions, however there are more [26].

The counterpart to the *call* instruction is *retn*. When a *call* instruction is reached during program execution, a return pointer is pushed on the stack, before jumping to the *call* instruction target. This return pointer in the stack will point to the address of the instruction immediately following the end of the *call* instruction itself. Therefore a *call* instruction can be seen as the combination of a *jmp* and *push*; a *retn* instruction, on the other hand, is the combination of *pop* and *jmp*.

The *retn* instruction pops the last value pushed to the stack and jumps to it; it is therefore typically used to return from a function call, but it could also be used for other purposes. When used for such other reasons the disassembler is generally fooled, because it still will interpret it as a return from a function call. Therefore it won't show any code cross-reference to the target being jumped to. As added benefit the disassembler will also prematurely terminate the function being analysed.

Misusing Structured Exception Handlers Another powerful anti-disassembly technique exploits the Structured Exception Handling (**SEH**) mechanism. Performing program flow control using this mechanism is able to fool both disassemblers and debuggers [26].

SEH provides programs a way to handle error conditions intelligently. *C++* and other programming languages heavily rely on exception handling (and therefore on **SEH**) when compiled for x86 systems.

Exceptions can be triggered for numerous reasons: for example when dividing by zero or accessing an invalid memory region. Moreover, software exception can also be raised by the code itself by calling the *RaiseException* function.

When an exception is raised it makes its way through the **SEH** chain, which is a list of functions specifically designed to handle exception, until it is caught by one exception handler in the chain. Each function in the list can either handle the exception (a.k.a. *catch* it) or pass it to the next handler in the list. *Unhandled exceptions* are the ones that make their way to the last handler. The last exception handler is the code responsible for triggering the 'unhandled exception' message to the user.

Exception handling is used in almost all programs and exceptions happen regularly in most processes (and are handled silently). A malicious actor could, however, exploit this mechanism to achieve covert flow control by adding his own specially crafted handler on top of the **SEH** chain.

This can be done at runtime simply pushing some specific values on the stack, effectively adding a new entry in the Exception handling chain. This procedure, however, is subject to the constraints imposed by the Software Data Execution Prevention (**Software DEP**), which is a

security feature that prevents the addition of third-party exception handlers at runtime. However various workarounds to this protection can be used in the case of handwritten assembly code.

2.2.5 Anti-debugging

This part needs to be summarized since it is too long and too specific.

Another popular anti-analysis technique, besides anti-disassembly, is *anti-debugging*. Malware authors use anti-debugging techniques to recognise when their malicious program is under the control of a debugger or to interfere with the debugger behaviour. This is done in an attempt to slow down the malware analysts who use debuggers to understand how the malware operates. A malware using these techniques usually alter its normal control flow paths or causes crashes if it detects it is running in a debugger, thus interfering with analysis [26].

Windows Debugger Detection

In Windows OS various techniques can be used to detect if a process is being run in a debugger: from exploiting the Windows API itself, to manually checking memory structures for debugging artefacts [26].

Using the Windows API One of the most obvious, and simple, ways to know if a debugger is attached to a process is by using Windows API functions. Inside the Windows API there are, in fact, functions that were specifically designed to detect debuggers; moreover some functions that were originally created with other purposes can also be used for debugger detection [26].

Malware analysts can counter this technique by manually modifying the malware code during execution modifying the resulting flag after the call to make sure the desired path is taken, or by straight up removing/skipping the function call.

Here are some examples of common Windows API functions used for *anti-debugging*:

- **IsDebuggerPresent** This is the simplest API function that can be used for debugger detection. It determines whether the **current** process is being debugged by a user-mode debugger. It does so by getting the value of the field *IsDebugged* from the Process Environment Block (**PEB**) structure. In particular this function returns zero if the process is not running within a debugger context and a non-zero value otherwise.
- **CheckRemoteDebuggerPresent** This API function is similar to the previously described one (*IsDebuggerPresent*). This function checks for a 'remote' debugger on the specified process. The term 'remote' in the name CheckRemoteDebuggerPresent does not imply that the debugger necessarily resides on a different machine; instead, it indicates that the debugger resides in a separate and parallel process. This function takes a process handle as argument, and will check if that process has a debugger attached. It can however be used also to check the current process by passing its handle.
- **NtQueryInformationProcess** This function can retrieve different kinds of information from a process. The first argument for this function is the process handle, the second one is the ProcessInformationClass parameter which specifies the information you want to get. When using the value *ProcessDebugPort* for this parameter, for example, the function will return a zero if the process is not currently being debugged; a non-zero value representing the debugger port number will instead be returned otherwise.
- **OutputDebugString** This function, originally designed to just send a string to a debugger for display, can be used to detect the presence of a debugger. In fact, if there is no debugger attached, the function will internally set the last-error code. In a few lines of code it is thus possible to know if a debugger is present or not:

```

1  DWORD errorValue = 12345;
2  // set custom last error code
3  SetLastError(errorValue);
4
5  // try outputting string on debugger;
6  // if no debugger is present, it will set
7  // the last-error code to a new value
8  OutputDebugString("Test for Debugger");
9
10 if(GetLastError() == errorValue){
11     // a debugger is present
12     ExitProcess();
13 }
14 else{
15     // no debugger was detected
16     RunMaliciousPayload();
17 }
18

```

Listing 2.4: OutputDebugString debugger detection

Manually Checking Structures Malware authors usually don't exploit the Windows API functions for detecting the presence of a debugger, but they prefer checking the PEB structure (and others) by themselves. One of the reasons why they usually don't like using Windows API functions is that API calls can be easily hooked by a rootkit to return false information, thus thwarting this technique [26].

- **Checking the BeingDebugged Flag** The Windows PEB structure contains all user-mode parameters associated with a process, including the process's environment data such as environment variables, addresses in memory and debugger status, among other things. Malware can 'manually' check the *BeingDebugged* flag within the PEB structure to understand if a debugger is attached its process. More precisely if this flag is zero it means that no debugger is attached.

Example of code listing performing 'manual' *BeingDebugged* check:

```

1  mov    eax, dword ptr fs:[30h] ; get PEB address
2  mov    ebx, byte ptr [eax+2]    ; get BeingDebugged flag value
3  test   ebx, ebx              ; test if the value is 0
4  jz     NoDebuggerDetected    ; if 0, no debugger was detected
5

```

Listing 2.5: BeingDebugged manual check

Malware analysts can counter this technique detecting this code sequence in the code and then either manually changing the *BeingDebugged* flag to zero, or forcing the jump to be taken (or not) by manually modifying the zero flag before the jump instruction.

- **Checking the ProcessHeap Flag** The *ProcessHeap*, which is an undocumented location within a reserved array inside the PEB structure, contains the location of the first heap of a process allocated by the loader. This heap can be used for debugger detection since it contains some information telling if it was created within a debugger or not. In particular malware usually check the values of the fields called *ForceFlags* and *Flags*.

To overcome this technique, malware analysts can change the *ProcessHeap* flags manually or use a hide-debug plug-in for their debugger.

- **Checking NTGlobalFlag** Processes started within a debugger run slightly differently than others, therefore they create memory heaps differently. The information needed to determine how to create heap structures is stored at an undocumented location in the PEB. Practically, a value of *0x70* at this location indicates that the process is running within a debugger.

Again, in order to counter this technique, malware analysts can change the flags manually or use a hide-debug plug-in for their debugger.

Checking for System Residue Debugging tools typically leave traces of their presence on the system. Malicious programs can therefore be designed to search for these traces in the system in order to determine when it is being analysed. For example malware can search for references to debuggers in the registry keys [26].

Moreover, malware can also be designed to search the system for files and directories commonly related to debuggers, such as debugger program executables.

Furthermore, malware can also detect debugger residues in live memory, by viewing the current process listing or, more commonly, by performing a *FindWindow* in search for a debugger.

Identifying Debugger Behaviour

Debuggers are very useful to malware analysts because they can be used to set breakpoints in the code or even to single-step through a process running code to ease the reverse-engineering process. These operations, however, modify the process code and are therefore easily detectable [26].

INT Scanning A common anti-debugging technique used by malware authors consists in making the process scan its own code in search for an **INT 3** (opcode *0xCC*). **INT 3** is, in fact, a software interrupt used by debuggers: when setting a breakpoint the debugger replaces the target instruction in the running program with the opcode *0xCC* (INT 3) which causes the process to call the debug exception handler [26].

Malware analysts can counter this technique exploiting hardware breakpoints instead of software ones.

Performing Code Checksums Another anti-debugger technique consists in calculating the checksum of a section of the process' own code. This has the same net effects as scanning the code for software interrupts. However, instead of explicitly searching for a specific opcode (*0xCC*) in the process code, this check performs a cyclic redundancy check (CRC) or a MD5 checksum of the malware code [26].

Again this technique can be countered by using hardware breakpoints instead of software ones, or by modifying the program's control flow at runtime with a debugger.

Timing Checks One of the most widespread techniques for debugger detection is to perform *timing checks*. Processes, in fact, tend to run substantially slower when executed within a debugger context. Moreover analysts usually run programs in single steps in order to better understand the code behaviour, this in turn greatly increases execution time [26].

Using timing checks it is possible to detect a debugger in different ways:

1. Recording 2 timestamps before and after the execution of some operations and then comparing them. If the lag is greater than a specified threshold then a debugger is probably being used.
 2. Recording 2 timestamps before and after raising an exception. If the current process is being debugged then the exception will be handled by the debugger itself more slowly than normal. Moreover, by default, debuggers ask for human intervention when an exception occurs, thus causing huge delays.
- **Using the *rdtsc* Instruction** The most common timing check method uses the *rdtsc* instruction. This instruction returns the number of ticks since the last system reboot. Malware authors thus use it as described above: rdtsc is called twice, once before and once after some other operations, and then the difference between the results is calculated. If too much time has elapsed between the two calls it means that a debugger is probably being used.

- **Using *QueryPerformanceCounter* and *GetTickCount*** These are two Windows API functions that can be used similarly to *rdtsc* for debugger detection.

More precisely *QueryPerformanceCounter* can be called to query a high-resolution counter available to processors which stores counts of activities performed by the processor.

The function *GetTickCount*, on the other hand, returns the number of milliseconds that have elapsed since the last reboot, very similarly to what the *rdtsc* instruction does.

Both of those functions, when used as described above, allow the malware to detect the presence of a debugger.

Anti-debugging through the use of timing checks can be discovered by malware analysts during debugging or static analysis by identifying specific sequences of instructions. Moreover, these checks usually detect debuggers only when the analyst is single-stepping through the code or setting a breakpoint between the two time related instruction calls. This implies that, to counter this technique, malware analysts could avoid setting breakpoints and single-stepping in those regions of code, or modify the result of the timestamps comparison as needed.

2.2.6 Anti-virtual machine

This part needs to be summarized since it is too long and too specific.

Malware analysts often use virtual machines (VMs) or other isolated environments like sandboxes, to analyse malware samples. With the purpose of evading analysis and bypassing security systems malware authors often design their code to detect isolated environments. The techniques used with such purpose are called *Anti-virtual machine* techniques (Anti-VM). Once a virtual machine is detected the evasion mechanism may alter the malware's behaviour, or it may even prevent the malicious code from running altogether [26].

VMware Artefacts

Virtual machines are designed to emulate real hardware functionality. To achieve that, however, some artefacts inevitably remain on the system, which can reveal that a virtual machine is indeed being used. These kind of artefacts can be specific files, processes, registry keys, services, network device adapters etc. [26].

Here are some examples of anti-virtual machine techniques applied to detect VMware virtualization software:

- **Checking for Processes Indicating a VM.** When a VMware virtual machine is running and VMware tools are installed, three VMware-related processes can be found in the system process listing: *VMwareService.exe*, *VMwareTray.exe* and *VMwareUser.exe*. A malicious software can therefore easily detect if VMware is being run searching through the process listing for the *VMware* string.
- **Checking for Existence of Files Indicating a VM.** The VMware default installation path usually also contains artefacts. Searching for the string *VMware* in such location may reveal the use of a virtualized environment.
- **Checking for Registry Keys.** VMware Tools may leave some artefacts also in the registry. More specifically the presence of specific registry entries may reveal the use of VMware.
- **Checking for Known Mac Addresses.** In order to connect a virtual machine to a network it needs to have its own virtual network interface card (NIC). This implies that VMware needs to create a MAC address for the virtual machine, to associate to its NIC. However, depending on VMware configuration, this may lead to the network adapter being able to identify VMware usage. VMware utilises, in fact, addresses with a specific starting sequence which depends on its current version. Therefore a malicious software just needs to check the system MAC address against common VMware values.

In order to counter anti-virtual machine techniques, malware analysts need to apply a two step process: identify the check for VMware artefacts and then 'manually' patch it. For example, depending on the anti-VM technique used, they may patch the malware code while debugging to artificially make all checks pass, or modify the name of VMware processes in order to make them undetectable by the malicious software.

Vulnerable Instructions

The virtual machine monitor program, which monitors the virtual machine execution, has some security weaknesses that may allow malware to detect its usage. In particular, in order to avoid performance issues deriving from fully emulating all instructions, VMware allows certain instructions to execute without being properly virtualized. This in turn means that certain instruction sequences may return different results when running within a VMware virtualized environment than they do on native hardware. This discrepancy can be used by malware authors to detect VMware usage [26].

However, those instructions previously mentioned are not typically used within a malicious program unless it is specifically performing VMware detection, because they are useless if executed in user mode. Therefore avoiding this type of anti-VM technique can be as easy as patching the malicious code to prevent it calling these instructions.

2.2.7 Packers and unpacking

This part needs to be summarized since it is too long and too specific.

Packing programs, commonly known as *packers*, are software programs that take an *executable file* or *dynamic link library (DLL)*, compresses and/or encrypts its contents and then packs it into a new executable file [26].

When packers are used on malicious programs, the malicious code appearance is changed as a consequence of the compression and/or encryption. The packed file will thus hinder basic static analysis and malware detection. Moreover, a packer specifically designed to make the file difficult to analyse may even employ anti-reverse-engineering techniques, such as anti-disassembly, anti-debugging or anti-VM on the resulting compressed version; on top of that some packers, using randomization, are also able to generate different variants of a single file every time it is packed [27].

Malware authors have thus increasingly been using these tools to hide their creations from anti-malware solutions and malware analysts. In order to analyse packed malware, in fact, it must be unpacked first. Properly unpacking a packed program generally is, however, not easy.

A packed file usually contains two basic components:

- A number of data blocks containing the compressed and/or encrypted original executable file.
- An unpacking stub able to dynamically recover the original executable file at runtime.

When the packed file is executed, the unpacking stub is loaded by the OS and begins unpacking the original executable code in memory. When the unpacking has been fully completed the control flow is transferred, with a *jmp*, *call* or the more stealthy *retn* instruction (also referred to as the *tail jump*), to the original file entry point (OEP). This implies that someone attempting to perform static analysis on the packed program, would actually analyse the unpacking stub and not the original code.

Packer types

Commercial and custom made packers can be divided in several levels of complexity depending on the packing techniques used and the additional features they have. The authors of [28] identified 6 packer main types with increasing complexity. Packer types from 1 to 5 allow, sooner or later at runtime, to have a complete view over the original (unpacked) malicious code, meaning that the unpacker stub unpacks all the code at once. However, what makes them differ is the amount and complexity of the encryption (and obfuscation) methodologies used during packing. On the other hand, type 6 packers unpack only a slice of code at a time in memory, never revealing the whole original code altogether. This implies that malware analysts need to take several memory dumps, instead of only one, if they want to get the complete unpacked code.

Another possible classification of packers can be made based on their purposes and behaviours. Following this idea packers can be broadly classified into four categories [29]:

- **Compressors** utilise compression to shrink files while exploiting few or no anti-unpacking tricks. Popular compressors include the Ultimate PE Packer (UPack), Ultimate Packer for Executables ([UPX](#)), and [ASPack](#).
- **Crypters** encrypt and obfuscate the original file contents. No compression is usually done. Malware developers widely use crypters such as [Yoda's Crypter](#) and PolyCrypt PE.
- **Protectors** combine features from both compressors and crypters. Some popular commercial protectors are Armadillo and [Themida](#).
- **Bundlers** are used to pack a software package of multiple executable files into a single bundled executable file. These files within the package can then be unpacked and accessed without extracting them to disk. Some common PE bundlers are [PEBundle](#) and [MoleBox](#).

Packers detection

Packed executables can be detected through a heuristic approach known as *Shannon Entropy Calculation*. Entropy is, generally speaking, a measure of uncertainty, disorder, in a system or program. The idea behind this approach is that compressed or encrypted executables tend to resemble random data, thus they have higher entropy than unencrypted/uncompressed programs. This approach, however, does not tell any information about the packer used to obtain the packed sample [26].

One common way to tackle this problem is through packer signatures checking. Tools like [PEiD](#) and Sigbuster use such method. These tools are, however, not always successful due to the huge number of packer variations and evolutions present in the wild, and the fact that malware authors frequently modify commercially available packers code or create their own packers so that their packed malicious programs do not match any known signature.

Unpacking

Unpacking is the process of restoring the original contents from packed executables in order to allow AV programs and security researchers to analyse the original executable code. There are three different techniques to unpack a packed executable: *automated static unpacking*, *automated dynamic unpacking* and *manual unpacking* ([26], [29]).

Automated static unpacking programs are dedicated routines designed to decompress and/or decrypt executables packed by specific packers, without actually executing the suspicious programs. This method, when it works, is the fastest and most secure method to unpack an executable. Automatic static unpackers are, however, specific to a single packer. Moreover, they are not able to unpack packed samples that were created with the intention to hinder analysis.

Automated dynamic unpackers, instead, use programs to run or emulate the packed executable allowing the unpacking stub to unpack the original executable code in memory. Once the original

executable is unpacked, the in-memory program's code is written on disk, and the automated unpacker reconstructs the original import table.

Most often security researchers prefer to perform manual unpacking. The two most common approaches used to manually unpack a program are:

- Discover what packing algorithm has been used to pack a sample and then write a program/script to revert it. This process is however time consuming.
- Manually run the packed program to allow the unpacking stub to unpack the original code in memory, then dump the process on disk and finally manually modify the PE header so that the program is complete. This process is more efficient than the previous one.

2.2.8 Code Obfuscation

This part needs to be summarized since it is too long and too specific.

Obfuscation is a technique that generally makes programs harder to understand [30], both for humans and automatic tools. To do so, it transforms a program into a new (structurally different) and more difficult to analyse version while retaining the same functionality as the original (the new version of the program is said to be *computationally equivalent* to the original one) [31].

Originally, this technology was conceived for legitimate purposes to protect the intellectual property of software developers; however it has been widely exploited by malware authors to evade detection [32]. Particularly, in order to elude anti-malware scanners, malware can, using obfuscation techniques, evolve their body into new generations [33], which eventually can be even harder to disassemble and analyse.

Obfuscation techniques can be broadly divided into 2 main sub-categories:

- *Data-based* obfuscation
- *Control-based* obfuscation

However, malware authors usually combine those 2 types of obfuscation techniques in complex and difficult ways to strengthen the resulting obfuscation [34].

Data-Based Obfuscation

Data-based obfuscation techniques focus on modifying data values and non-control computations. In the following paragraphs some common data-based obfuscation techniques will be discussed.

Constant Unfolding *Constant folding* is a technique commonly used by compilers to optimize a program's code. It does so by replacing expressions with results known at compile time with the results themselves [34].

For example, a compiler usually transforms the following expression 2.6, into 2.7.

```
1 x = 4 * 5;
```

Listing 2.6: Before constant folding

```
1 x = 20;
```

Listing 2.7: After constant folding

Constant unfolding is, instead, an obfuscation technique that performs the exact inverse operation: it replaces the constants in the program's code with some expressions having the constant as a result.

For example, the listing 2.8, after *Constant unfolding* may become 2.9. The two listings are equivalent. Moreover, there is an infinite amount of listings equivalent to 2.8 that can be generated following this principle.

```
1 push 0h
```

Listing 2.8: Before constant unfolding

```
1 push 0F9CBE47Ah
2 add  dword ptr [esp], 6341B86h
```

Listing 2.9: After constant unfolding

Data-Encoding Schemes The previously described technique is, however, easily defeated by simply applying the standard compiler’s constant folding optimization. This is possible because both the data encoding and decoding functions ($f(x) = x - 6341B86h$ and $f_{-1}(x) = x + 6341B86h$ respectively) were present in the code one after the other. The use of *fully Homomorphic* mappings (operation-preserving mappings) allow us to perform some operations on the encoded data before decoding it back, thus overcoming the previous technique’s flaw. *Fully Homomorphic* mappings are however still not widely used because still too inefficient [34].

Dead Code Insertion *Dead code elimination* is another common compiler optimization technique. Its objective is to remove program statements/expressions that have no real effects on the program operation and final results [34].

For example, the listing 2.10, using *dead code elimination* would become 2.11.

```
1 int f(){
2     int x, y;
3     x = 1;    // this assignment is useless, here x is dead
4     y = 2;    // y is never used, it is thus dead.
5     x = 3;
6     return x; // x is live
7 }
```

Listing 2.10: Before dead code elimination

```
1 int f(){
2     return 3;
3 }
```

Listing 2.11: After dead code elimination

Obfuscators, on the other hand, use the so-called *dead code insertion* technique in an attempt to make the code harder to follow. This technique performs the inverse operation with respect to *dead code elimination*, adding dead code in the original program’s code.

However, when used alone, this techniques produces an obfuscated program that can be efficiently de-obfuscated by using the compiler’s dead code elimination optimization.

Arithmetic Substitution via Identities This technique aims at replacing certain operators with combinations of other operators with equal net result. Exploiting the equivalence between different combinations of different operators the code can be changed arbitrarily without changing the effective program operation and final result [34]. Here are some examples of operators equivalences:

```
1 -x == ~x + 1
2
3 x-1 == ~-x
4
5 x+1 == -~x
6
7 rotate_left(x,y) == (x << y) | (x >> (bits(x) - y))
8
9 rotate_right(x,y) == (x >> y) | (x << (bits(x) - y))
```

Listing 2.12: Operators equivalences

Register Reassignment Another simple obfuscation technique is called *register reassignment*. An obfuscator using this technique switches the registers used throughout the code at every application, while keeping the same program code and behaviour [33].

An analyst/attacker using wildcard searching, however, easily defeats this technique.

Instruction Substitution *Instruction substitution* creates variants of a program's original code by replacing some instructions with other equivalent ones [33].

Pattern-Based Obfuscation *Pattern-based* obfuscation is another commonly used technique similar in principle to *instruction substitution*, but more complex. It consists in constructing *patterns* (transformations) that map single or multiple adjacent instructions into a more complex, computationally equivalent, sequence of instructions [34].

For example, the sequence 2.13 might be converted into 2.14, as well as into 2.15 or even 2.16.

```
1 push reg32
```

Listing 2.13: Original sequence

```
1 push imm32
2 mov dword ptr [esp], reg32
```

Listing 2.14: Obfuscation using pattern 1

```
1 lea esp, [esp-4]
2 mov dword ptr [esp], reg32
```

Listing 2.15: Obfuscation using pattern 2

```
1 sub esp, 4
2 mov dword ptr [esp], reg32
```

Listing 2.16: Obfuscation using pattern 3

Moreover, patterns can be arbitrarily complicated. For example a listing such as 2.17, could be substituted by the more complex 2.18.

```
1 sub esp, 4
```

Listing 2.17: Original sequence #2

```
1 push reg32
2 mov reg32, esp
3 xchg [esp], reg32
4 pop esp
```

Listing 2.18: Obfuscation of sequence #2

Malware authors (and also software developers wishing to protect their intellectual property) can use hundreds of patterns in the same program. Moreover, most protections randomly apply patterns so that obfuscating the same program multiple times yields different results. On top of that, patterns can also be applied iteratively: after transforming the original code **C** into **C'** using pattern **P**, another pattern **P'** can be applied to **C'** in order to obtain **C''**, and so on.

Some patterns preserve *semantic equivalence*, meaning that the CPU state will be the same when executing them or the original code. Some other patterns, however, do not. Therefore, depending on the code logic, some substitutions are safe (meaning that the program behaviour and final results are preserved) while others are not. This makes the job of an obfuscator challenging.

Control-Based Obfuscation

Standard static analysis tools generally make assumptions similar to the ones human reverse engineers make when analysing code. Compilers, in fact, predictably translate control flow constructs and data structures. As a result, reverse engineers (and static analysis tools) can easily recognise the original code high level control flow. *Control-based obfuscation* transforms the code control flow structures in non standard ways in order to complicate both static and dynamic code analysis [34].

Some examples of standard static analysis tools assumptions are:

- The *CALL* instruction is always used with the sole purpose of invoking functions.
- Both sides of a conditional branch may feasibly be taken at runtime.
- Function calls almost always return.
- All control transfers target code locations, not data locations.
- Exceptions are used in standard and predictable ways.
- etc..

By violating these assumptions, *control-based obfuscation* techniques confuse disassemblers and other static analysis tools making the analysis more difficult.

Functions In/Out-Lining Reverse engineers frequently rely on control-flow and call graphs to better understand a program's high-level logic. In particular a call graph represents calling relationships between subroutines (functions) in a computer program. Each node of a call graph represents a procedure and each edge (f, g) indicates that procedure f calls procedure g . By making the call graph harder to interpret, obfuscators can hinder the reverse engineer's capability of understanding the program behaviour [34]. To do so one could:

- **Inline functions.** The code belonging to a subfunction is merged into the code of its caller. If a subfunction is called multiple times, however, the code size can quickly grow.
- **Outline functions.** A subpart of a function is extracted and transformed into an independent function and replaced by a call to the newly created function.

Using these two operations in combination on a program's code results in a degenerated call graph with no clear logic. Moreover, also the functions' prototypes can be modified adding extra fake arguments, reordering arguments and so on, to further hide the high-level logic.

Destruction of Sequential and Temporal Locality Usually, in non-obfuscated code, the instructions of a single basic block lie one after the other (*sequential locality*), and basic blocks related to one another (such as successive blocks) are close to each other (*sequential locality of temporally related code*). This is done in order to maximize the instruction cache locality and reduce the number of branches in the final code. Reverse engineers thus can usually rely on the fact that all the code responsible for a specific operation will reside in a single region [34].

Violating this assumption introducing unconditional branches that break sequential locality and temporal locality of multiple basic blocks makes manual analysis more difficult. However, by constructing the control-flow graph and removing spurious unconditional branches the original control flow can be restored.

Processor-Based Control Indirection Instructions like *JMP* (branch) and *CALL* (save instruction pointer and branch) are, for most processors, the 2 essential control flow transfer primitives. In order to make analysis more difficult, one could obfuscate these primitives for example using dynamically computed branch addresses or by emulating them.

For example the instruction *JMP* instruction 2.19, can be replaced by the (almost) semantically equivalent listing 2.20.

```
1 jmp target_addr
```

Listing 2.19: Processor-based control indirection before

```
1 push target_addr
2 ret
```

Listing 2.20: Processor-based control indirection after

Operating System-Based Control Indirection As already seen when talking about anti-disassembler techniques, obfuscation can also exploit operating system primitives and structures. For example, the Structured Exception Handler (*SEH*), Vectored Exception Handler (*VEH*) and Unhandled Exception Handler are commonly used to obfuscate the control flow of Microsoft Windows executables (in Unix-like systems the signal handlers *setjmp* and *longjmp* are commonly used instead) [34].

Subroutine Reordering *Subroutine reordering* is an obfuscation technique that randomly changes the order of a program's subroutines in the original code. This technique can thus generate $n!$ code variations, where n is the number of subroutines [33].

Opaque Predicated An *opaque predicate* is a non-trivial boolean expression with a constant result (always true or always false) known only at compilation/obfuscation time. Combining it with a conditional *jmp* instruction introduces an additional branch in the control flow graph (*CFG*). We already briefly talked about this specific combination when talking about the *Jump instruction with a constant condition* anti-disassembly technique. The added branch should look as real as possible in order to elude detection, and it can be used to insert junk code or to form cycles in the control-flow graph to better hide the original program's logic [34].

Simultaneous Control-Flow and Data-Flow Obfuscation

Data-flow obfuscation and *Control-flow obfuscation* techniques are commonly used together to complicate analysis.

Inserting Junk Code This technique consists in introducing a dead code block (meaning that it will never be executed at runtime) between two other code blocks. Typically used in conjunction with *opaque predicates*, this technique is used to hinder a disassembler that is disassembling an invalid path. Moreover, the junk code typically contains partially invalid instructions, or branches to invalid addresses with the objective of over-complicating the *CFG* [34].

```
1 push eax
2 xor eax, eax
3 jz 9
4 ;<junk code start>
5 jg 4
6 inc esp
7 ret
8 ;<junk code end>
9 pop eax
```

Listing 2.21: Junk Code example

The listing 2.21 presents an example of this technique. More precisely the instruction at line 2 (`xor eax, eax`) zeroes the *EAX* register setting clearing the zero flag (it is set to 0); therefore the conditional jump (`jz 9`) at line 3 is always taken at runtime. The immediately next instructions are therefore junk code.

Control-Flow Graph Flattening *Control-flow graph flattening* consists in replacing all control structures within a sub-part of the control flow graph with a single switch statement commonly called *dispatcher*. This is done to hide the true basic blocks relationships within the dispatcher. When using this technique, first a subpart of the program's control flow graph is selected to be substituted by the dispatcher. Some transformations may then be applied to the basic blocks inside the chosen sub-graph (they split or merged) to further complicate analysis and finally each basic block updates the dispatcher's context to reflect the relative basic block relationships. The final resulting graph offers no clues about the structure of the algorithm, but has the same logic [34].

CFG flattening is frequently used, together with *opaque predicates*, to insert dead code paths in the CFG.

Code Transposition An obfuscator using the technique called *code transposition* effectively reorders the sequence of a program's original code instructions without changing its behaviour [33]. To achieve this two approaches are commonly followed:

- Randomly shuffling the instruction and then recovering the original execution order by inserting unconditional branches. This is easily defeated restoring the original program by removing (and following) the unconditional branches.
- Choosing and reordering independent instructions that have no impact on the others. This approach is harder to implement given the complexity of finding independent instructions, but it is more effective.

Code Virtualization *Code virtualization* consists in transforming a program's binary code (compiled for a specific machine) into a different binary code that is understood by a virtual machine. More specifically, the instruction set from the source machine is converted into a new instruction set, understood by the target virtual machine. This can be done using multiple types of virtual machines with different instruction sets. This means that a specific block of, for example, Intel x86 instructions can be converted into a different instruction set for each machine, preventing an analyst/attacker from recognizing any generated virtual opcode after the transformation from x86 instructions [34].

Usually, some specific blocks of the program's code are virtualized (and not the whole program) and inserted back into the program alongside the associated interpreter. At run time, the interpreter assumes execution control and translates the virtualized code back to the original byte code.

When an analyst/attacker tries to decompile a virtualized block of code, however, he will not find the original x86 instructions. Instead, he will find a completely new instruction set which is not recognized by him or any other special decompiler. This will force the attacker to identify how each opcode is executed and how the specific virtual machine works for each protected application.

Some examples of code virtualization tools include [VMProtect](#) and [CodeVirtualizer](#).

Code Integration *Code integration*, one of the most sophisticated obfuscation techniques, was first introduced by *Win97/Zmist* malware. A malware using this technique first decompiles the target program into a set of manageable objects, it then inserts itself between them and finally it reassembles the code [33].

2.2.9 Obfuscated Malware

The huge amount of malware released in the wild since the creation of the first virus in 1960s can be split into two generations. More specifically, the first generation malwares were static, their code and behaviour did not change. The more sophisticated second generation malwares, on the other hand, change their internal structure between one variant and the other maintaining the same malicious behaviour in order to avoid detection.

Encrypted Malwares

The first second-generation malwares ever existed exploited encryption in order to evade detection by signature-based antivirus scanners.

In this approach, an encrypted malware typically consists of two parts: the encrypted main body and a decryption code (also called *decryptor*). The *decryptor*'s objective is to recover the original malware code from the encrypted body whenever the infected file is run [31].

Moreover, to hide from signature-based scanners, encrypted malware encrypts its code using a different key at each infection, thus creating a unique encrypted body. The decryption routine (*decryptor*), however, remains the same from one generation to another. This means that encrypted malwares can be detected with signature-based scanners by searching for the decryptor's code pattern [22].

The first known malware to exploit encryption for detection evasion was CASCADE which spread in the 1980s and early 1990s.

Encryption of malware code is often used in conjunction with the use of packers.

Packed Malwares

Malware authors are nowadays increasingly exploiting packers (or even multiple packers at once) to produce numerous variants of the same original malware code [5].

As stated by Perdisci, et al [35], more than 80% of the new malware currently discovered are actually packed versions of already existing malware.

Packers are used to compress the original file into a smaller size and, moreover, encryption is sometimes applied to the compressed version of the file in order to make the unpacking process more difficult.

However, it is not uncommon to see malware authors writing and using custom packers. This fact can be used by analysts to detect if a file is malicious, without further analysis, based on the fact that benign software vendors would almost never use custom packers. On the other hand, many malware authors frequently use commercial, and readily available, packers to generate malware variants.

Oligomorphic Malwares

Malware authors tried to overcome the short comings of encrypted malware developing malware that can mutate the used decryptor from one variant to another. Initially the decryptor could only be changed slightly. However, a common method used by *oligomorphic malware*, also called '*semipolymorphic*', to provide more diverse decryptors is, in practice, to randomly select one decryption routine at infection time from a set of pre-defined different decryptors [22].

However, this type of malware is able to generate at most few hundreds different decryptors. For example the virus called *Win95/Memorial* was capable of constructing up to 96 different decryptor patterns. This means that signature-based detection techniques are still able to detect *oligomorphic malwares* by generating the signature of all the decryptors utilised by the malware strain [31]. Still, signature based techniques are not an effective approach to detect *oligomorphic malware*.

The virus named *Whale* was the first known malware to make use of this technique. It carried a few dozens of different decryptors and picked one randomly at infection time [5].

Polymorphic Malwares

The *oligomorphic* malware limitations lead malware authors to develop a more advanced type of malware called *polymorphic*.

In *polymorphic* malware countless numbers (millions) of distinct decryptors can be generated by using obfuscation methods including, for example, *dead-code insertion*, *register reassignment*, *subroutine reordering*, *instruction substitution*, *code transposition/integration*, etc. to avoid signature based detection [31].

Polymorphic malware, like *oligomorphic* malware, consists of two parts: the malware encrypted main body and the decryptor. The decryptor is again run once the malware is executed and it enables the execution of the original malware code decrypting the encrypted body. When replication occurs, the malware encrypts its code with a different key, generates the new associated decryptor and encloses it in the new malware variant code. The malware appearance is thus changed at each infection [22].

In order to have a wide range of decryptors, *polymorphic* malware typically use a powerful toolkit called 'the *Mutation Engine (MtE)*'. In particular, during malware replication, the mutation engine is used to create the new decryptor which is appended to the new malware variant code. The mutation engine is, in fact, responsible for rearranging the decryptor code using different obfuscation techniques in order to prevent signature based detection.

Even though *polymorphic* malware can create a large number of different decryptors effectively hindering signature matching techniques, still the constant malware body, which appears after decryption, can be used for detection. In particular, by using emulation techniques, the tool execute the malware in a '*Sandbox*' without resulting in any harm to the system. As soon as the constant malware body is decrypted and loaded into memory, the common detection techniques, such as signature based scanning, can be applied [5].

Various armouring techniques are thus used by malware authors to prevent detection by emulation, however most antivirus scanners are now capable of addressing also these techniques effectively defeating *polymorphic* malware.

The first known malware to exhibit *polymorphism* is called *1260* and was written in 1990.

Metamorphic Malwares

After the *oligomorphic* and *polymorphic* malware types were effectively defeated, malware authors designed a new and more advanced approach: *metamorphic* malware. This, similarly to *polymorphic* malware, uses obfuscation techniques to create new variants of the original malware in order to evade detection [31].

However, in this case, instead of generating new decryptors, it is the malware body itself to be mutated through generations to appear different while having the same behaviour and functionality. *Metamorphic* malware is in fact said to be *body-polymorphic*. In practice the malware code logic is maintained while its appearance is changed using obfuscation techniques such as *dead-code insertion*, *register reassignment*, *code transposition* and more. This way, every generated malware variation appears different making signature based detection ineffective [22].

However, *metamorphic* malware, in order to efficiently evolve its code it need to be able to recognise, parse and mutate its own body during propagation. This is far from being easy. Moreover, creating a true *metamorphic* malware without arbitrary increasing its code size is also challenging [5].

Moreover, *metamorphic* malware is also capable of interleaving its own code inside host programs, thus making detection even harder.

The first malware to exhibit metamorphic behaviour was called *Win95/Regswap* and was developed in 1998.

Chapter 3

Detection Techniques

Malware detection is the process of identifying malicious code from benign code. This is done in order to protect systems and being able to recover from the malicious code effects [5].

In order to counter malware attacks and threats, in recent years many anti-malware tools have been developed. Many of these are based on static features (such as signatures) with the assumption that most malware is static, it doesn't mutate/change significantly at infection/replication time [22].

However, attackers are nowadays increasingly using the more sophisticated second generation malwares, which strongly mutate at each infection. Researchers and anti-malware software developers are thus focusing their attention on the creation of more advanced tools capable of detecting this type of evolving malware.

Commercial Portable Executable (PE) malware detectors consist of a hybrid of static and dynamic analysis engines. Static detection - which is fast and effective at detecting a large fraction of malware - is usually first employed to flag suspicious samples. Static detection involves analysing the raw PE image on disk and can be performed very quickly, but it is vulnerable to code obfuscation techniques.

Dynamic detection, by contrast, requires running the PE in an emulator and analysing behaviour at run time. When dynamic analysis works, it is less susceptible to code obfuscation, but takes substantially greater computational capacity and time to execute than static methods. Moreover, some files are difficult to execute in an emulated environment, but can still be statically analysed. Consequently, static detection methods are typically the most critical part of an endpoint's malware prevention pipeline.

3.1 Integrity Checker

When compromising a computer system or network some changes are inevitably made within the target environment. This implies that systems, like *integrity checkers*, that rely on actively monitoring changes made to existing files within the target operating system, can be used to perform intrusion detection [5].

Generally, *integrity checkers*, use hashing functions like the *md5* sum, *Sha1* or *Sha256* to calculate the digest of files and/or executables which are then stored in a database of digests. Programs and files digests are then periodically re-calculated and compared against the ones in the database looking for modifications. If the digest of a file is different and no software updates nor patches were applied, then the file was probably tampered with.

Integrity checkers present a number of challenges:

- The system state in which the initial file digests are calculated has to be considered clean. However this is difficult to be guaranteed.

- The application of system (and software) updates and patches, which modify system files and programs, must be followed by an update of the digests database, otherwise there will be a very high false positive rate.
- The digests database needs to be stored securely and there has to be an offline (and safe) backup, otherwise there would be a single point of failure.

Integrity checking can be considered as an important tool for detecting any system modifications, but it is more an incident recovery method rather than a malware intrusion/infection prevention method.

3.2 Signature-based Detection

Signature-based detection is the simplest and most widely used method in commercial anti-virus software (together with *heuristic-based* techniques) but is becoming less and less effective as the number of malware variants and second generation malwares increases [36].

Signature-based detection relies on *signatures*, represented by specific unique byte code sequences/strings extracted from malware samples, to detect the presence of malicious files in a system. *Signatures* are typically created using static analysis techniques and are selected to be long enough to uniquely characterize a specific malware families with respect to benign programs.

The *signatures*, which are created by malware experts from a significant number of already identified malware samples, are saved in a *signature database* and deployed in anti-malware tools. Anti-malware tools in turn scan the files in the target system and consider as malicious any file that matches one of the known signatures [5]. This implies that the database of signatures must be maintained and frequently updated, especially whenever new malware variants are identified and new signatures are generated in order to detect them.

Some *signature-based* algorithms require an exact match between the signature of the analysed sample and one of the known signatures, others instead make use of *wildcards* characters to detect slight variations. Some second generation (evolving) malwares have been detected in the past by using wildcards, e.g. *W32/Regswap*.

This approach is fast, easy to use and has a high positive rate, however, since the number of known malwares is increasing so fast, it is quickly becoming time-consuming, expensive and impractical. Moreover, this is a completely reactive technique which is unable to counter threats/attacks from new malwares families/variants until they cause damages. Additionally, most second-generation malwares are able to escape this type of detection [22].

3.2.1 Yara Rules

This part needs to be summarized since it is too long and too specific.

YARA is a widely accepted open-source *signature-based* malware analysis tool which has emerged in recent years thanks to its flexible and customisable nature. It allows malware analysts/researchers to develop malware "descriptions" based on text or binary patterns, commonly referred to as *Yara rules*. *Yara rules*, which combine simple regular expressions matching with logic rules, can be used to identify specific malware families, the presence of *CVEs*, specific functionality signatures or even generic maliciousness indicators. Given the success obtained by this technique, many commercial malware analysis tools nowadays support *Yara rules* natively [37].

Yara rules can be generated either manually or automatically. Generating rules manually obviously requires high expertise, whereas generating them automatically using tools is a relatively easy task. However, automatically generated rules are not guaranteed to be effective and may require post-processing operations for their optimization [38].

Malware analysts typically create *Yara rules* manually by reverse engineering malware samples looking for common *Indicator of Compromise (IoC)* strings. This is followed by the development

and iterative refinement of the rules which are considered effective based on their coverage and false positive rate on a dataset of malicious, benign and out-of-family samples. Developing effective *Yara rules* can therefore be challenging and very time consuming, even for expert users with years of experience [39].

Yara Rules syntax

Listing 3.1 presents an example of the syntax of a simple *Yara rule*.

```

1  rule RuleName
2  {
3      meta:
4          description = "description of rule"
5          author = "name"
6          date = "dd/mm/yyyy"
7          reference = "url"
8
9      strings:
10         $text_string1 = "text1 you wish to find in malware"
11         $text_string2 = "text2 you wish to find in malware"
12
13        $hex_string1 = {hex1 you wish to find in malware}
14        $hex_string2 = {hex2 you wish to find in malware}
15
16        $reg_exp_string1 = /regular expression1 you wish to find in malware/
17        $reg_exp_string2 = /regular expression2 you wish to find in malware/
18
19      condition:
20         $text_string1 or $text_string2 or
21         $hex_string1 or $hex_string2 or
22         $reg_exp_string1 or $reg_exp_string2
23 }
```

Listing 3.1: YARA Rules Syntax

As it can be seen in the above example, *Yara rules* must start with the keyword '*rule*', followed by the actual *RuleName*, which is the rule identifier. The *RuleNames* follow the same lexical conventions of the *C* programming language. They are, in fact, case sensitive, they cannot exceed 128 characters and they can contain only alphanumeric characters (with the addition of the underscore character), with the exception of the first character which cannot be a digit. Furthermore there is a list of *YARA* reserved keywords that cannot be used as identifiers [40].

Yara rules main body contains three sections: *meta*, *strings* and *condition*.

Meta section The rule author can include additional information about the rule as a list of attribute-value pairs, also called *metadata*, in the *meta* section, at the top of the rule. The values can be strings, integers or boolean values. The metadata, however, cannot be used in the condition section since that is not their purpose [41].

Some commonly used meta tags are, for example, "author" and "description", which convey information about the author and purpose of the rule. Moreover, malware analysts sometimes also leave tags with the hashes of the malicious files used for the creation of the rule, or references to blog posts with similar information [37].

Strings section This section contains the strings/patterns/signatures that a file must contain to 'trigger' the rule. This section is optional and can be omitted if it is not necessary. *YARA* supports searching for 3 string types: *Hexadecimal Strings*, *Text (ASCII) Strings* and *Regular Expressions*.

- *Hexadecimal Strings*: *Hexadecimal Strings* will match hexadecimal characters/sequences of raw bytes in the file being analysed. Example:

```

1 rule ExampleRule
2 {
3   strings:
4     $my_text_string = "text here"
5     $my_hex_string = { E2 34 A1 C8 23 FB }
6
7   condition:
8     $my_text_string or $my_hex_string
9
10 }
```

Listing 3.2: YARA Hexadecimal

Three special flexible formats, namely *wildcards*, *jumps* and *alternatives*, can be used to complement the search.

- **Wildcards** are represented by the ‘?’ symbol. They indicate that some bytes in the pattern are unknown and should match anything. For example:

```

1 rule WildcardExample
2 {
3   strings:
4     $hex_string = { E2 34 ?? C8 A? FB }
5
6   condition:
7     $hex_string
8
9 }
```

Listing 3.3: YARA Hexadecimal Wildcard

- **Jumps** are used in circumstances when the values of the pattern are known but their length varies. For example:

```

1 rule JumpExample
2 {
3   strings:
4     $hex_string = { F4 23 [4-6] 62 B4 }
5
6   condition:
7     $hex_string
8
9 }
```

Listing 3.4: YARA Hexadecimal Jump

In particular, in listing 3.4, the value ‘[2-3]’ indicates that any arbitrary sequence from 2 bytes to 3 bytes long can occupy the sequence at that position.

- **Alternatives**, whose syntax resembles regular expressions, are used in situations in which the author wants to provide different alternatives for a given fragment of the hex string. For example:

```

1 rule AlternativesExample
2 {
3   strings:
4     $hex_string = { F4 23 ( 62 B4 | 56 ) 45 }
5
6   condition:
7     $hex_string
8
9 }
```

Listing 3.5: YARA Hexadecimal Alternatives

In particular, in listing 3.5, the value ‘(62 B4 | 56)’ indicates that one sequence between ‘62 B4’ and ‘56’ can occupy the sequence at that position.

- **Text Strings:** Text strings are generally readable sequences of ASCII characters which are then matched in the condition section [37].

Example of an ASCII-encoded, case-sensitive string:

```

1   rule TextExample
2   {
3     strings:
4     $text_string = "foobar"
5
6     condition:
7     $text_string
8   }
9

```

Listing 3.6: YARA Text Strings example

Additionally, to specify how *YARA* should search for strings, some modifiers can be added at the end of a string definition. Moreover, even more than one modifier can be used in combination to keep rule as simple and readable as possible. Here are described some of the available modifiers:

- *nocase*: Text strings in *YARA* are, by default, case-sensitive. However it is possible to search for strings in case-insensitive mode by appending the modifier '*nocase*' at the end of the string definition, in the same line. Example:

```

1   rule CaseInsensitiveTextExample
2   {
3     strings:
4     $text_string = "foobar" nocase
5
6     condition:
7     $text_string
8   }
9

```

Listing 3.7: YARA nocase example

- *wide*: The '*wide*' modifier can be used to search for strings encoded with two bytes per character (also known as *wide character strings*), something which is typical in many executable binaries.

For example, if the string "Borland" appears in the file encoded as two bytes per character, then the following rule will match:

```

1   rule WideCharTextExample1
2   {
3     strings:
4     $wide_string = "Borland" wide
5
6     condition:
7     $wide_string
8   }
9

```

Listing 3.8: YARA Wide Character Strings

- *xor*: *YARA* can also encode text before searching it in the analysed file. The '*xor*' modifier, for example, can be used to search for strings with a single byte XOR applied to them.

The following rule will search for every string resulting from a single-byte XOR applied to the string "This program cannot":

```

1   rule XorExample1
2   {
3     strings:
4     $xor_string = "This program cannot" xor
5
6     condition:
7     $xor_string
8   }
9

```

Listing 3.9: YARA XOR-ed Strings

- **base64**: The ‘*base64*’ modifier can be used to search for strings that have been base64 encoded.

The following rule will search for all the possible base64 permutations of the string “This program cannot”:

```

1  rule Base64Example1
2  {
3    strings:
4      $a = "This program cannot" base64
5
6    condition:
7      $a
8  }
9

```

Listing 3.10: YARA Base64 encoded Strings

- **Regular Expressions**: Starting from version 2.0 YARA has its own regular expression engine, which is one of its most powerful features. *Regular expressions* are defined in the same way as text strings, but enclosed in forward slashes instead of double-quotes [38].

Example:

```

1  rule RegExpExample
2  {
3    strings:
4      $re1 = /md5: [0-9a-fA-F]{32}/
5      $re2 = /state: (on|off)/
6
7    condition:
8      $re1 and $re2
9
10

```

Listing 3.11: YARA Regular Expression

Conditions section The last section of *YARA rules*, which is the only required one, contains the rule conditions that determine when the rule gets triggered. These conditions are Boolean expressions similar to those used in programming languages [41]. Through the use of all the usual logical and relational operators, conditions can be made arbitrary complex in order to accommodate for the specific author needs [37].

Inside the *Conditions* section, among other things, it is possible to:

- **Count strings** Sometimes it is necessary to know how many times a string appears in the analysed file, not only if it is present or not. The number of occurrences of each string defined in the string section can be retrieved by using a variable whose name is the string identifier but with a # character in place of the initial \$ character.

For example:

```

1  rule CountExample
2  {
3    strings:
4      $a = "dummy1"
5      $b = "dummy2"
6
7    condition:
8      #a == 6 and #b > 10
9
10

```

Listing 3.12: YARA Count Example

- **Check String at specific offset/in offset range**: Sometimes we need to know if a particular string is available at some specific offset of the file or at some virtual address within the process address space. In such situations it is possible to use the ‘*at*’ operator.

The '*in*' operator, on the other hand, allows to search for a specific string within a range of offsets or addresses, rather than at an exact one.

Examples:

- The rule in listing 3.13 will find whether the '*a*' string is located at offset 100 and '*b*' at offset 200 of the running process.

```

1  rule AtExample
2  {
3    strings:
4    $a = "dummy1"
5    $b = "dummy2"
6
7    condition:
8    $a at 100 and $b at 200
9  }
10

```

Listing 3.13: YARA At Example

- The rule in listing 3.14 will find the '*a*' in the memory location between *0* and *100* and '*b*' between *100* and *filesize* of the running process.

```

1  rule InExample
2  {
3    strings:
4    $a = "dummy1"
5    $b = "dummy2"
6
7    condition:
8    $a in (0..100) and $b in (100..filesize)
9  }
10

```

Listing 3.14: YARA In Example

- **Check file size:** '*filesize*' is a special variable that can be used in rules conditions, which holds the size of the file being scanned in bytes.

Example:

```

1  rule FileSizeExample
2  {
3    condition:
4    filesize > 200KB
5  }
6

```

Listing 3.15: YARA Filesize Example

- **Check a set of strings:** When it is necessary to know if a file contains a certain number of strings from a given set the '*of*' operator can be used.

Example:

```

1  rule OfExample
2  {
3    strings:
4    $a = "dummy1"
5    $b = "dummy2"
6    $c = "dummy3"
7
8    condition:
9    2 of ($a, $b, $c)
10
11

```

Listing 3.16: YARA Of Example

Additional modules YARA's code functionality can be extended through the use of modules. Some modules like the *PE module* and the *Cuckoo module* are officially distributed with YARA, however additional ones can also be created.

Here are mentioned some useful (in this document context) Yara modules:

- **YARA with PE** Starting with version 3.0, YARA can parse Portable Executable (PE) files [40]. For example, the rule in listing 3.17 will check for the string "abc", will parse the PE file and look for "CreateProcess" and "httpsendrequest" function names in the import sections 'Kernel32.dll' and 'wininet.dll', respectively.

```

1 Import "PE"
2
3 rule PE_Parse_Check
4 {
5     strings:
6         $string_pe="abc" nocase
7
8     condition:
9         pe.imports("Kernel32.dll", "CreateProcess") and
10        pe.imports("wininet.dll", "httpsendrequest") and
11        $string_pe
12
13 }
```

Listing 3.17: YARA with PE

- **YARA with PEiD** YARA can also be integrated with *PEiD* to check what packer was used to compile the malicious/suspicious executable [40].

Yara Rules Advantages and Disadvantages

1. **Advantages** of *Yara rules*: *Yara rules* offer several advantages over other malware analysis techniques. Here are some of the most notable ones [38]:

- *Yara rules* allow malware analysts to write flexible and custom rules in an easy and efficient way.
- *Yara rules* are an open standard which work on most of the major operating systems such as Windows, Linux and Mac OS.
- *Yara rules* can be easily integrated into Python and C/C++ programming languages.
- *Yara rules* can be used both for static and dynamic malware analysis.
- Several automatic tools have been developed, and are readily available, to generate *Yara rules* easily and efficiently.
- There are various public repositories of *Yara rules* which offer readily available rules for malware analysis.

2. **Limitations** of *Yara rules*: *Yara rules*, however, have also some limitations. Here are some of the most notable ones [38]:

- *Yara rules* are commonly written based on *IoC* (*Indicator of Compromise*) strings, however, malware authors can easily obfuscate, replace or encrypt these *IoC* strings in their creations in order to evade detection. This could make these rules less effective.
- *IoC* strings are usually extracted from existing malware samples/families through the use of reverse engineering techniques. The use of these techniques in manually creating effective rules, however, requires a highly specialised skill-set and years of experience.
- The effectiveness of *Yara rules* is generally influenced by the types and number of *IoC* strings included in the rules. However, achieving the right balance of both is a challenging task.
- *Yara rules* are effective in detecting malware which matches known malware signature. It may, however, completely miss new and unique malware variants.

Yara Rules Automatic Generators

There are various automatic *Yara rules* generator tools available. In the following the most notable ones will be briefly described:

***YarGen* Tool** *YarGen* python-based tool exploits some smart techniques, namely fuzzy regular expressions, Naive Bayes classifier and Gibberish Detector, to generate *Yara rules*.

The produced rules include features (strings and opcodes) common to malware samples that don't match with the provided goodware databases. A predefined number of features (generally up to 20 strings) are selected, based on their potential utility and a number of heuristics, to be combined and used by the rule in order to maintain a reasonable operation speed.

This tool is able to generate two types of rules: *basic rules* and *super rules*. *Basic rules* can generally target specific malware samples, where *super rules* are able to target a set of malware samples or a whole malware family [38].

The *yarGen* authors encourage its use as a starting point for rule construction, followed by manual adjustments and to refine *yarGen*'s output [39].

1. *YarGen* tool advantages:

- It allows generation of *Yara rules* based on both opcodes and strings.
- It supports the use of PE (portable executable) modules, which are used to interpret Windows operating system executables such as *DLL* and *COM* files.
- It can be integrated with other anti-malware software in order to improve its effectiveness.
- It reduces the false positive rate by checking all strings against databases of goodware samples.
- It is deployed as a simple and easy-to-use python script that can be run through a command-line interface.

2. *YarGen* tool disadvantages:

- It requires post-processing of the generated rules for increasing their effectiveness.
- It requires significant resources for generating opcode-based rules and for loading goodware files.
- The rule generation process is slow.
- The creation of super rules may cause redundancy and duplication of rules.
- All dependencies and built-in databases have to be installed in order for the tool to work successfully.

***YaraGenerator* Tool** This python-based tool uses string prioritization logic and code refactoring to generate *Yara rules* with a completely different signature for different file types, such as *EXEs*, *PDFs*, and *Emails*.

The generated *Yara rules* contain only strings (opcodes are not supported) extracted from malware samples that do not match with the provided database of strings from blacklisted files. In particular 30,000 blacklisted strings are contained in such database, arranged based on the different file formats.

The produced *Yara rules* contain a large number of strings which are selected randomly. In fact, no score computation takes place in order to weight the different strings [38].

1. *YaraGenerator* tool advantages :

- It can generate specialised rules for specific file formats.

- It supports the use of PE (portable executable) modules, which are used to interpret Windows operating system executables such as *DLL* and *COM* files.
- It reduces the false positive rate by checking all strings against databases of blacklisted files.
- It is deployed as a simple and easy-to-use python script that can be run through a command-line interface.

2. *YaraGenerator* tool *disadvantages*:

- It requires post-processing of the generated rules for increasing their effectiveness.
- It generates rules based on a random selection of features (strings). This implies that the most appropriate strings may not be selected in many cases, thus making the produced rules less effective on average.
- It does not support the use of opcodes.
- It was developed as a work-in-progress project and has not been updated for a while now.

Yabin Tool This is another *python-based* tool, developed by the *Alien Vault Open Threat Exchange (OTX)* community, for the automatic generation of *Yara rules*.

In this case *Yara rules* are created by finding rare functions in specific malware samples or families. Functions are recognised by checking specific bytes sequences called *function prologues*, which define the start of the code of a function. For example, the byte sequence '*55 8B EC*' usually specifies the start of a function in programs compiled by Microsoft Visual Studio.

The generated *Yara rules* include those strings common to malware samples that don't match with the provided whitelist of commonly used library functions. Such whitelist was obtained from 100 Gb of non-malicious software in order to exclude common library functions.

The *Yara rules* produced contain a list of hexadecimal strings to be compared against suspicious files looking for similarities in their byte-sequences [38].

1. *Yabin* tool *advantages*:

- It can be used to cluster malware samples based on the reuse of their code.
- The list of patterns to search for can be extended during the rule post-processing phase.
- With the purpose of excluding commonly used library functions in the produced rules, a large whitelist obtained from numerous non-malicious executable files is provided with the tool.
- It is deployed as a simple and easy-to-use python script that can be run through a command-line interface.

2. *Yabin* tool *disadvantages*:

- It requires post-processing of the generated rules for increasing their effectiveness.
- Some specific file types/formats may not be supported.
- The created rules contain only function prologues. No other string types are used.
- Since it relies on function prologues, it works only with unpacked executables.
- It is not designed to work on *.NET* executables, *Java* files and *Microsoft* documents.
- It was mainly developed for research and testing purpose, not for production use.

Auto Yara Tool Compared to the previously mentioned tools like *YarGen*, which rely on a number of heuristics and string features, *Auto Yara* tool makes larger rules using the redundancy and conjunction of components to achieve extremely low false-positive rates [39].

The two primary concerns of the *Auto Yara* authors while designing this tool were:

1. Yara rules that generate a lot of false positives could slow the investigation
2. Malware analysts often have few samples (≤ 10) when creating a Yara rule

Auto Yara authors thus developed a workflow composed of two steps: the first step leveraged recent works in finding frequent larger n-grams, for $n \leq 1024$, to find several candidate byte strings that could become features. In the second step a bi-clustering method, which consists of simultaneously clustering the rows and columns of an input data matrix, is used on those strings to construct the output rules. Most bi-clustering algorithms require the specific number of bi-clusters to be known in advance, and enforce no overlaps between bi-clusters. The *Auto Yara* authors exploited an already existing bi-clustering algorithm extending it to work when the number of bi-clusters is not known *a priori* (the number of bi-clusters gets determined automatically) and to allow overlapping bi-clusters, discarding rows and columns that do not fit in any bi-cluster [39].

Auto Yara uses *bi-clustering* because it allows to easily produce complex and effective logic rules that enable the creation of signatures with low false positive rates.

To build a good Yara rule, in fact, one needs to know:

1. which features should be used at all
2. which features should be combined into '*and*' statements (which reduce the False Positive Rate), and which should be placed into '*or*' statements (which increase the True Positive Rate)

Bi-clustering provides a simple approach to do this jointly over the features, rather than considering the features one at a time. In particular, the features within a bi-cluster are combined into an '*and*' statement since they co-occur; moreover the '*and*' statements from multiple bi-clusters are placed into an '*or*' statement resulting in a "disjunction of conjunctions" rule formulation.

1. *Auto Yara* tool advantages:

- It is fast, allowing it to be deployed even on low-resource equipment (like remote networks).
- It was designed with the intent of producing *Yara rules* with low false positive rates.
- It was designed to be able to generate *Yara rules* from as few as ≤ 10 available samples.

2. *Auto Yara* tool disadvantages:

- It requires post-processing of the generated rules for increasing their effectiveness.
- It is a very recent tool, mainly developed for research purposes and not for production use.

Consider if to add something about "Manually generated Yara Rules + Open Source Yara Rules databases".

3.3 Semantic Based Detection

Semantic-based malware detection aims at identifying malware by deducing the analysed code logic and comparing it to a database of already known malicious logic patterns. This technique, differently from signature-based detection which looks at the code syntactic properties, tracks the semantics of the program code instructions. This implies that *semantic-based* detection approaches are capable of overcoming obfuscation attempts and even detecting unknown malware variants [5].

3.4 Behavioural Based Detection

Behavioural-based malware detection is based on the use of behavioural patterns for the identification of malicious software. This is done by dynamically analysing malware samples and extracting specific system/application behaviours and activities in order to form a '*behavioural signature*' of a malware strain. New samples are then analysed in the same way and identified as malware if their behavioural pattern is similar to the *behavioural signature* of a known malware [5].

Behavioural-based detection is for the most part immune to obfuscation attempt. However, being based on the time consuming dynamic analysis and on the challenging task of determining the unsafe activities and behaviours to consider within the environment, its applicability is limited.

3.5 Heuristics-based Detection

As opposed to traditional *signature-base* detection methods which identify malware by looking in the code for specific bytes/strings, *heuristic-based* detection uses rules and/or algorithms to search for commands or instructions not commonly found in harmless applications, thus indicating possible malicious intents [42].

Heuristic-based anti-malware tools may exploit different scanning techniques such as:

- *File analysis (static heuristic analysis)*: the suspicious program is disassembled and its source program is examined looking for known malware patterns (stored in a heuristic database). If the percentage of matched code exceeds a predefined threshold then the code is marked as probably infected [43].
- *File emulation (dynamic heuristic analysis)*: in this approach, the suspicious piece of code is examined in a virtual machine (or sandbox) looking for suspicious operations such as attempts at executing other executables, at changing the Master Boot Record, at concealing themselves etc. that are uncommon in benign programs.
- *Genetic signature detection*: this technique is designed to spot different malware variations within the same family using previous malware definitions [44].

Heuristic analysis is a promising technique for the detection of unknown malware, particularly for encrypted and polymorphic variants [22].

Nowadays *heuristic* analysis can be found in most mainstream antivirus solutions in the market, combined with signature-based scanners in order to improve detection rate while reducing false alarms [5].

3.6 Machine Learning

In recent years, the rapid proliferation and increased sophistication of malicious software, coupled with the rising popularity of machine learning techniques in many fields, led to the adoption of more general ML-based approaches to malware detection on top of the use of manually generated signatures [3].

In particular, the application of machine learning for Information Security (ML-Sec) methods to perform malware detection generally consists in training a highly parametrized ML classifier to reliably (as much as possible) predict a binary label (malicious or benign) using features extracted from sample files. In order to do this the classifier parameters are numerically optimized to learn general concepts of *malware* and *benignware*, by minimizing misclassification as measured by some loss criterion, which measures the deviation of predictions from their actual ground truth. This is based on the assumption that, if the samples are well labelled and malware/benignware samples in the training set are similar enough to those seen at test/deployment, the learned detection function should work well on unseen samples [2].

Most static ML-Sec classifiers work on learned embeddings over portions of files (e.g. headers), learned embeddings over full files, or most commonly, on pre-engineered numerical *feature vectors* designed to summarize the content from each file. Learned embeddings generally are the result of convolutional architectures which do not presume a fixed file structure. However, the process of embedding features directly from inputs is expensive, and does not scale gracefully. Moreover, generic bytes do not present structural localities/hierarchies typical of images and text inputs that can be exploited by convolutional filters. Pre-engineered feature vector representations, on the other hand, quickly distil content useful for classification from each file [45]. There are various possible ways to statically craft feature vectors, for example:

- tracking per-byte statistics over sliding windows
- byte histograms
- n-gram histograms
- treating bytes as pixel values in an image
- opcode and function call graph statistics
- symbol statistics
- hashed/numerical metadata values
- hashes of delimited tokens
- etc.

The process of transforming a sample file to its numerical feature representation is called *feature extraction* and consists of some numerical transformation that preserves aggregate and fine-grained information throughout each sample [2].

ML methods generally require many high quality samples in order to train effective models. When creating datasets for these models labels are often collected from vendor aggregation feeds, which combine detection results from various vendors for each malware sample. This can be done, for example, by using a 1-/5+ criterion or by using statistical estimation methods. The 1-/5+ criterion works as follows: if a file has one or fewer vendors reporting it as malicious, the file is labelled as 'benign'; on the other hand, if a sample has five or more vendors reporting it as malicious, the file is labelled as 'malicious'. Moreover, it is common practice to introduce a time lag to let vendors update their models to account for new malware samples. When deployed, classifiers are periodically re-trained on new data/labels to reflect the current malware trends [45].

The advantage of machine learning techniques with respect to signature engines, where the aim is to reactively blacklist/whitelist samples that hard-match manually-defined patterns (signatures), is that, by being more general, they are able to detect not only known malwares but also novel malware strains/variants, providing some degree of proactive detection [36].

Commercial anti-malware solutions/engines have nowadays integrated ML on top of standard detection methods (without replacing them) with the aim of enhancing the detection results, especially for second generation malwares and novel malware strains. Popular ML techniques employed by such tools are, for example, deep neural networks (DNN), boosted decision tree ensembles, Naïve Bayes models, Data Mining approaches and Hidden Markov Models. Moreover, multiple vendors in the IT security industry nowadays have dedicated ML-Sec teams [22].

In the following I will focus on the following recent static ML-based malware detection methods:

- ALOHA: Auxiliary Loss Optimization for Hypothesis Augmentation
- Automatic Malware Description via Attribute Tagging and Similarity Embedding
- Learning from Context: Exploiting and Interpreting File Path Information for Better Malware Detection

3.6.1 ALOHA: Auxiliary Loss Optimization for Hypothesis Augmentation

In a recent work called **ALOHA: Auxiliary Loss Optimization for Hypothesis Augmentation**, [2], Rudd et al. observed that, although ML-based malware detection is frequently framed as a binary classification task (using a simple binary cross-entropy loss function), there are often a number of other sources of contextual metadata for each input sample available at training time, beyond just aggregate malicious/benign labels. Such metadata might include malicious/benign labels from multiple sources (e.g. from various security vendors), malware family information, temporal information, counts of affected endpoints, and associated tags.

However, this metadata is, in many cases, not available at deployment time thus making it difficult to include it as input features. Therefore, the authors proposed exploiting the metadata from threat intelligence feeds as auxiliary targets in the multi-target learning approach. Simultaneously optimizing classifier parameters for multiple targets (labels) while training a model may, in fact, have a regularizing effect leading to better generalization, particularly if the auxiliary targets are related to the main target of interest.

In practice, in their work [2] the authors fit a deep neural network with multiple additional targets derived from metadata in a threat intelligence feed for Portable Executable (PE) malware and benignware. The additional losses include a multi-source malicious/benign loss, a count loss on multi-source detections, and a semantic malware attribute tag loss. The authors ultimately stated that this approach yielded a considerable improvement in performance on the main detection task.

Inner Workings

The model presented in [2] (fig. 3.1) is composed of a base feed-forward neural network consisting of 5 blocks, each composed of Dropout, a dense layer, batch normalization, and an exponential linear unit (ELU) activation, with 1024, 768, 512, 512, and 512 hidden units respectively. This base topology applies the function $f(\cdot)$ to the input feature vector \mathbf{x} (of size 1024 in this case) to produce an intermediate 512 dimensional representation of the input file $\mathbf{h} = f(\mathbf{x})$. An additional block for each output of the model, consisting of one or more dense layers and activation functions, is then appended on top of the base net. This composition of the base topology and the target-specific 'heads' is denoted as $f_{target}(\mathbf{x})$.

The output for the main malware/benign prediction task - $f_{mal}(\mathbf{x})$ - is always present (it constitutes the baseline model) and consists of a single dense layer followed by a sigmoid activation function on top of the base shared network. On top of that one or more auxiliary outputs are added with similar structure as described above: one fully connected layer (two for the *tag* prediction task) with a task-specific activation function. Finally, multi-task losses are produced by computing the sum, across all tasks, of the per-task loss multiplied by a task-specific weight (1.0 for the malware/benign task and 0.1 for all other tasks) [2].

Malware Loss The task of predicting if a given binary file, represented by its features $\mathbf{x}^{(i)}$, is malicious or benign is optimized by minimizing the binary cross-entropy loss between the malware/benign output of the network $\hat{y}^{(i)} = f_{mal}(\mathbf{x}^{(i)})$ and the malicious label $y^{(i)}$. This results in the following loss for a dataset with M samples:

$$\begin{aligned} L_{mal}(X, Y) &= \frac{1}{M} \sum_{i=1}^M l_{mal}(f_{mal}(\mathbf{x}^{(i)}), y^{(i)}) \\ &= -\frac{1}{M} \sum_{i=1}^M y^{(i)} \log(\hat{y}^{(i)}) + (1 - y^{(i)}) \log(1 - \hat{y}^{(i)}). \end{aligned} \tag{3.1}$$

A "1-/5+" criterion was used for labelling a given file as malicious or benign [2].

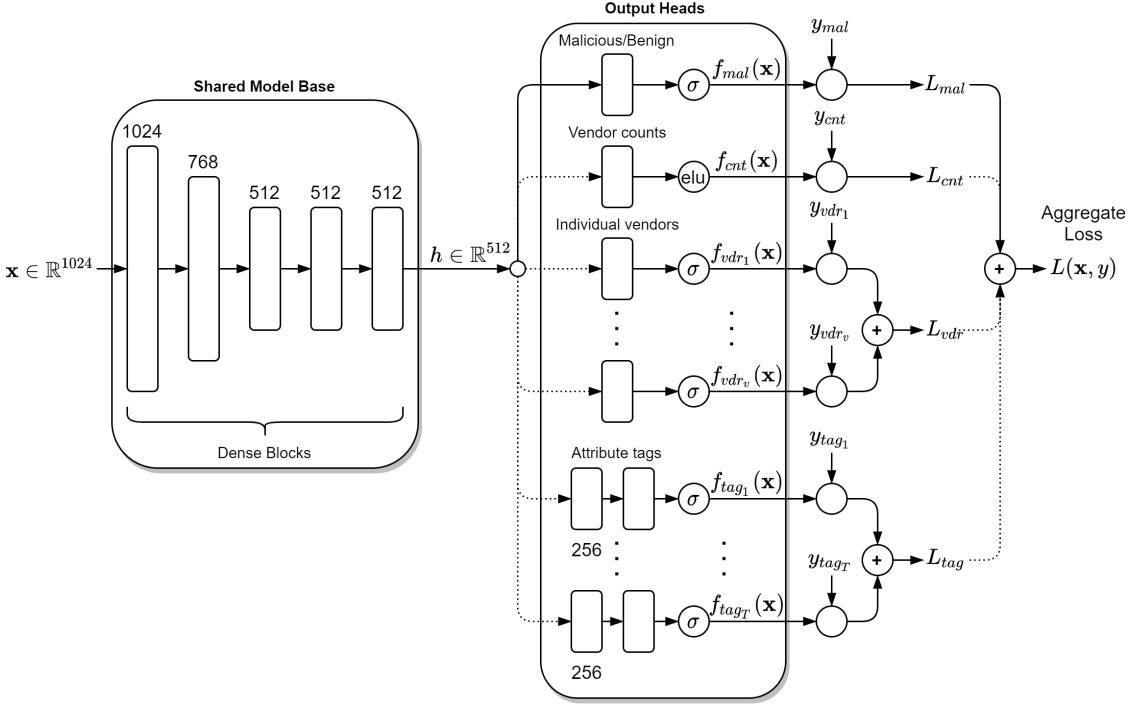


Figure 3.1: ALOHA model architecture

Vendor Count Loss The first additional target the authors of [2] tried using is the count of ‘malicious’ reports for a given sample from the vendor aggregation service. This is based on the assumption that the more a sample gets reported by vendors, the more likely it is to be malicious. To model count data, the authors used the Poisson noise model parametrized by the parameter μ , where μ is the mean and variance of the Poisson distribution. The probability of an observation of y counts is therefore:

$$P(y|\mu) = \mu^y e^{-\mu} / y!. \quad (3.2)$$

The model is then trained to estimate μ for each sample $\mathbf{x}^{(i)}$ such that the likelihood of $y^{(i)}|\mu^{(i)}$ is maximized (or, equivalently, the negative log-likelihood is minimized). The output of the model vendor count head, for sample i , is denoted as $f_{cnt}(\mathbf{x}^{(i)})$. Thereafter, an activation function $a(\cdot)$, which maps $f_{cnt}(\mathbf{x}^{(i)})$ to the non-negative real numbers, is used so the final approximation for parameter μ is: $\mu^{(i)} = a(f_{cnt}(\mathbf{x}^{(i)}))$. In particular, the activation function used in [2] is an exponential linear unit activation (ELU).

Letting $y^{(i)}$ denote the actual number of vendors that deemed sample $\mathbf{x}^{(i)}$ malicious, the corresponding negative log-likelihood loss over the dataset is:

$$\begin{aligned} L_p(X, Y) &= \frac{1}{M} \sum_{i=1}^M l_p(a(f_{cnt}(\mathbf{x}^{(i)})), y^{(i)}) \\ &= \frac{1}{M} \sum_{i=1}^M \mu^{(i)} - y^{(i)} \log(\mu^{(i)}) + \log(y^{(i)}!), \end{aligned} \quad (3.3)$$

which is referred to in [2] as the *Poisson* or *vendor count* loss. In practice, the term $\log(y^{(i)}!)$ can be ignored when minimizing this function loss since it does not depend on the network parameters.

Per-Vendor Malware Loss The authors of [2] identified a subset $\mathcal{V} = \{v_1, \dots, v_V\}$ of 9 vendors that each produced a result for (nearly) every sample in their data. Each vendor result was added as additional target by adding an extra fully connected layer per vendor followed by a sigmoid activation function to the end of the shared baseline architecture. A binary cross-entropy loss

per vendor was then employed during training. The aggregate *vendors* loss L_{vdr} for the $V = 9$ selected vendors is simply the sum of the individual vendor losses:

$$\begin{aligned} L_{vdr}(X, Y) &= \frac{1}{M} \sum_{i=1}^M \sum_{j=1}^V l_{vdr}(f_{vdr_j}(\mathbf{x}^{(i)}), y_{v_j}^{(i)}) \\ &= -\frac{1}{M} \sum_{i=1}^M \sum_{j=1}^V y_{v_j}^{(i)} \log(\hat{y}_{v_j}^{(i)}) + (1 - y_{v_j}^{(i)}) \log(1 - \hat{y}_{v_j}^{(i)}), \end{aligned} \quad (3.4)$$

where l_{vdr} is the per-sample binary cross-entropy function and $f_{vdr_j}(\mathbf{x}^{(i)}) = \hat{y}_{v_j}^{(i)}$ is the output of the network that is trained to predict the label $y_{v_j}^{(i)}$ assigned by vendor j to input sample $\mathbf{x}^{(i)}$ [2].

Malicious Tags Loss Further additional targets were provided in the form of malicious tags. This was done in an attempt to exploit information contained in family detection names provided by different vendors. In particular, the tags used as auxiliary targets in [2] are: *flooder*, *downloader*, *dropper*, *ransomware*, *crypto-miner*, *worm*, *adware*, *spyware*, *packed*, *file-infecto* and *installer*.

These tags were created by parsing individual vendor detection names, using a set of 10 vendors which provided high quality detection names. After having extracted the most common tokens, the tokens not related to well-known malware family names were filtered out. Finally, a mapping between tokens and tags was created based on the authors experience.

The tag prediction task was then defined as a multi-label binary classification, since zero or more tags from the set of possible tags $\mathcal{T} = \{t_1, \dots, t_T\}$ can be present at the same time for a given sample. This was implemented as a *multi-headed* architecture: two additional layers per tag were added to the end of the shared baseline architecture, a fully connected layer of size 512-to-256, followed by a fully connected layer of size 256-to-1, followed by a sigmoid activation function. Each of the possible $T = 11$ tags has its own loss term computed with binary cross-entropy. Finally, the aggregate tag loss is the sum of the individual tag losses [2]. For the dataset with M samples the loss is:

$$\begin{aligned} L_{tag}(X, Y) &= \frac{1}{M} \sum_{i=1}^M \sum_{j=1}^T l_{tag}(f_{tag_j}(\mathbf{x}^{(i)}), y_{t_j}^{(i)}) \\ &= -\frac{1}{M} \sum_{i=1}^M \sum_{j=1}^T y_{t_j}^{(i)} \log(\hat{y}_{t_j}^{(i)}) + (1 - y_{t_j}^{(i)}) \log(1 - \hat{y}_{t_j}^{(i)}), \end{aligned} \quad (3.5)$$

where $y_{t_j}^{(i)}$ indicates if sample i is annotated with tag j , and $\hat{y}_{t_j}^{(i)} = f_{tag_j}(\mathbf{x}^{(i)})$ is the prediction of the network for that value.

Final Loss Finally, in [2] each model used a loss weight of 0.1 on the aggregate malicious/benign loss and 0.1 on each auxiliary loss. Therefore, when adding K targets to the main loss, the final loss that gets back-propagated through the model becomes:

$$L(X, Y) = L_{mal}(X, Y) + 0.1 \sum_{k=1}^K L_k(X, Y). \quad (3.6)$$

3.6.2 Automatic Malware Description via Attribute Tagging and Similarity Embedding

As explained by Ducau et al. in [3], in order to counter and remediate a malware infection on a system/network it is of vital importance to understand the nature (malware family and variant) of the attack in progress. In fact, knowing the malicious capabilities associated with each suspicious file found on a system/network gives important clues to the system end user, administrator or

security operator that help define a remediation procedure, identify possible root causes, and evaluate the severity and potential consequences of the attack.

Machine learning detection methods, which have the potential to identify even new malware samples/variants as malicious, however, generally produce a simple binary malicious/benign label with no further information about the type of threat posed by malicious samples, which in turn does not allow the identification of relationships between malware samples. On the other hand, most commercial anti-malware solutions provide, when they alert about potentially harmful files detected in a machine, *detection names* coming from specific hand-written signatures created by reverse engineers to identify particular threats, that are theoretically useful for categorizing known malware variants. However, even these detection names are problematic. In fact, the increasing number of feature-rich malware and the fact that different vendors use differing malware naming conventions led to inconsistent and highly vendor-specific detection names. Moreover, it is now common to see detection names which only act as unique identifiers without providing actionable information about the type of potential harm the malicious sample is capable of doing.

The authors of [3] therefore proposed the use of SMART (Semantic Malware Attribute Relevance Tagging) tags, which are human interpretable, high level descriptions of the capabilities of a given malware sample, to approach malicious software description. These SMART tags, which were derived by leveraging the underlying knowledge encoded in detection names from different anti-malware vendors in the industry, are non-exclusive, meaning that one malware family can be associated with multiple tags and a given tag can be associated with multiple malware families. In [3], the authors defined a set of malicious tags \mathcal{T} , with $|\mathcal{T}| = 11$ different tags (or descriptive dimensions) of interest that they then used to describe malicious PE files: *adware*, *crypto-miner*, *downloader*, *dropper*, *file-infector*, *flooder*, *installer*, *packed*, *ransomware*, *spyware* and *worm*.

These tags were then used to train a Joint Embedding neural network to learn a low dimensional Euclidean representation (embedding) space in which malware samples with similar characteristics are close to each other, having access only to the files' static binary representations; this in turn is used at test/deploy time for automatically predicting tags for new (unseen) files in real time.

The representation of a malware sample in the embedding space can also be taught as an implicit 'signature' describing its capabilities.

Interestingly, in [3] the authors state that the proposed representation of PE files in joint embedding space, along with the distance function defined in this space, allows to measure not only similarities between files' and tags' embeddings, but between files' representations as well.

In particular, because of how the joint embedding model is designed and trained, two pieces of malware that are close to each other in latent (embedding) space should theoretically exhibit similar malicious capabilities. This makes it possible to measure how close two pieces of malware are in terms of their capabilities by using their latent representations. Moreover, the joint embedding representation, given that its size (32 dimensions) is 32 times smaller than the original file feature representation (1024 dimensions), is particularly efficient for storing, indexing and querying large malware databases [3].

Malware Family Categorization Given the extremely high number of malware samples which are nowadays received by security vendors' labs, in the order of millions per month, it has become unfeasible to manually and consistently group each malicious file into a well defined hierarchy of families.

The detection names typically used by modern security vendors can be divided into four categories, which contain varying amounts of information about the specific malware families [3].

- *Traditional family based*: Detection names are associated to unique and distinctive attributes of malware and its variants. More effort in analysing malware inner-workings is required. Detection names of this type are often of high-quality and have more consistency across vendors.

- *Technique based*: Detection names group together malware that may come from different origins and/or have multiple authors but share a common method or technique. The difference in detection methods employed by different vendors often results in less consistency in detection names across vendors.
- *Method based*: Detection names denote the specific detection technology used to detect the malware sample. Some detection names can be that of a patented technology, project, or internal code name specific to the AV vendor, indicating the use of heuristics, ML, or real-time detection technologies. In this case the detection name does not refer to a particular malware family, but to the specific method that was used to detect the sample.
- *Kit based*: Detection names are associated to kits/tools used to generate the specific malware sample. These kits can be used both offensively by penetration testing teams and by malware authors and are therefore often referred to as grey hat tools. Detection names in this category tend to identify methods used by the kit or tool to obfuscate or hide the payload of a specific malware sample, rather than to describe its origin or functionality.

Multi-Label Classification The authors of [3] applied multi-label classification in order to perform semantic attribute tagging. The most trivial way of implementing a multi-label classification model is by learning one classifier per-label. This approach, however, is far from being efficient since the single per-tag classifiers are independently optimized. In this scenario, it is more common to use a single classifier with multiple outputs (multi-label learning) and multiple target losses which are combined and jointly optimized during training (multi-objective loss). This approach yields a more compact representation while also improving classification performance with respect to using independent classifiers. In [3], the authors used a multi-label deep neural network architecture as baseline architecture.

An alternative approach to multi-label classification is to learn a compact shared vector (embedding) space representation to which map both input samples and labels - a joint embedding - where similar content across modalities are projected into similar vectors in the same low dimensional space [3]. Then, at test/deploy time, in order to determine likely labels a similarity comparison between vectors belonging to this learned latent space is performed, e.g. via inner product. In [3], the authors used a joint embedding model that maps malware tags and executable files into the same low dimensional Euclidean joint-embedding space for the malware description problem.

Tag Distillation from Detection Names Ducau et al. [3] relied on semi-automatic strategies, even if they are noisier than manual labelling, because they allowed them to label millions of files that can be then used to train a classifier. In particular, they designed a labelling function which annotates PE files using the previously mentioned set of tags \mathcal{T} by combining information extracted from the detection names got from ten reputable anti-malware vendors.

The labelling process consisted of the following three main stages:

1. *token extraction*
2. *token-to-tag mapping*
3. *token relationship mining*

The token extraction phase consisted of normalizing and parsing the multiple detection names and converting them in sets of sub-strings. In a similar way to what the malware labelling tool AVClass, which was proposed by Marcos et al. in [46], does, the token-to-tag mapping stage uses rules created from expert knowledge by a group of malware analysts, that associate relevant tokens with the set of tags of interest. Finally, this mapping is extended by mining statistical relationships between tokens to improve tagging stability and coverage.

Tags Prediction - Inner Workings In [3], in order to predict, in a multi-label classification manner, zero or more tags per sample from the set of T possible tags $\mathcal{T} = \{t_1, t_2, \dots, t_T\}$, the authors proposed two different neural network architectures, represented in 3.2 and 3.3, which they referred to as *Multi-Head* (top) and *Joint Embedding* (bottom).

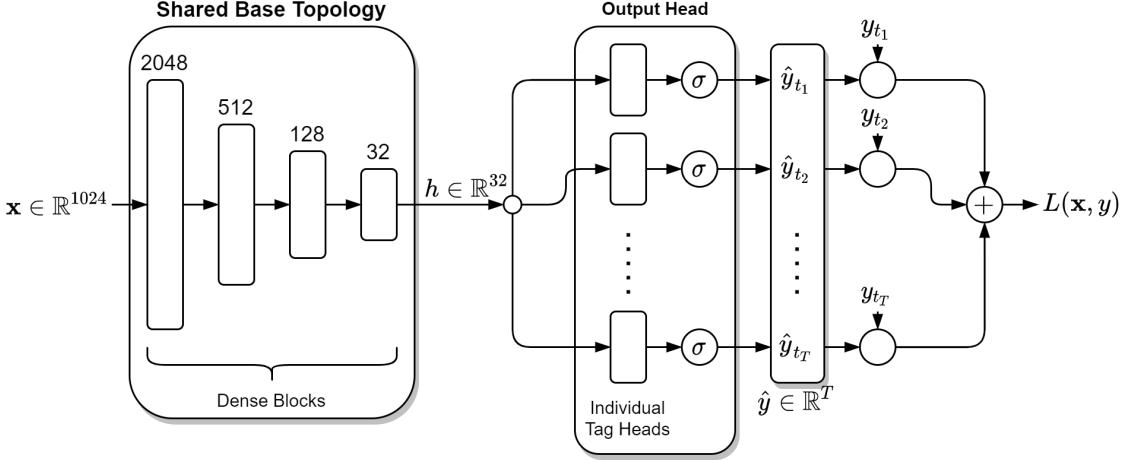


Figure 3.2: Multi Head model architecture

In particular, the *Multi-Head* (fig. 3.2) consisted of a base topology common to the prediction of all tags, and one output (or "head") per tag. The shared base topology can be thought of a feature extraction (or embedding) network that transforms the input features \mathbf{x} into low dimensional hidden vector \mathbf{h} and consisted of an input feed-forward layer of output size 2048, followed by a batch normalization layer, an ELU non-linearity and three blocks, each composed by dropout, a linear layer, batch normalization and ELU of output sizes 512, 128, and 32 respectively. Each head, on the other hand, is a binary classifier that predicts the presence or absence of each tag and it consists of a linear layer (the same for each head) composed of the same type of basic blocks as are in the main base architecture, but with output size 11 (the number of tags being predicted) and a sigmoid non-linearity instead of the ELU to compute the predicted probability for each label. Binary cross-entropy loss is computed as the output of each head and then added together to form the final loss.

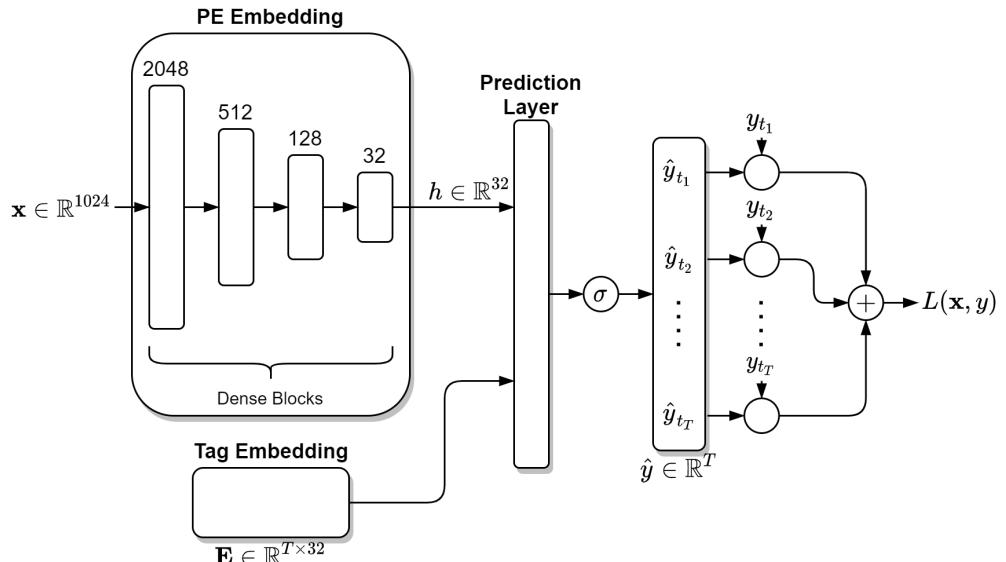


Figure 3.3: Joint Embedding model architecture

The *Joint Embedding* model (fig. 3.3), instead, mapped both the labels (malware tags) and the binary file features \mathbf{x} to vectors in a joint Euclidean latent-space in such a way that, for a given

similarity function, the transformations of semantically similar labels are close to each other, and the embedding of a binary file should be close to that of its associated labels in the same space. In practice, the *Joint Embedding* model consisted on a PE *embedding* network, a *tag embedding* matrix \mathbf{E} , and a *prediction layer*.

The PE embedding network, using the same base topology as the Multi-Head model, learns a non-linear function $\phi_\theta(\cdot)$, with parameters θ that maps the input binary representation of the PE executable file $\mathbf{x} \in \mathbb{R}^d$ into a vector $\mathbf{h} \in \mathbb{R}^D$ in low dimensional Euclidean space (with $D = 32$),

$$\phi_\theta(\mathbf{x}) : \mathbb{R}^d \rightarrow \mathbb{R}^D. \quad (3.7)$$

The tag embedding matrix $\mathbf{E} \in \mathbb{R}^{T \times D}$ of learnable parameters learns a mapping from a tag $t_n \in \mathcal{T} = \{t_1, \dots, t_T\}$, to a distributed representation $\mathbf{e} \in \mathbb{R}^D$ in the joint embedding space (with $D = 32$).

$$\phi_E(t) : \{t_1, \dots, t_T\} \rightarrow \mathbb{R}^D. \quad (3.8)$$

In practice, the embedding vector for the tag t_n is simply the n -th row of the tag embedding matrix, i.e. $\phi_E(t_n) = \mathbf{E}_n$.

Finally, the prediction layer compares both tag and sample embeddings (\mathbf{e} and \mathbf{h} respectively) and produces a similarity score. This is later run through a sigmoid non-linearity to estimate the probability that sample \mathbf{x} is associated with tag t for each $t \in \mathcal{T}$. In the final model implementation, the similarity score is the dot product between the embedding vectors. The output of the network $f_n(\mathbf{x}|\theta, \mathbf{E})$ then becomes,

$$\begin{aligned} \hat{y}_n &= f_n(\mathbf{x}|\theta, \mathbf{E}) = \sigma(\langle \phi_E(n), \phi_\theta(\mathbf{x}) \rangle) \\ &= \sigma(\langle \mathbf{E}_n, \mathbf{h} \rangle), \end{aligned} \quad (3.9)$$

where σ is the sigmoid activation function, and \hat{y}_n is the probability estimated by the model of tag t_n being a descriptor for \mathbf{x} .

Furthermore, the authors of [3] constrained the embedding vectors for the tags such that:

$$\|\mathbf{E}_n\|_2 \leq C, n = 1, \dots, T, \quad (3.10)$$

which has a regularizing effect and they fixed the value of C to 1.

During training, the parameters of both embedding functions $\phi_\theta(\cdot)$ and $\phi_E(\cdot)$ are jointly optimized to minimize the binary cross-entropy loss for the prediction of each tag via back-propagation and stochastic gradient descent. The loss function to minimize for a mini-batch of M samples is:

$$\begin{aligned} \mathcal{L} &= -\frac{1}{M} \sum_{i=1}^M \sum_{n=1}^T f_n(\mathbf{x}^{(i)}|\theta, \mathbf{E}) \log(t_n^{(i)}) + (1 - f_n(\mathbf{x}^{(i)}|\theta, \mathbf{E})) \log(1 - t_n^{(i)}) \\ &= -\frac{1}{M} \sum_{i=1}^M \sum_{n=1}^T \hat{y}_n^{(i)} \log(t_n^{(i)}) + (1 - \hat{y}_n^{(i)}) \log(1 - t_n^{(i)}) \end{aligned} \quad (3.11)$$

where $t_n^{(i)} = 1$ if sample i is labelled with tag t_n or zero otherwise, and $\hat{y}_n^{(i)}$ is the probability predicted by the network of that tag being associated with the i -th sample.

In order to get a vector of tag similarities for a given sample \mathbf{x} with PE embedding vector \mathbf{h} , the matrix of tag embeddings $\mathbf{E} \in \mathbb{R}^{T \times D}$ is multiplied by $\mathbf{h} \in \mathbb{R}^D$; the output is then scaled to obtain a prediction vector $\hat{\mathbf{y}} = \sigma(\mathbf{E} \cdot \mathbf{h}) \in \mathbb{R}^T$, where σ is the element-wise sigmoid function for transforming the similarity values into a valid probability. Each element in $\hat{\mathbf{y}}$ is then the predicted probability for each tag.

Evaluation of Tagging Algorithms The performance evaluation of tagging algorithms can be done along two orthogonal dimensions: *per-tag* or a *per-sample*. The former aims at quantifying the tagging algorithm performance at identifying each tag, while the latter focuses on evaluating the tagging algorithm performance for a given sample, across all tags.

One suitable way to evaluate the *per-tag* performance of a model is by measuring the *Area Under the Receiver Operating Characteristic* curve (*AUC-ROC*, or simply *AUC*) for each of the tags being predicted. A *ROC* curve is created by plotting the *True Positive Rate (TPR)* against the *False Positive Rate (FPR)*. Furthermore, given the binary True/False nature of the target value for the n -th tag of a given sample ($t_n \in \{0, 1\}$), the typical binary classification evaluation metrics such as 'Accuracy' (eq. 3.12), 'Precision' (eq. 3.13), 'Recall' (eq. 3.14), and 'F-score' (eq. 3.15) can also be computed. However, in order to compute these metrics, the output probability prediction needs to be binarized first. This can be done simply choosing a threshold independently for each tag so that the *FPR* on the validation set is 0.01 and then using the resulting 0/1 predictions.

$$\text{accuracy} = \frac{TP + TN}{TP + FP + TN + FN} \quad (3.12)$$

$$\text{precision} = \frac{TP}{TP + FP} \quad (3.13)$$

$$\text{recall} = \frac{TP}{TP + FN} \quad (3.14)$$

$$f_\beta \text{score} = (1 + \beta^2) \cdot \frac{\text{precision} \times \text{recall}}{\beta^2 \cdot \text{precision} + \text{recall}} \quad (3.15)$$

where TP , TN , FP and FN are the number of *True Positives*, *True Negatives*, *False Positives* and *False Negatives* respectively.

On the other hand, the *Mean Jaccard* similarity (eq. 3.16) and *Mean per-sample Accuracy* (eq. 3.17) metrics can be used to assess the per-sample performance of a tagging algorithm. In particular, the Jaccard similarity (or index) can be used as a figure of how similar the set of tags associated with a specific sample is with respect to the set of tags predicted for the same sample after binarizing the predictions. The per-sample accuracy is instead defined as the percentage of samples for which the target vector is equal to the prediction vector, i.e. all tags correctly predicted. For an evaluation dataset with M samples the equations of the per-sample evaluation metrics are:

$$\text{Mean Jaccard similarity} = J(T^{(i)}, \hat{T}^{(i)}) = \frac{1}{M} \sum_{i=1}^M J(T^{(i)}, \hat{T}^{(i)}) = \frac{1}{M} \sum_{i=1}^M \frac{T^{(i)} \cap \hat{T}^{(i)}}{T^{(i)} \cup \hat{T}^{(i)}} \quad (3.16)$$

$$\text{Mean per-sample accuracy} = \frac{1}{M} \sum_{i=1}^M \mathbb{I}(\mathbf{y}^{(i)} = \hat{\mathbf{y}}^{(i)}) \quad (3.17)$$

where $T^{(i)}$ is the set of tags associated with sample i , $\hat{T}^{(i)}$ is the set of tags predicted for the same sample after binarizing the predictions, $\mathbf{y} \in \{0, 1\}^T$ is the binary target vector for a PE file (where y_n indicates whether the n -th tag applies to the file), $\hat{\mathbf{y}}$ is the binarized prediction vector from a given tagging model and \mathbb{I} is the indicator function which is 1 if the condition in the argument is true, and 0 otherwise.

3.6.3 Learning from Context: Exploiting and Interpreting File Path Information for Better Malware Detection

Recent static portable executable (PE) malware detection techniques, including [2], typically employ ML-Sec classifiers with a single numerical feature vector derived from each file as input, having as output one or more target labels/tasks during training. However, as noted by Kyadige et. al. in [45], there is still much unused orthogonal information that could be exploited regarding the samples files, namely the file paths. The authors in [45] thus proposed utilizing the paths of

the PE files as static source of contextual information as auxiliary data to the classifier in order to augment static ML detectors. File paths, which are already commonly used by malware analysts to correct and investigate detection errors, are available statically with very little overhead, and can seamlessly be integrated into a multi-view static ML detector.

File paths are not inherently malicious or benign; however, given that many malware strains use specifically crafted file paths to perpetrate their malicious intents (a file path may in fact be chosen to increase the odds of the malware being executed, to avoid disk scans, or to hide from the user's view), they provide much instrumental information that can be used to enhance the overall detection.

In [45], the authors thus proposed the use of a *multi-view* neural network which combines, in input, information about the PE file content, via feature vectors, with information about how likely it is to see such file in a specific location, through file paths, and outputs a detection score.

To compare their results, they actually focused their experiments on three models:

- A baseline file-content-only PE model, which takes only the PE features as input and outputs a malware confidence score.
- Another baseline file-path-only FP model, which takes only the file's paths as input and outputs a malware confidence score.
- Their proposed multi-view PE file-content + contextual file-path (PE + FP) model, which takes both the PE file content features and file paths as inputs, and also outputs a malware confidence score.

The experiments were conducted on a dataset of files and file paths collected from actual scans on customer endpoints from a large anti-malware vendor.

The authors of [45] finally stated that their proposed multi multi-view classifier trained on both file content and the contextual file paths yielded statistically better results across the ROC curve and particularly in low false positive rate (FPR) regions.

Feature Engineering In order to be able to use file paths in a feed-forward neural network along with PE file content feature vectors, the file paths, which are strings of variable length, needed to be converted into numerical vectors of fixed size. To do this the authors of [45] created a lookup table keyed on each character with a numeric value (between 0 and the character set size) representing each character. Moreover, the conversion required the file paths to be trimmed to a fixed size, therefore the authors considered just the first 100 characters of each file path.

As features for the content of the PE files, they used floating point 1024-dimensional feature vectors consisting of four distinct feature types, similar to [47]:

1. A 256-dimensional (16x16) 2D histogram of windowed entropy values per byte with a window size of 1024.
2. A 256-dimensional (16x16) 2D logarithmically scaled string length/hash histogram.
3. A 256-dimensional bin of hashes of metadata from the PE header, including PE metadata, including imports, exports, etc.
4. A 256-dimensional (16x16) byte standard deviation/entropy histogram.

In total, they represented each sample as two feature vectors: A PE-content feature vector of 1024 dimensions and a contextual file-path feature vector of 100 dimensions [45].

Inner Workings The model proposed in [45] (fig. 3.4) has two inputs, the 1024 element PE-content feature vector, \mathbf{x}_{PE} , and the 100 element file-path integer vector, \mathbf{x}_{FP} . The two distinct inputs are fed to two different base sub-networks, each composed by a series of layers with input-specific parameters: θ_{PE} for the PE content part and θ_{FP} parameters for file-path part. The two sets of parameters (θ_{PE} and θ_{FP}) are jointly optimized during training. The outputs of these base sub-networks are then concatenated and passed through a series of final hidden layers - a joint output sub-network with parameters θ_O terminating with a final dense layer followed by a sigmoid activation function. The final sigmoid activation is used to have as output a detection score between 0 (benign) and 1 (malicious). However, the threshold for determining if a sample is malicious or benign can be set anywhere along the (0.0, 1.0) range according to false positive rate (FPR) and detection rate (TPR) trade-offs for the application at hand - a reasonable threshold is typically at or below 10^{-3} FPR.

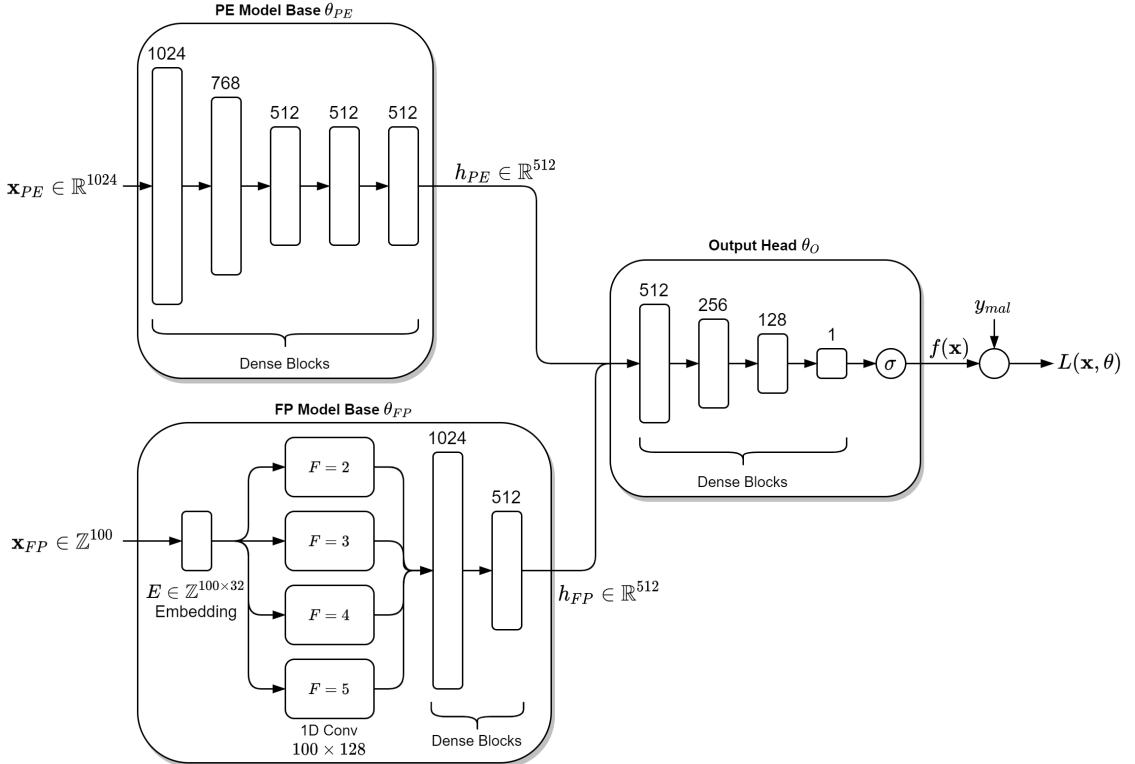


Figure 3.4: PE + FP model architecture

The PE base sub-network (with parameters θ_{PE}) passes its input \mathbf{x}_{PE} through a series of 5 blocks with sizes 1024, 768, 512, 512, and 512, each consisting of four layers: a Fully Connected layer, a Normalization layer, a Dropout layer with dropout probability of 0.05, and a Rectified Linear Unit (ReLU) activation function.

The FP base sub-network (with parameters θ_{FP}), on the other hand, passes \mathbf{x}_{FP} first into an Embedding layer that converts the integer input vector into a (100, 32) embedding. This embedding is then fed into 4 parallel convolution blocks with filters of size 2, 3, 4 and 5 respectively, each composed by a 1-D convolution layer with 128 filters, a Layer Normalization layer and a 1-D sum layer to flatten the output to a vector. The flattened outputs of these convolution blocks are then concatenated and serve as input to two dense blocks similar to those found in the PE input arm.

Finally, the outputs from the two base sub-networks are then concatenated and passed into the joint output path with parameters θ_O . This output sub-network is composed by dense blocks (same form as in the PE input arm) of layer sizes 512, 256 and 128. Finally, a single fully-connected layer is employed to project the 128-D output from the previous blocks into a 1-D output, followed by a sigmoid activation function that provides the final output score of the model.

The PE-only model can be easily derived from the PE+FP model removing the FP arm,

taking input \mathbf{x}_{PE} and fitting parameters θ_{PE} and θ_O . Similarly, the FP-only model can be constructed by using the PE+FP model but without the PE sub-network, taking input \mathbf{x}_{FP} and fitting parameters θ_{FP} and θ_O . Obviously the first layer of the output sub-network needs to be modified to match the output from the previous layer.

In [45], the authors fit all models using a binary cross entropy loss function. Therefore, denoting the output of the model as $f(\mathbf{x}; \theta)$, where \mathbf{x} is the model input, θ are the model parameters and $y \in \{0,1\}$ are the ground truth labels, the loss is:

$$L(\mathbf{x}, y; \theta) = -y \log(f(\mathbf{x}; \theta)) + (1 - y) \log(1 - f(\mathbf{x}; \theta)). \quad (3.18)$$

During training, the equation is optimized for $\hat{\theta}$: the optimal set of parameters that minimize the combined loss over the dataset.

$$\hat{\theta} = \arg \max_{\theta} \sum_{i=1}^M L(\mathbf{x}^{(i)}, y_i; \theta), \quad (3.19)$$

where M is the number of samples in the dataset, and y_i and $\mathbf{x}^{(i)}$ are the label and the feature vector of the i th training sample respectively [45].

3.7 Malware Normalization

In order to improve the detection rate of existing anti-malware techniques also against malware produced by advanced packers and toolkits, code *normalization* techniques can be exploited. These techniques consist of a *normalizer* which accepts obfuscated code as input and tries to eliminate obfuscation producing as output the normalized executable.

After *normalization*, the usual signature-based techniques can be applied on the normalized sample [22].

Chapter 4

Workflow and Datasets used

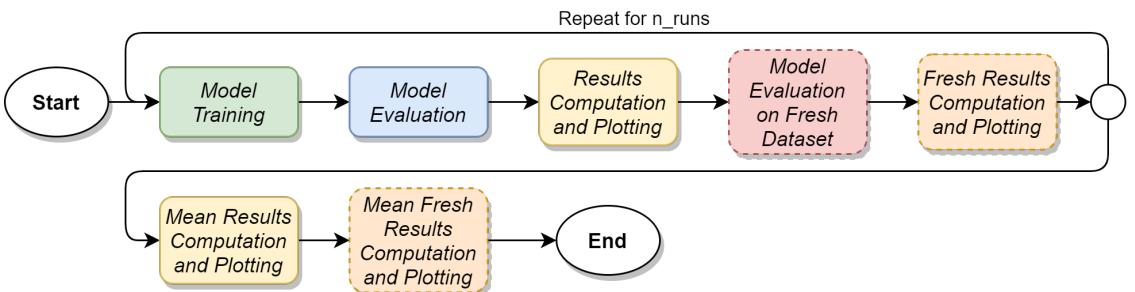


Figure 4.1: Overall workflow

The workflow presented in image 4.1 summarizes the steps taken throughout the experiments conducted and reported in this document. In particular, for each different model architecture and combination of parameters, the model is first trained and evaluated on the large-scale dataset called ***Sorel 20M dataset*** [6] on the task of SMART tagging and/or malicious/benign label prediction. Next, the model evaluation results are computed and plotted. Depending on the model type, the model is then optionally evaluated also using a smaller dataset, referred to as '***Fresh dataset***' in this document, specifically crafted for the task of predicting the malware family each specific sample belongs to. The results of this second evaluation are computed and plotted in a separate step. The whole process presented until now is repeated for n_runs times and finally the mean results for both evaluation types are computed, aggregating the results of the single runs.

4.1 Sorel 20M Dataset

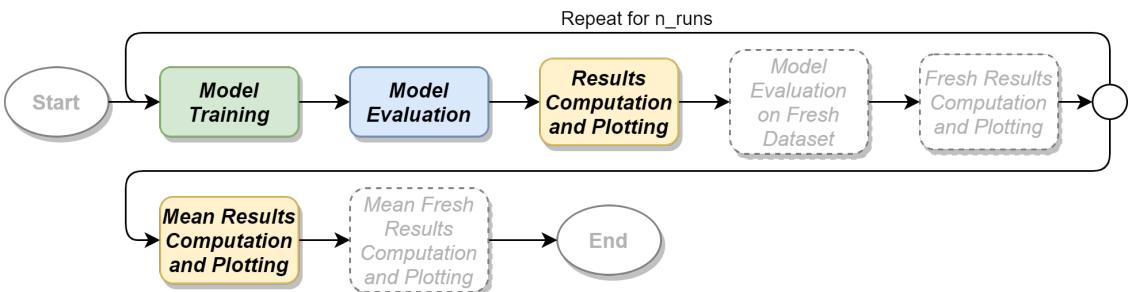


Figure 4.2: Workflow steps which use the Sorel20M dataset

The ***Sorel20M*** large-scale dataset, which was released in 2020 and is used in the workflow steps shown in image 4.2, consists of nearly 20 million files, each represented by the corresponding

pre-extracted features, metadata, malicious/benign label derived from multiple sources, vendor detections count information and additional SMART 'tags' to serve as additional targets [6].

SOREL20M dataset builds upon EMBER dataset [48], the first standard and open dataset specifically designed to be used for static malware detection, addressing some of its drawbacks that limited its utility as a malware benchmark set. In particular, EMBER contains 900,000 training samples (300K malicious, 300K unlabelled) and 200,000 test samples (100K malicious, 100K benign) and is therefore of limited size compared to the private/proprietary datasets containing from tens to hundreds of millions of samples commercial malware models are usually trained on. Additionally, the small number of validation samples makes it difficult to evaluate the model performance at low false positive rates due to high variance. Moreover, the EMBER dataset only provides a single binary malware/benignware label with no additional information.

SOREL20M, in fact provides an order of magnitude more samples for analysis: when using the recommended time splits to establish training, validation and test sets, there are 12,699,013 training samples, 2,495,822 validation samples and 4,195,042 test samples, respectively. Furthermore, Sorel20M [6] provides 9,919,251 samples of malware (7,596,407 training samples, 962,222 validation samples and 1,360,622 test samples), which have been 'disarmed' by setting both the *optional_headers.subsystem* and *file_header.machine* to 0 in order to prevent execution. Additionally, complete PE metadata, obtained via the Python *pefile* module using the *dump_dict* method, is provided for each sample.

Finally, the dataset provides for each sample a number of additional targets (SMART tags) for the model that describe behaviour inferred from vendor family labels, together with the vendor detection count and the malicious/benign label.

4.1.1 Sorel 20M Dataset Description

The complete dataset, which is available as a AWS bucket, consists of the following items:

- 9,919,251 original (disarmed) malware samples compressed via the Python *zlib.compress* function
- A SQLite3 and two LMDB databases:
 - The SQLite3 "meta.db" database containing malware labels, tags, detection counts, and first/last seen times
 - The "ember_features" LMDB database containing the EMBER features (EMBER features version 2)
 - the "pe_metadata" LMDB database containing the PE metadata extracted through the *pefile* module, as described above
- Moreover, some Pre-trained baseline models (a Pytorch feed-forward neural network (FFNN) model and a LightGBM gradient-boosted decision tree model) and their results are also provided, but will not be used in this document.

All samples are identified by their sha256 hash which serves as the primary key for the SQLite3 database, and the key to be used to access the two LMDB databases. LMDB entries are stored as arrays or dictionaries that are then serialized with *msgpack* and compressed with *zlib*.

The data was collected from January 1st, 2017 to April 10th, 2019. In [6], Harang et al. suggest to use the following time-splits of the data (based on the first-seen time in RL telemetry): training data from the beginning of collection until November 29th, 2018; validation data from then until January 12th, 2019; and testing data from January 12th, 2019 through the end of the data.

LMDB database, what is it? Lightning Memory-Mapped Database (LMDB) is a B+-trees-based database management library that provides a high-performance embedded transactional database with full ACID semantics in the form of a key-value store (it is not a relational database).

The entire database is exposed in a memory map, and all data fetches return data directly from the mapped memory, so no `'malloc`'s nor `'memcpy`'s occur during data fetches. Therefore, the library is extremely simple as it requires no page caching layer of its own (the OS is responsible for managing the pages), and it is extremely high performant and memory-efficient.

The memory map can be used as a read-only or read-write map. It is read-only by default as this provides total immunity to corruption. On the other hand, using read-write mode offers much higher write performance [49]. LMDB may also be used concurrently in a multi-threaded or multi-processing environment, with read performance scaling linearly by design. In particular it uses shared memory copy-on-write semantics with a single writer; however unlike many similar key-value databases, write transactions do not block readers, nor do readers block writers.

4.1.2 Ember Features

As previously mentioned, Sorel20M dataset consists of a bunch of databases (Sqlite3 and LMDB), indexed by the sha256 hash of the files, which contain the feature vectors representing the samples together with the corresponding labels/metadata. In particular the feature vectors were derived from the samples' code through the use of an open source feature extraction code released by Anderson et al. [48] in 2018 as an attachment to the Endgame Malware BEncmark for Research (EMBER) dataset. In particular, Sorel20M uses version 2 of this feature extraction code which adds information on data directories to the feature representation with respect to version 1.

The PE file format and the features extracted by the EMBER feature extraction code will be described in the next two sections. Before that, here is a brief description of the EMBER dataset useful for context.

The EMBER dataset, which was extracted from a large corpus of Windows Portable Executable (PE) malicious and benign files, consists of a collection of JSON lines files, where each line contains a single JSON object. Each object includes the following types of data:

- the sha256 hash of the original file as a unique identifier;
- coarse time information (month resolution) about when the file was first seen;
- a malicious/benign label, which may be 0 for benign, 1 for malicious or -1 for unlabelled;
- 8 groups of raw features that include both parsed values as well as format-agnostic histograms.

A notable difference with respect to Sorel20M, however, is that the EMBER dataset is comprised of raw features that are human-readable instead of directly having the feature vectors required for model building. This was done to allow researchers to decouple raw features from the vectorizing strategy and to improve model interpretability. Anyway, the authors of [48] provided also the code for producing numeric feature vectors from those raw features. The raw feature extraction code and the numeric feature vector generation code were used in succession by Sorel20M authors to directly obtain the feature vector representations from samples.

PE File Format

The Portable Executable (PE) format is a file format for executables, object code, dynamically-linked libraries (DLLs), FON font files and more, used in 32-bit and 64-bit versions of Windows operating systems (it is currently supported on Intel, AMD and variants of ARM instruction set architectures).

The PE format structure consists of a number of standard headers (4.3) followed by one or more sections and it encapsulates the information necessary for the Windows OS loader to manage

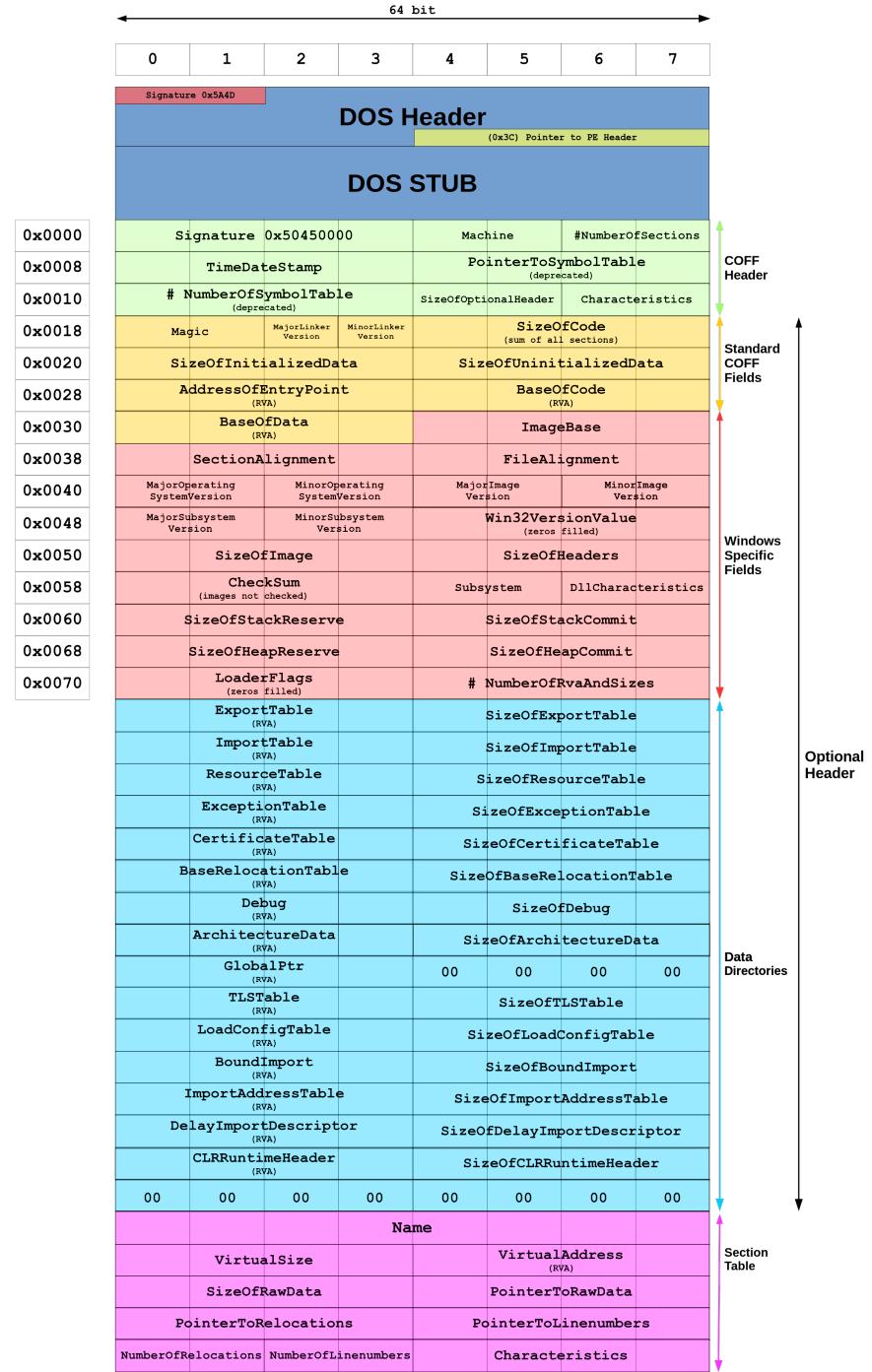


Figure 4.3: PE file structure

the wrapped executable code: for example, among other things, it tells the dynamic linker how to map the file into memory. In particular, the PE structure usually consist of the following 3 headers:

- *Common Object File Format (COFF)* file header: it contains important information about the file such as the type of machine for which it is intended, its nature (DLL, EXE, OBJ), the number of sections, the number of symbols, etc.
- *Optional Header*: it can be further divided into:
 - *Standard COFF fields*: they identify the linker version, the size of the code, the size of initialized and uninitialized data, the address of the entry point, etc.
 - *Windows Specific fields*: they provide windows-specific information such as minor and major operating system, subsystem and image versions, stack and heap sizes, section and file alignment, etc.
 - *Data Directories*: they provide pointers to the sections that follow it, which include tables for exports, imports, resources, exceptions, debug information, certificate information, and relocation tables.
- *Section Table*: it outlines the name, offset and size of each section in the PE file.

PE sections, on the other hand, contain code and initialized data that will be mapped into executable or readable/write-able memory pages, respectively, by the Windows loader at execution time, as well as imports, exports and resources defined by the file. Each section contains a header that specifies its size and address. A windows executable typically has the nine predefined sections named *.text*, *.bss*, *.rdata*, *.data*, *.rsrc*, *.edata*, *.idata*, *.pdata*, and *.debug*. Some applications however do not need all of these sections, while others may define still more sections to suit their specific needs.

Among the different sections a PE file may contain here are described some of the most common:

- *Executable code section, .text*: it holds the program code. The *.text* section also contains the code entry point and the *Import Address Table (IAT)* (which is used as a lookup table when the application is calling a function in a different module).
- *Data sections, .bss, .rdata, .data*: The *.bss* section contains uninitialized data for the application, the *.rdata* section contains read-only data. All other application/global variables are stored in the *.data* section.
- *Resources section, .rsrc*: it contains resource information for a module, such as the ones required for user interfaces: cursors, fonts, bitmaps, icons, menus, etc.
- *Export data section, .edata*: it contains export data for the application or DLL.
- *Import data section, .idata*: it contains import data, including the import directory and import address name table.
- *.reloc section*: it stores *Relocation tables* which are used by the Windows loader to rebase the PE file if it cannot be loaded at its preferred base address. In fact, PE files normally do not contain position-independent code but are compiled to a preferred base address, and all addresses emitted by the compiler/linker are fixed ahead of time.
- *.tls section*: it contains special thread local storage (TLS) structures for storing thread-specific local variables.

A basic PE file normally contains a *.text* code section and one or more data sections (*.data*, *.rdata* or *.bss*).

It is interesting to notice how packers sometimes create new sections, for example, the UPX packer creates a section *UPX1* to store packed data and an empty section *UPX0* that reserves an address range for runtime unpacking [48].

Feature Set Description

EMBER feature extraction code extracts 8 groups of raw features that include both parsed features and format-agnostic histograms and string counts.

In particular, the EMBER dataset authors [48] made a distinction between human-readable *raw features* and numerical *model features* (or *vectorized features*) derived from the dataset. *Model features* consist of a feature matrix of fixed size used for training models, representing the numerical summary of raw features, wherein strings, imported names, exported names, etc., are captured using the feature hashing trick [50].

Parsed features The first 5 groups of features are extracted after parsing the PE file. Anderson et al. [48] leveraged the *Library to Instrument Executable Formats (LIEF)* [51] as a convenient PE parser. *LIEF* names are used for strings that represent symbolic objects, such as characteristics and properties.

- **General file information** - The set of raw features belonging to the general file information group includes the file size and basic information obtained from the PE header: the virtual size of the file, the number of imported and exported functions, whether the file has a debug section, thread local storage, resources, relocations, and the number of symbols etc.
- **Header information** - From the *COFF* header, the timestamp, the target machine (string) and a list of image characteristics (list of strings) are extracted. From the optional header, instead, the EMBER feature extracting code extracts the target subsystem (string), DLL characteristics (a list of strings), the file magic as a string (e.g. "PE32"), major and minor image versions, linker versions, system versions and subsystem versions, and the code, headers and commit sizes. Then, in order to create the model features, string descriptors such as DLL characteristics, target machine, subsystem, etc. are summarized using the feature hashing trick, with 10 bins allotted for each indicator vector.
- **Imported functions** - Imported functions and libraries are extracted from the parsing of the import address table. To create numerical features, the feature vector generation code collects the set of unique library names and uses the hashing trick with 256 bins. Similarly, the hashing trick (with 1024 bins) is used to capture individual function names, by representing each as a string in the *library:FunctionName* format (e.g. *kernel32.dll>CreateFileMappingA*).
- **Exported functions** - The raw features include a list of the exported functions. These strings are summarized into model features using the hashing trick with 128 bins.
- **Section information** - Specific section properties are extracted from each section. They include the section name, size, entropy, virtual size, and a list of strings representing the section characteristics. Then the entry point is captured and specified by name. To convert them to model features, the hashing trick is again used on (section name, value) pairs to create vectors containing section size, section entropy, and virtual size (50 bins each). The hashing trick is also used on the entry point characteristics (list of strings).

Format-agnostic features The next 3 groups of features are instead format agnostic, meaning that they do not require the PE file to be parsed for their extraction.

- **Byte histogram** - The byte histogram contains 256 integer values, representing the counts of each byte value within the file. When generating model features, this byte histogram is normalized to a distribution, since the file size is represented as a feature in the general file information.
- **Byte-entropy histogram** - The byte entropy histogram, on the other hand, approximates the joint distribution $p(H, X)$ of entropy H and byte value X . This is done as described by Saxe et al. in [47], by computing the scalar entropy H for a fixed-length window and

pairing it with each byte occurrence within the window. This is repeated as the window slides across the input bytes. In particular, EMBER feature extraction code uses a window size of 2048 and a step size of 1024 bytes, with 16×16 bins that quantize entropy and the byte value. These counts are then normalized to sum to 1.

- **String information** - EMBER features also include simple statistics about printable strings (consisting of characters in the range `0x20` to `0x7f`, inclusive) that are at least five printable characters long. In particular, the code extracts information like the number of strings, their average length, a histogram of the printable characters within those strings, and the entropy of characters across all printable strings. In addition, the string feature group also provides the number of strings that begin with specific character sequences such as `C:\` (case insensitive) that may indicate a path, `http://` or `http://` (case insensitive) that may indicate a *URL*, `HKEY_` that may indicate a registry key, etc.

4.1.3 Trying to improve Dataset Loading Speed

Dataset Pre-Processing

Harang et al. [6] provided, together with the Sorel20M dataset, the python/pytorch code they used to load it in memory and pass it as input to their model. This code, however, is particularly I/O bounded. In fact, in order to load the data related to a single sample i it first gets the corresponding sha256 hash from the Sqlite3 database (which is already loaded in memory) and then uses it as key to access the LMDB database of features. The extracted features are then decompressed and deserialized. To load the entire dataset, one batch of data of size `batch_size` (= 8192 by default) is loaded at a time, repeating this process for `batch_size` randomly chosen samples i for each batch until the end of the epoch. The code is therefore especially dependant to the hard disk random access speed. The authors of [6] managed to train their model for 10 epochs in under 90 minutes, by training it on an AWS instance with high I/O speeds (exploiting a NVMe SSD), not suffering from any bottleneck. However, when using a less powerful instance not optimized for I/O performance the time needed for loading the dataset and subsequently for training the model increased excessively. This made the use of the original data loading code for the purposes of this project unfeasible.

In order to speed up the dataset loading code the dataset is therefore pre-processed and saved in a easier (and faster) to read format. In particular the entire dataset is sequentially loaded into a set of 3 parallel memory mapped `numpy` array (for the features, labels and sha256 hashes respectively) using the original data loading code after being pre-processed (the features are decompressed and deserialized), similarly to what was done in the original code every cycle. The resulting `numpy` arrays are then saved 'as-is' to file. During training (or evaluation) the pre-processed dataset files are read back into memory mapped `numpy` arrays 'as-is' and then used in conjunction with the default Pytorch Dataloader class, with no additional processing needed. This greatly speeded up the overall training. The main drawback of this approach is that the dataset is saved with its features being decompressed, thus taking up a much larger disk portion.

When using an instance with limited disk size the amount of space occupied by the dataset is a concern. In particular, the instance used for this project had up to 130 GB of disk space and this made it possible to train the models with only approximately half of the samples provided by Sorel20M. In fact for all the experiments only the first 6.000.000 of the 12.699.013 training samples, the first 1.153.846 of 2.495.822 validation samples and the first 1.846.154 of 4.195.042 test samples of Sorel20M dataset were considered. The models presented in this document therefore suffered from the smaller dataset size meaning that they cannot be directly compared with other models which used Sorel20M as base dataset. However, the code was designed to work with any dataset size so additional experimentations with more powerful instances can easily be made in the future.

Generator (Dataloader) versions

Pytorch Dataloader In the first implementation of the data loading code the original *Pytorch* Dataloader class was used by passing it a specially crafted class called '*Dataset*' (derived from

Pytorch ‘`torch.utils.data.Dataset`’ class) which is responsible for loading the pre-processed version of the dataset and for retrieving the data (features, labels and sha256) for a specific sample ‘*i*’.

In particular, as it can be seen in algorithms 1 and 2, the *Dataset* class consists of 2 main member functions (although other less important member functions such as the *Len* function are also implemented): *Init* and *GetItem*. In the *Init* constructor function (alg. 1) the dataset, consisting of features (*X*), labels (*y*) and sha256 hashes (*S*), gets loaded into a set of 3 parallel memory mapped arrays. In the *GetItem* member function (alg. 2), instead, the data (features, labels and/or sha256) corresponding to a specific sample with index ‘*index*’ is retrieved and returned to the function caller (that will be the Pytorch original Dataloader itself).

Algorithm 1 Dataset class, *Init*

```

1: class DATASET
2:   function INIT(self, ds_root, mode, n_samples, return_shas, ...)
3:     self.return_shas  $\leftarrow$  return_shas
4:     ...
5:     X_path  $\leftarrow$  os.path.join(ds_root, 'X_{\}.\{}.dat').format(mode, n_samples)'
6:     y_path  $\leftarrow$  os.path.join(ds_root, 'y_{\}.\{}.dat').format(mode, n_samples)'
7:     S_path  $\leftarrow$  os.path.join(ds_root, 'S_{\}.\{}.dat').format(mode, n_samples)'
8:
9:     self.S  $\leftarrow$  load_as_memmap(S_path, dtype=np.dtype('U64'), mode="r+")
10:    self.y  $\leftarrow$  load_as_memmap(y_path, dtype=np.float32, mode="r+")
11:    self.X  $\leftarrow$  load_as_memmap(X_path, dtype=np.float32, mode="r+")
12:  end function

```

Algorithm 2 Dataset class, *.GetItem*

```

13:  function GETITEM(self, index)
14:    features  $\leftarrow$  self.X[index]
15:
16:    labels  $\leftarrow$  {}
17:    labels['malware']  $\leftarrow$  self.y[index][0]
18:    labels['count']  $\leftarrow$  self.y[index][1]
19:    labels['tags']  $\leftarrow$  self.y[index][2:]
20:
21:    if self.return_shas then
22:      sha  $\leftarrow$  self.S[index]
23:      return sha, features, labels
24:    else
25:      return features, labels
26:    end if
27:  end function
28: end class

```

Before model training/evaluation the dataset generator (dataloader) is defined by passing this dataset class to the *Pytorch* Dataloader implementation (*torch.utils.data.Dataloader*) together with additional arguments specifying the batch size, the number of workers and whether to shuffle the data during loading (alg. 3). Then, during training, the *Pytorch* Dataloader class will load the batches of data by iteratively calling the *Dataset*’s *GetItem* function *batch_size* times for each batch concatenating the extracted samples together using the *torch.cat* function.

As previously mentioned, this implementation is faster than the original Sorel20M dataloader code while being also relatively simple, however given how the pre-processed dataset is saved (consecutive samples are saved on consecutive locations on disk) it is still somewhat inefficient. In fact, the time needed for completing 10 epochs training on a less powerful instance is still unreasonable. One possible optimization would be to read batches in one go as big chunks rather than calling the *GetItem* function on the Dataset once per sample and then concatenating the resulting data. However, random sampling during model training introduces a regularization

effect which generally improves model generalization on unseen samples, therefore it would be better to find an optimization combining both data loading efficiency and random sampling.

Algorithm 3 Pytorch Dataloader definition

```

1: ds  $\leftarrow$  Dataset(ds_root, mode, n_samples, return_shas, ...)
2: generator  $\leftarrow$  Pytorch_Dataloader(ds, batch_size, shuffle, n_workers)
  
```

Generator alt1/alt2 The first attempts at optimizing the generator were inspired by the '*index select*' and '*shuffle in-place*' versions of **FastTensorDataLoader** suggested in [52].

In particular, the alternative generator 1 ('*alt1*'), again loads the pre-processed dataset as a set of memory mapped arrays and assigns them to a set of tensors (X , y and S) from which batches of data are randomly sampled using the *Pytorch 'index_select'* function in *multithreading*.

In the alternative generator 2 ('*alt2*'), on the other hand, the tensors created by loading the dataset as was done in *alt1* are randomly shuffled in place at each iteration. Then the batches of data are sampled (in order) from them in *multithreading*.

Algorithm 4 Alt1/Alt2 FastTensorDataLoader class, Init function

```

1: function INIT(self, tensors, batch_size, shuffle, n_workers, ...)
2:   self.tensors  $\leftarrow$  tensors
3:   self.batch_size  $\leftarrow$  batch_size
4:   self.shuffle  $\leftarrow$  shuffle
5:   self.n_workers  $\leftarrow$  n_workers
6:   self.dataset_len  $\leftarrow$  tensors[0].shape[0]
7:
8:   if num_workers  $> 1 then
9:     self.async_results  $\leftarrow$  []
10:    self.pool  $\leftarrow$  ThreadPool()
11:   end if
12:
13:   self.n_batches  $\leftarrow$   $\lceil \text{self.dataset\_len} / \text{self.batch\_size} \rceil$ 
14: end function$ 
```

Algorithm 5 Alt1 FastTensorDataLoader class, Iter function

```

1: function ITER(self)
2:   if self.shuffle = true then
3:     self.indices  $\leftarrow$  randperm(self.dataset_len)
4:   else
5:     self.indices  $\leftarrow$  None
6:   end if
7:
8:   self.i  $\leftarrow$  0
9:   return self
10: end function
  
```

In practice, both versions of the generator share (for the most part) the *Init* and *Next* functions. The main differences are, instead, in the *Iter* and *get_batch* functions.

The *Init* function (alg. 4) is used for setting up some **FastTensorDataLoader** variables and optionally initializing the **ThreadPool** used for *multithreading*.

The *Iter* function purpose is to set up the parameters for the current cycle at the beginning of each epoch. The two generator alternatives have, however, different *Init* functions. In particular, generator *alt1* *Iter* function (alg. 5) resets the count of extracted samples and, if shuffling is enabled, randomly generates the current epoch sample indices (which define the order the samples

Algorithm 6 Alt2 FastTensorDataLoader class, Iter function

```

1: function ITER(self)
2:   if self.shuffle = true then
3:     r  $\leftarrow$  randperm(self.dataset_len)
4:     for i, t  $\in$  enumerate(self.tensors) do
5:       self.tensors[i]  $\leftarrow$  t[r]
6:     end for
7:   end if
8:
9:   self.i  $\leftarrow$  0
10:  return self
11: end function

```

Algorithm 7 Alt1/Alt2 FastTensorDataLoader class, Next function

```

1: function NEXT(self)
2:   if self.i  $\geq$  self.dataset_len then
3:     if (self.n_workers = 1 or len(self.async_results) = 0) then
4:       raise StopIteration
5:     end if
6:   end if
7:
8:   if self.num_workers = 1 then
9:     batch  $\leftarrow$  get_batch(self.tensors, self.batch_size, self.i, (self.indices))
10:    self.i  $\leftarrow$  self.i + self.batch_size
11:    return batch
12:   else
13:     while self.i < self.dataset_len and len(self.async_results) < self.n_workers do
14:       arguments  $\leftarrow$  (self.tensors, self.batch_size, self.i, (self.indices))
15:       async_task  $\leftarrow$  self.pool.apply_async(get_batch, arguments)
16:       self.async_results.append(async_task)
17:       self.i  $\leftarrow$  self.i + self.batch_size
18:     end while
19:
20:     current_result  $\leftarrow$  self.async_results.pop(0)
21:     return current_result.get()
22:   end if
23: end function

```

are extracted from the dataset tensors). On the other hand, generator *alt2 Iter* function (alg. 6) again resets the count of extracted samples, but if shuffling is enabled, it randomly permutes *in-place* the dataset tensors themselves.

In order to get a batch of data, the *Next* function (alg. 7), which is similar for both generator versions, is used. In particular, this function monitors the number of samples already extracted from the dataset tensors and raises a *StopIteration* exception when it has cycled through all the data. Moreover, this function is also responsible for extracting and returning a batch of data from the dataset tensors. This can be done in sequence or in *multithreading*. Specifically, when the number of workers selected is 1 the function sequentially gets one single batch and returns it after having updated the count of already extracted samples for the current epoch. On the other hand, if more than 1 workers are used the function prepares a number of asynchronous batch extraction tasks to be run in parallel by the threads in the thread pool, then it waits for the first result and returns it. The next call to the function will prepare other tasks and wait for the next first result. The exact number of asynchronous tasks prepared at each call is chosen dynamically to keep an array of async task results of size fixed to *num_workers* always full.

Finally, in algorithm 8 is shown the *get_batch* function used by generator *alt1*: if the sample *indices* are provided they are used with the *index_select* function to select the data samples for

Algorithm 8 Alt1 get_batch function

```

1: function GET_BATCH(tensors, batch_size, i, indices, ...)
2:   batch  $\leftarrow$  []
3:   if indices is provided then
4:     indices  $\leftarrow$  indices[i:(i + batch_size)]
5:
6:     for all t  $\in$  tensors do
7:       batch_data  $\leftarrow$  index_select(t, indices)
8:       batch.append(batch_data)
9:     end for
10:   else
11:     for all t  $\in$  tensors do
12:       batch_data  $\leftarrow$  t[i:(i + batch_size)]
13:       batch.append(batch_data)
14:     end for
15:   end if
16:
17:   return batch
18: end function

```

Algorithm 9 Alt2 get_batch function

```

1: function GET_BATCH(tensors, batch_size, i, ...)
2:   batch  $\leftarrow$  []
3:   for all t  $\in$  tensors do
4:     batch_data  $\leftarrow$  t[i:(i + batch_size)]
5:     batch.append(batch_data)
6:   end for
7:
8:   return batch
9: end function

```

the current batch of data, otherwise a set of consecutive samples is drawn from the dataset. The *get_batch* function used by generator *alt2* (alg. 9) is simpler: it simply selects a set of consecutive samples from the dataset tensors since those were already shuffled if required.

Unfortunately, the presented two generator alternatives did not improve the data loading process enough. More specifically, when considering 6M samples, generator *alt1* was only slightly faster than the original Pytorch Dataloader, while *alt2* was even significantly slower. Generator *alt1*, in fact, was still fairly similar to the Pytorch Dataloader implementation having used the *index_select* function instead of the slower *torch.cat*. On the other hand, generator *alt2* worked by shuffling the entire dataset tensors prior to batch extraction. If this solution surely can be more efficient for small to moderately sized dataset, it can become a bottleneck in case of huge datasets.

Generator alt3 The generator alternative 3 ('*alt3*') managed to considerably speed up the dataset loading process making a tradeoff between loading speed and samples dispersion. In fact, the time needed for 1 training epoch (considering 6M samples) passed from being of approximately 6 hours to \sim 15 minutes.

This optimization uses a new FastTensorDataLoader class which exploits a pool of threads to asynchronously load the dataset into memory in chunks of consecutive data. In particular each thread loads into memory '*n_chunks*' randomly chosen chunks. Each thread then proceeds to concatenate together its '*n_chunks*' chunks, which contain '*chunk_size*' malware samples each, generating each a '*chunk_aggregate*', which is then randomly shuffled and returned by the thread. The '*chunk_aggregates*' are asynchronously inserted in a queue (the order depends on the threads instantiation order) of fixed length equal to the number of workers used. The main

dataloader thread manages the *chunk_aggregate* queue, instantiating the parallel threads such that the queue is always full. Moreover, the main dataloader thread sequentially extracts, when possible, one *chunk_aggregate* at a time from the queue and then proceeds to return one batch of data at a time from it, when required.

Algorithm 10 Alt3 FastTensorDataLoader class, Init

```

1: class FASTTENSORDATALOADER
2:   function INIT(self, tensors, batch_size, chunk_size, chunks, shuffle, n_workers, ...)
3:     self.tensors  $\leftarrow$  tensors
4:     self.batch_size  $\leftarrow$  batch_size
5:     self.chunk_size  $\leftarrow$  chunk_size
6:     self.chunks  $\leftarrow$  chunks
7:     self.shuffle  $\leftarrow$  shuffle
8:     self.n_workers  $\leftarrow$  n_workers
9:     self.dataset_len  $\leftarrow$  tensors[0].shape[0]
10:
11:    if num_workers > 1 then
12:      self.async_results  $\leftarrow$  []
13:      self.pool  $\leftarrow$  ThreadPool()
14:    end if
15:
16:    self.n_batches  $\leftarrow$   $\lceil \text{self.dataset\_len} / \text{self.batch\_size} \rceil$ 
17:    self.n_chunks  $\leftarrow$   $\lceil \text{self.dataset\_len} / \text{self.chunk\_size} \rceil$ 
18:    self.last_chunk_size  $\leftarrow$  self.dataset_len % self.chunk_size
19:  end function

```

Algorithm 11 Alt3 FastTensorDataLoader class, Iter

```

20:  function ITER(self)
21:    if self.shuffle = true then
22:      self.chunk_indices  $\leftarrow$  randperm(self.dataset_len)
23:    else
24:      self.chunk_indices  $\leftarrow$  arange(self.n_chunks)
25:    end if
26:
27:    self.chunk_agg  $\leftarrow$  None
28:    self.chunk_agg_size  $\leftarrow$  0
29:    self.chunk_i  $\leftarrow$  0
30:    return self
31:  end function

```

More specifically, in the *Init* function (alg. 10), which is similar to the *Init* of the previous alternatives, some **FastTensorDataLoader** variables are set/computed, and a **ThreadPool** is initialized if necessary.

The *Iter* function (alg. 11), instead, is used to reset the current *chunk aggregate* (and its size) and the number of already processed chunks for the current cycle (or *epoch*) at the beginning of each *epoch*. Moreover, it also initializes the *chunk indices*, which define the order in which the data chunks are retrieved, randomly or sequentially (linearly) depending on whether shuffling is required.

In algorithm 12 is presented the **FastTensorDataLoader alt3** *Next* member function, which is used for managing the asynchronous extraction of chunks of consecutive samples from the dataset and for returning a single batch of data. In particular, when the selected number of workers is 1 then if all the data of the current *chunk aggregate* has been cycled through (or it is the first function call) the function simply extracts a single new *chunk aggregate* (which is a concatenation of multiple chunks selected randomly or linearly depending on the value of *self.shuffle* variable) through the **get_chunks** function and updates the number of already processed chunks for the

Algorithm 12 Alt3 FastTensorDataLoader class, Next

```

32:   function NEXT(self)
33:     if self.n_workers = 1 then
34:       if self.chunk_agg is None or self.i  $\geq$  self.chunk_agg_size then
35:         if self.chunk_i  $\geq$  self.n_chunks then
36:           raise StopIteration
37:         end if
38:
39:         start_i  $\leftarrow$  self.chunk_i
40:         end_i  $\leftarrow$  start_i + self.chunks
41:         arguments  $\leftarrow$  (self.tensors, self.chunk_indices[start_i:end_i],
42:                           self.chunk_size, self.last_chunk_size,
43:                           self.n_chunks, self.shuffle)
44:         self.chunk_agg, self.chunk_agg_size  $\leftarrow$  get_chunks(arguments)
45:         self.chunk_i  $\leftarrow$  end_i
46:         self.i  $\leftarrow$  0
47:       end if
48:     else
49:       while self.chunk_i < self.n_chunks and len(self.async_results) < self.n_workers do
50:         start_i  $\leftarrow$  self.chunk_i
51:         end_i  $\leftarrow$  start_i + self.chunks
52:         arguments  $\leftarrow$  (self.tensors, self.chunk_indices[start_i:end_i],
53:                           self.chunk_size, self.last_chunk_size,
54:                           self.n_chunks, self.shuffle)
55:         async_task  $\leftarrow$  self.pool.apply_async(get_chunks, arguments)
56:         self.async_results.append(async_task)
57:         self.chunk_i  $\leftarrow$  end_i
58:       end while
59:
60:     if self.chunk_agg is None or self.i  $\geq$  self.chunk_agg_size then
61:       if len(self.async_results) = 0 then
62:         raise StopIteration
63:       end if
64:
65:       current_result  $\leftarrow$  self.async_results.pop(0)
66:       self.chunk_agg, self.chunk_agg_size  $\leftarrow$  current_result.get()
67:       self.i  $\leftarrow$  0
68:     end if
69:   end if
70:
71:   batch  $\leftarrow$  get_batch(self.chunk_agg, self.batch_size, self.i)
72:   self.i  $\leftarrow$  self.i + batch[0].shape[0]
73:   return batch
74: end function
75: end class

```

current *epoch* and the number of extracted samples from that specific *chunk aggregate*. This is however done only if the number of chunks already processed is less than the total amount of chunks, otherwise the *StopIteration* exception is raised. When the number of workers is greater than 1, on the other hand, until the number of extracted chunks is less than the total amount the function prepares a number of asynchronous tasks of *chunk aggregate* extraction such that the queue (of size equal to the number of selected number of workers) in which they are inserted is always full, and also updates the number of extracted chunks appropriately. Then when the current *chunk aggregate* data has been explored completely (or when it is the first function call), if the async task queue is empty it raises the *StopIteration* exception because it means there are no more *chunk aggregates* for the current *epoch*; otherwise the function gets the first async task

Algorithm 13 Alt3 get_chunks function

```

1: function GET_CHUNKS(tensors, chunk_indices, chunk_size, last_chunk_size, n_chunks, shuffle)
2:   if n_chunks - 1 in chunk_indices then
3:     chunk_agg_size  $\leftarrow$  (len(chunk_indices) - 1) · chunk_size + last_chunk_size
4:   else
5:     chunk_agg_size  $\leftarrow$  len(chunk_indices) · chunk_size
6:   end if
7:
8:   chunk_agg  $\leftarrow$  []
9:   for all t in tensors do
10:    chunk_agg.append(emptyTensor)
11:   end for
12:
13:   c_start  $\leftarrow$  0
14:   for all idx  $\in$  [0, ..., len(chunk_indices)] do
15:     t_start  $\leftarrow$  chunk_indices[idx] · chunk_size
16:     if chunk_indices[idx]  $\neq$  n_chunks - 1 then
17:       c_end  $\leftarrow$  c_start + chunk_size
18:       t_end  $\leftarrow$  t_start + chunk_size
19:     else
20:       c_end  $\leftarrow$  c_start + last_chunk_size
21:       t_end  $\leftarrow$  t_start + last_chunk_size
22:     end if
23:
24:     for i, t in enumerate(tensors) do
25:       chunk_agg[i][c_start:c_end]  $\leftarrow$  t[t_start:t_end]
26:     end for
27:     c_start  $\leftarrow$  c_end
28:   end for
29:
30:   if shuffle then
31:     r  $\leftarrow$  randperm(chunk_agg_size)
32:     for i, t in enumerate(chunk_agg) do
33:       chunk_agg[i]  $\leftarrow$  t[r]
34:     end for
35:   end if
36:
37:   return chunk_agg, chunk_agg_size
38: end function

```

from the queue and waits for its result (which is the extracted *chunk aggregate*). Independently from the number of workers used, the function then proceeds to efficiently get a single batch of data from the current already-in-memory *chunk aggregate* and returns it.

The **GetBatch** function used in generator *alt3* is the same one already exploited in version *alt2* (9), however, *alt3* sees the addition of the **GetChunks** function (13) which is used to extract a number of chunks of consecutive data from the dataset and combine them into an *aggregate chunk*. In particular, the function first computes the size of the resulting *chunk aggregate* depending on the number of chunk indices specified and the chunk size (considering as a special case the last chunk of the dataset which may have a smaller size). Then, the *chunk aggregate* itself is initialized as a series of properly sized empty tensors (one for samples features, one for labels and one for sha256 hashes) which are later filled sequentially by the data chunks got from the dataset tensors. Finally, the *chunk aggregate* data gets shuffled in place, if needed, and returned.

As it can be seen from the algorithms provided, this version of the generator depends on the value a number of parameters among which the most important are *chunk_size* and *n_chunks*. Selecting different values for *chunk_size* and *n_chunks*, in fact, has an impact on the speed of the generator and on the samples dispersion. In fact the samples are not anymore randomly

chosen from the whole dataset, but from a random sub-part of it (the `chunk_aggregate`). This effectively decreases the amount of dispersion (and randomness) of batches, possibly affecting the final model generalization. It is therefore better to consider a higher value for `n_chunks` possibly decreasing `chunk_size`. Furthermore, the values for `chunk_size` and `n_chunks`) should be chosen in conjunction since their product results in the number of samples loaded into memory (RAM) at once for a single worker thread. Increasing too much this number (together with the number of workers used) can potentially saturate main memory (RAM) of the instance used.

Generator parameters optimization To understand the behaviour of the FastTensorDataLoader `alt3` on the target instance and choose the best combination of values for the parameters `chunk_size` and `n_chunks` when using 8 threads (which is double the amount of cores available in the instance used for this project), the code speed was cross evaluated on using powers of 2 for both values. In particular the data loading code was evaluated using values for `chunk_size` and `n_chunks` got from two intervals and the corresponding average *speed* and average *elapsed time* heatmap plots were generated.

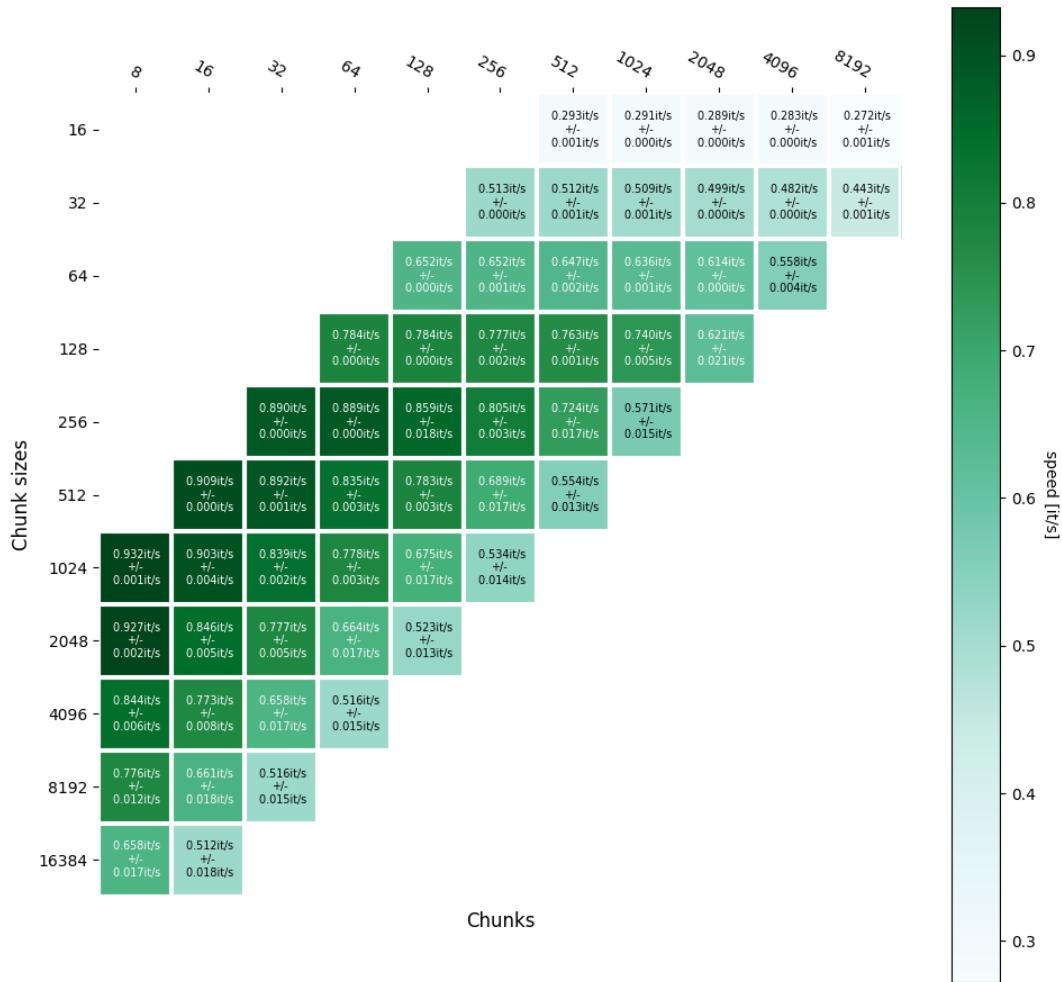


Figure 4.4: Generator Alt.3 Speeds Heatmap, Higher is Better

In practice `chunk_size` and `n_chunks` values were chosen to be the powers of 2 ranging from 2^4 to 2^{14} (included) and from 2^3 to 2^{13} , respectively. Moreover, in order to constrain the evaluation to meaningful values only, two additional parameters have to be set: `min_mul` and `max_mul`. In fact, as previously mentioned, the product of `chunk_size` \times `n_chunks` gives the total number of samples in one `chunk_aggregate` residing in main memory; this quantity must be constrained to a certain range to avoid using too much main memory while being able to retrieve at least one *batch* of data from the `chunk aggregate`. The parameters `min_mul` and `max_mul`

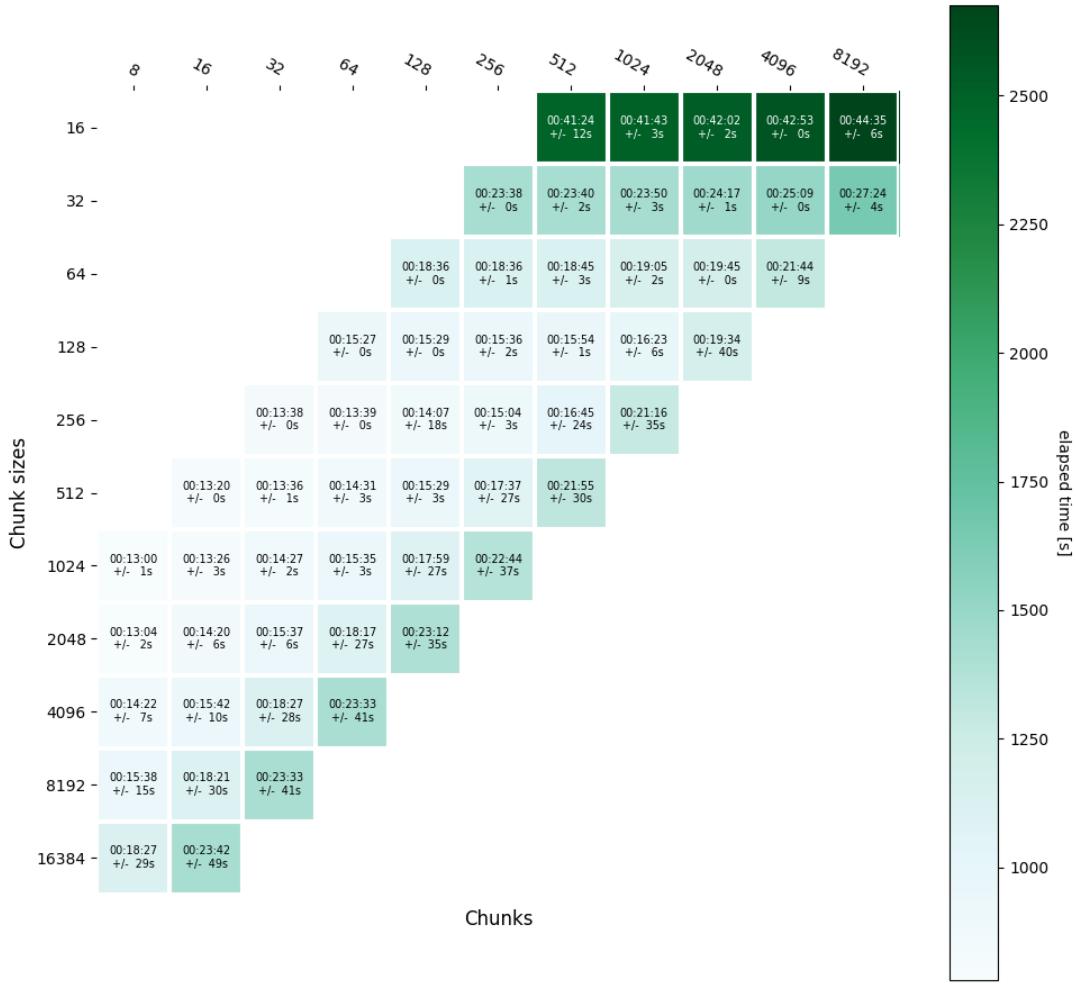


Figure 4.5: Generator Alt.3 Elapsed Times Heatmap, Lower is Better

are used exactly to indicate the minimum and maximum number of batches retrievable from the resulting *chunk_aggregates* and therefore indirectly pose a constrain on the product *chunk_size* \times *n_chunks*.

While effectively evaluating the data loading code (*alt3*) performance the *min_mul* was set to a value of 1, while *max_mul* to 32. Those values were chosen in order to always be able to retrieve at least one *batch* from the *chunk_aggregate* without saturating the RAM available on the instance used for this project (in fact, having for *max_mul* a value greater than $32 \times$ *batch_size* was too much).

Given the resulting heatmaps (4.4, 4.5), *chunk_size* was permanently set to **256** and *n_chunks* to **256** for all the experiments conducted for this project since these values seemed to be a good compromise between data loading speed and samples dispersion.

In any case, with respect to the previous solutions (original Pytorch dataloader and the generators *alt1* and *alt2*) the speed up was huge. This new version took, in fact, approximately 15 minutes, on average, to complete 1 training epoch while the previous solution took, at best, ~ 6 hours. Therefore this was the generator used for loading the dataset in all the experiments done for this project. However, it has to be reiterated that this generator version works by approximating the random sampling step during training which in turn may hinder the model generalization on unseen samples. The results presented in this document may therefore be slightly lower than what could be achieved on a more powerful instance using the original Pytorch dataloader.

4.2 Fresh Dataset

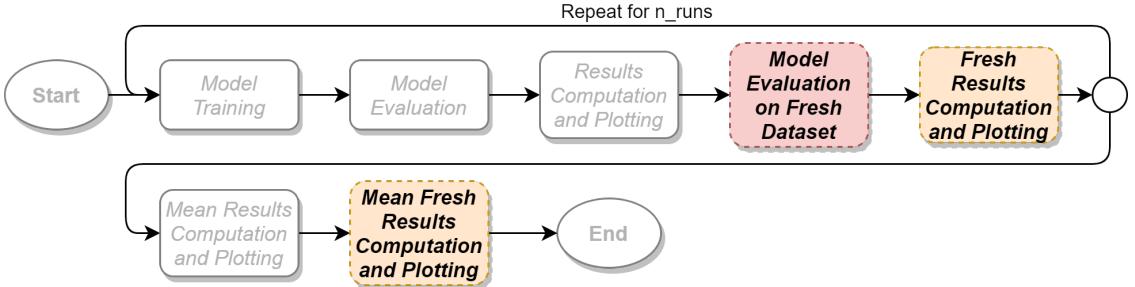


Figure 4.6: Workflow steps using the Fresh dataset

In order to evaluate the learned representation of PE files (embedding) of the proposed model on the sample family prediction task (fig. 4.6), a further dataset, also referred to as *fresh_dataset* throughout this document (and in the code), was created. It consists of a number of sample files' feature vectors along with the corresponding family labels and the sha256 hashes of the original files.

More specifically, a number ' m ' of PE malware families were selected from the list of the most prominent ones present in Italy at the time of writing (as reported by CERT-AGID summary [53]) of which Malware Bazaar provided at least ' x ' sample files. Then, ' x ' sample files per family were downloaded by Malware Bazaar [54] along with their metadata information (label, sha256 hash etc.) and the corresponding numerical feature vectors were extracted using the EMBER feature extraction and numerical feature generation codes [48]. Malware Bazaar, in fact, is a malware sample database maintained by malware analysts which provides examples of malware executables and high-quality manually crafted classifications/descriptions for different malware families and it is therefore a good source for creating the new (fresh) dataset.

A function called *build_fresh_dataset* was specifically implemented with the purpose of creating the *fresh_dataset* given the following arguments:

- a list of malware families (more than ' m ') in order of importance;
- the number ' x ' of samples per family to download;
- the number ' m ' of families to consider.

More specifically, the function downloads ' x ' samples per malware family from Malware Bazaar (if possible, otherwise it skips the malware family and considers the next in order of importance between the ones selected), extract the features from each sample and create the new dataset containing ' $n = x \times m$ ' samples which are then stored on disk as (*numpy*) memory mapped array (as it was done with the pre-processed Sorel20M dataset). Therefore, between the $\geq m$ malware families provided to the function, only the first m for which x PE samples can be retrieved from Malware Bazaar are considered.

For all the model evaluations on the *fresh dataset* done for this project the *fresh dataset* was composed of $x = 100$ samples per family, considering the following $m = 10$ families: *heodo*, *agenttesla*, *formbook*, *masslogger*, *loki*, *avemariarat*, *njrat*, *remcosrat*, *netwire* and *azorult*. Therefore the *fresh dataset* consists of 1000 PE samples.

4.2.1 Model Evaluation with Fresh Dataset

The actual model evaluation on the *fresh dataset* is computed by another function called *evaluate_fresh*. This function evaluates the model learned representation in both the malware family prediction and malware family ranking tasks.

Malware Family Prediction task evaluation For the malware family prediction task there are two similar solutions: one is called *soft-family prediction* evaluation and the other *hard-family prediction* evaluation. In both solutions the learned representation is tested by first selecting n anchor samples per family (for a total of $n \times n_{families} = 100$ anchor samples). The other $n_{samples} - 100 = 900$ non-anchor samples are then assigned, for each family:

- a probability score based on their Euclidean distance or Cosine similarity (depending on the similarity function used by the model) to the most similar anchor sample of that family for the *soft* variant
- a value of 1 if the most similar anchor sample is of that family, 0 otherwise, for the *hard* variant

notice that the *hard* variant does not mean that exactly one single family label is assigned to each test sample. In fact, in case two (or more) anchor samples from different families are at the same distance to the test sample, then a value of 1 is assigned to the test sample for each of those families. In any case, this step results in having for each non-anchor sample a per-family binary prediction output.

Finally, a number of common binary prediction statistics such as *TPR* (True Positive Rate), *Accuracy*, *Recall*, *Precision* and *F1 Score* are computed for each family at different *FPRs* (False Positive Rates) (particularly at $FPR = 1\%$), as well as the corresponding ROC (Receiving Operating Characteristic) curves and AUC-ROC (Area Under Curve) scores.

Malware Family Ranking task evaluation On the other hand, to evaluate the model learned representation on the family ranking task the function first randomly selects ' q ' query samples among the ones in the *fresh dataset*. Then, it computes the similarity (using the model's *get_similarity* function) between each query sample with the other ' $n - 1$ ' samples present in the *fresh dataset* ordering the latter by similarity. At this point, for each query sample there is a ranking of all the other $n - 1$ samples sorted from the most to least similar. Using those rankings the function finally calculates the **MRR** (*Mean Reciprocal Rank*) and **MAP** (*Mean Average Precision*) scores for the model learned representation.

The Mean Reciprocal Rank (**MRR**, 4.2) is the average of the Reciprocal Ranks (**RR**, 4.1) of a series of queries. In particular, the Reciprocal Rank (**RR**) of a query response is the multiplicative inverse of the rank (position in the ranking) of the single sample belonging to the family of interest which classified with the highest rank (it is the closest).

$$RR(i) = \frac{1}{rank_i} \quad (4.1)$$

$$MRR = \frac{1}{q} \sum_{i=1}^q RR(i) = \frac{1}{q} \sum_{i=1}^q \frac{1}{rank_i} \quad (4.2)$$

where $rank_i$ refers to the rank position of the first relevant sample (meaning the first sample with the same family as the current query sample) for the i -th query and q is the number of queries.

The **MAP** score (4.5), on the other hand, takes into account all the samples in the family of interest, with proper weights, instead of considering just the best classified one. Therefore, a model will have a higher **MRR** score if it classifies a single sample of the family of interest higher in the ranking; by contrast, it will have a higher **MAP** score if it classifies all the samples belonging to the family of interest higher in the ranking.

$$P(k) = \frac{|\text{Relevant Samples Retrieved } @k|}{k} \quad (4.3)$$

$$AvgP(k) = \sum_{k=1}^n \frac{P(k) \times rel(k)}{\text{Number of Relevant Samples}} \quad (4.4)$$

$$MAP = \frac{1}{q} \sum_{i=1}^q AvgP(i) \quad (4.5)$$

where $P(k)$ denotes the *precision* at position k in the ranking, $AvgP(k)$ indicates the *average precision* at position k in the ranking, n is the number of samples in the *fresh_dataset*, $rel(k)$ is an indicator function which equals 1 if the sample item at rank k is a relevant sample for the current query, zero otherwise, and finally q is the number of queries. Moreover, a sample is considered relevant if its family is the same as the current query sample one.

The *evaluate_fresh* function provides, together with the **MRR** and **MAP** scores, the complete rankings for the q query samples in a single *json* file. Moreover, for convenience, 5 particular rankings are also saved as *csv* files. These 5 rankings are those that produced:

- the maximum **RR** (Reciprocal Rank);
- the minimum **RR** (Reciprocal Rank);
- the maximum **AP** (Average Precision);
- the minimum **AP** (Average Precision);
- the last ranking is randomly sampled between all the other rankings.

In practice, for all the experiments presented in this document the number of query samples q was set to 100.

Chapter 5

Previous Methods

The state of the art malware detection/description methods presented above (section 3.6), namely *Auxiliary Loss Optimization for Hypothesis Augmentation* (**ALOHA** model, [2]), *Automatic Malware Description via Attribute Tagging and Similarity Embedding* (**Joint Embedding** model, [3]) and *Learning from Context: Exploiting and Interpreting File Path Information for Better Malware Detection* (**File content + contextual file path** (**PE + FP**) model, [45]), represent a good starting point for further research in this topic. Moreover, the models used in those methods can easily be modified to work with Sorel20M dataset, with the exception of the **PE + FP** model [45] since the mentioned dataset provides no file path information (and it is unfeasible to extend it with the needed additional information).

Both the **ALOHA** [2] and **Joint Embedding**, [3] models were therefore implemented in *Python/Pytorch* following the descriptions in the respective papers and having as a starting point the example code of Sorel20M dataset.

5.1 Implementation

Both the implementations developed for this project only loosely followed the original models' structure as described in the respective papers. Therefore here are presented the implementation details.

5.1.1 ALOHA model

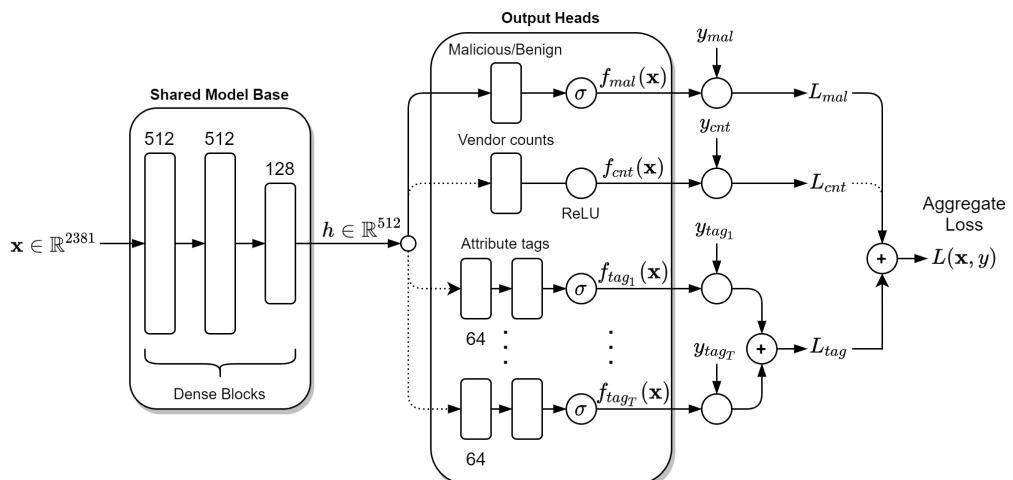


Figure 5.1: ALOHA model implementation architecture

The implementation of ALOHA model (fig. 5.1) consists of a shared model base topology and a number of output heads. The shared base topology is composed of 3 blocks, each consisting of Dropout, a dense layer, layer normalization, and an exponential linear unit (ELU) activation function, with 512, 512, and 128 hidden units respectively. The exact number of layers, along with the layers' dimensions, the dropout probability and the activation function used, however, can be set dynamically in the code at network definition providing a good amount of customizability. The shared model base, given an input sample feature vector \mathbf{x} (of size 2381) got from Sorel20M dataset, outputs an intermediate representation $h = f(\mathbf{x})$ of size 512 which is then used as input to the different parallel output heads.

The output heads used in this implementation are similar to those described in the original ALOHA paper ([2]) with the exclusion of the *Per-Vendor Malicious/Benign* prediction head which could not be reproduced given the lack of the needed information (namely the per-vendor malicious/benign labels) in the Sorel20M dataset. In particular, for each output head (*Malicious/Benign label*, *Vendor Count* and *Malicious tags* prediction heads) an additional block, consisting of one or more dense layers and activation functions was appended to the shared base topology. Moreover, the *Vendor Count* and *Malicious tags* prediction heads are optional, meaning that they can be turned on or off dynamically at network definition.

More specifically, for the *Malicious/Benign* label prediction head, which overall applies function $f_{mal}(x)$ to each sample feature vector \mathbf{x} , a single dense layer followed by the *sigmoid* activation function was used. The loss L_{mal} between the resulting $f_{mal}(x)$ and the ground truth y_{mal} is then computed exactly as described in paper [2] (and as presented in paragraph 3.6.1).

The *Vendor Count* output head, on the other hand, is composed of a single dense layer which outputs $f_{cnt}(x)$. Again, the loss L_{cnt} with respect to the ground truth value y_{cnt} is computed as described by Rudd et al. [2] (and as presented in paragraph 3.6.1).

Finally, the *Malicious (SMART) tags* output head consists of two additional dense layers of size 64 and 1 respectively, followed by a *sigmoid* activation function, per tag. The resulting structure therefore has $T = 11$ parallel paths, one for each tag, that combined compute the functions $f_{tag_i}(\mathbf{x}), \forall i \in 1,..,11$. The aggregate tag loss L_{tag} is then computed, as described in paper [2] (and as presented in paragraph 3.6.1), as the sum of the individual tag losses $L_{tag_i}, \forall i \in 1,..,11$ with respect to each individual ground truth tag y_{tag_i} .

The final aggregate loss $L(\mathbf{x}, y)$ is again computed as the weighted sum of all the output heads losses.

Net Definition

Algorithms 14, 15 and 16 shows some pseudo-code describing how the *ALOHA* model was actually implemented using *Python* language and *Pytorch* library. In particular alg. 14 describes how the model is constructed/initialized: first the model shared base topology is defined as a sequence of layers composed each of a *Pytorch* linear layer, layer normalization, an activation function dynamically chosen at network instantiation among the ones provided by *Pytorch* library, and Dropout. Then the *malware (Malicious/Benign)* head is defined as a single *Pytorch* linear layer followed by a *Sigmoid* activation. The *Vendor Count* head, on the other hand is simply composed of a single linear layer. Finally, the *SMART Tag* prediction head consists of a sequence of 3 linear layers interleaved by *ELU* activations and trailed by a final *Sigmoid* activation function.

Alg. 15, instead, describes how the data (PE file numerical feature vectors of size = 2381) is forwarded through the network to produce the final predictions. In particular, the data first passes through the shared base topology to produce an intermediate representation *base_out*, which is then used as input for all the output heads to produce their predictions.

Finally, in alg. 16 it is shown how the loss between the model predictions and the actual labels gets computed. In particular, the final total loss is the weighted sum of the losses of the different output heads which are enabled for the current run. Both the *Malicious/Benign* and the *SMART tags* prediction heads use the binary cross entropy loss function to compute the loss between their predictions and the ground truth labels. On the other hand, the *Vendor Count* prediction head uses the PoissonNLLLoss (Negative log likelihood loss with Poisson distribution) function.

Algorithm 14 ALOHA Net class, Init

```

1: class NET
2:   function INIT(self, use_malware, use_counts, use_tags, n_tags, feature_dimension,
   layer_sizes, dropout_p, activation_function)
3:     self.use_malware  $\leftarrow$  use_malware
4:     self.use_counts  $\leftarrow$  use_counts
5:     self.use_tags  $\leftarrow$  use_tags
6:     self.n_tags  $\leftarrow$  n_tags
7:     layers  $\leftarrow$  [ ]
8:
9:     if layer_sizes is None then
10:      layer_sizes  $\leftarrow$  [512, 512, 128]
11:    end if
12:
13:    for all i, ls in enumerate(layer_sizes) do
14:      if i = 0 then
15:        layers.append(nn.Linear(feature_dimension, ls))
16:      else
17:        layers.append(nn.Linear(layer_sizes[i - 1], ls))
18:      end if
19:
20:      layers.append(nn.LayerNorm(ls))
21:      layers.append(activation_function())
22:      layers.append(nn.Dropout(dropout_p))
23:    end for
24:
25:    self.model_base  $\leftarrow$  nn.Sequential(*tuple(layers))
26:
27:    self.malware_head  $\leftarrow$  nn.Sequential(nn.Linear(layer_sizes[-1], 1), nn.Sigmoid())
28:    self.count_head  $\leftarrow$  nn.Sequential(nn.Linear(layer_sizes[-1], 1), nn.ReLU())
29:    self.tag_head  $\leftarrow$  nn.Sequential(nn.Linear(layer_sizes[-1], 64), nn.ELU(),
30:                                nn.Linear(64, 64), nn.ELU(),
31:                                nn.Linear(64, n_tags), nn.Sigmoid())
32:  end function

```

Algorithm 15 ALOHA Net class, Forward

```

33:   function FORWARD(self, data)
34:     rv  $\leftarrow$  {}
35:     base_out  $\leftarrow$  self.model_base(data)
36:
37:     if self.use_malware then
38:       rv['malware']  $\leftarrow$  self.malware_head(base_out)
39:     end if
40:
41:     if self.use_counts then
42:       rv['count']  $\leftarrow$  self.count_head(base_out)
43:     end if
44:
45:     if self.use_tags then
46:       rv['tags']  $\leftarrow$  self.tag_head(base_out)
47:     end if
48:
49:   return rv
50: end function

```

Algorithm 16 ALOHA Net class, Compute Loss

```

51:   function COMPUTE LOSS(predictions, labels, loss_wts)
52:     if loss_wts is None then
53:       loss_wts  $\leftarrow \{\text{'malware': } 1.0, \text{'count': } 0.1, \text{'tags': } 1.0\}$ 
54:     end if
55:
56:     loss_dict  $\leftarrow \{\text{'total': } 0.\}$ 
57:     if 'malware'  $\in$  labels then
58:       malware_loss  $\leftarrow$  Binary_cross_entropy(predictions[‘malware’]), labels[‘malware’])
59:
60:       if 'malware'  $\in$  loss_wts then
61:         weight  $\leftarrow$  loss_wts[‘malware’]
62:       else
63:         weight  $\leftarrow 1.0$ 
64:       end if
65:
66:       loss_dict[‘malware’]  $\leftarrow$  malware_loss
67:       loss_dict[‘total’]  $\leftarrow$  loss_dict[‘total’] + malware_loss  $\times$  weight
68:     end if
69:
70:     if 'count'  $\in$  labels then
71:       count_loss  $\leftarrow$  PoissonNLLLoss(predictions[‘count’]), labels[‘count’])
72:
73:       if 'count'  $\in$  loss_wts then
74:         weight  $\leftarrow$  loss_wts[‘count’]
75:       else
76:         weight  $\leftarrow 1.0$ 
77:       end if
78:
79:       loss_dict[‘count’]  $\leftarrow$  count_loss
80:       loss_dict[‘total’]  $\leftarrow$  loss_dict[‘total’] + count_loss  $\times$  weight
81:     end if
82:
83:     if 'tags'  $\in$  labels then
84:       tags_loss  $\leftarrow$  Binary_cross_entropy(predictions[‘tags’], labels[‘tags’])
85:
86:       if 'tags'  $\in$  loss_wts then
87:         weight  $\leftarrow$  loss_wts[‘tags’]
88:       else
89:         weight  $\leftarrow 1.0$ 
90:       end if
91:
92:       loss_dict[‘tags’]  $\leftarrow$  tags_loss
93:       loss_dict[‘total’]  $\leftarrow$  loss_dict[‘total’] + tags_loss  $\times$  weight
94:     end if
95:
96:     return loss_dict
97:   end function
98: end class

```

5.1.2 Joint Embedding

The implementation of the Joint Embedding model (fig. 5.2) is composed of 3 main parts: the *Model Base/PE Embedding* topology, the *Tag Embedding* and the *Prediction Layer*.

The model *base topology* consists of a series of 3 dense blocks, each composed by dropout, a linear layer, layer normalization and ELU activation function, of output sizes 512, 512 and 128

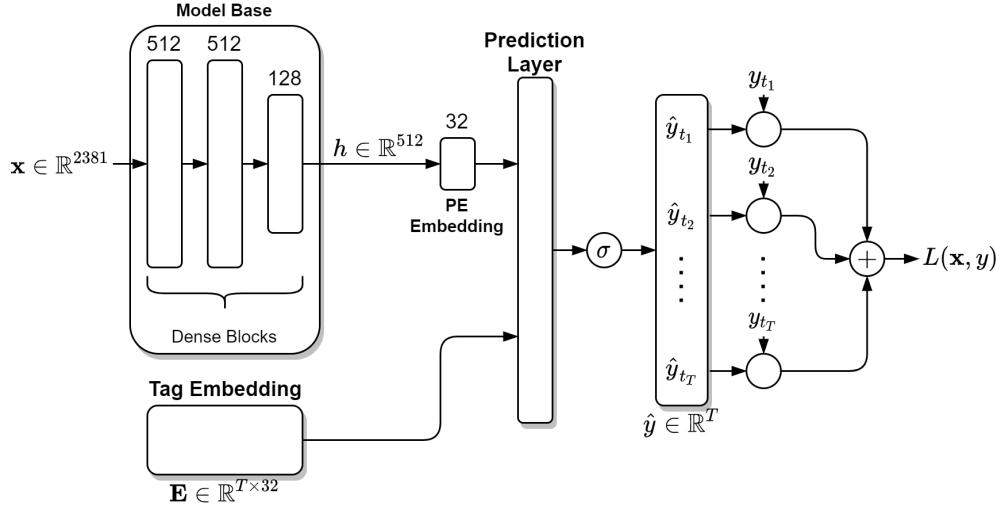


Figure 5.2: Joint Embedding model implementation architecture

respectively. Similarly to the ALOHA implementation, at network definition it is again possible to dynamically set the exact number of layers, together with the linear layers' sizes, the dropout probability and the activation used for the base topology. Anyway, the base PE embedding topology outputs an intermediate representation h of size 512 given an input sample feature vector \mathbf{x} (of size 2381). This intermediate representation is then used as input to a further linear layer with output size equal to the chosen **Joint Embedding** size (set to 32 throughout all the experiments). The output of this further layer is the one used as input to the *Prediction Layer* and corresponds to the representation of the input file features x in the Joint Embedding space.

On the other hand, the *Tag Embedding* matrix $\mathbf{E} \in \mathbb{R}^{T \times 32}$, which is the same as the original implementation [3], maps each tag t_n to the corresponding representation in the Joint Embedding space and is therefore used as second input to the *Prediction Layer*.

Finally, the *Prediction Layer* combines the sample (PE) and tags embeddings producing a similarity score matrix that is run through a *sigmoid* activation function (by value) resulting in the probabilities of each tag t being associated to sample \mathbf{x} . The function used to compute the similarity score between the two embedding vectors can be chosen dynamically at network definition between: the *dot product* (eq. 5.1), the *cosine similarity* (eq. 5.2) and the inverse of the *Euclidean distance*. In particular, there are 3 versions of inverted *Euclidean distance* that can be selected at network definition: *exp* (eq. 5.3), *inv* (eq. 5.4) or *inv-pow* (eq. 5.5).

$$\text{dot product}(\mathbf{a}, \mathbf{b}) = \mathbf{a} \cdot \mathbf{b} = \sum_{i=1}^D a_i b_i \quad (5.1)$$

$$\text{cosine similarity}(\mathbf{a}, \mathbf{b}) = \frac{\mathbf{a} \cdot \mathbf{b}}{\|\mathbf{a}\| \|\mathbf{b}\|} = \frac{\sum_{i=1}^D a_i b_i}{\sqrt{\sum_{i=1}^D a_i^2} \sqrt{\sum_{i=1}^D b_i^2}} \quad (5.2)$$

$$\text{euclidean similarity exp}(\mathbf{a}, \mathbf{b}) = e^{(-d(\mathbf{a}, \mathbf{b}))} \quad (5.3)$$

$$\text{euclidean similarity inv}(\mathbf{a}, \mathbf{b}) = \frac{1}{1 + \frac{d(\mathbf{a}, \mathbf{b})}{\alpha}} \quad (5.4)$$

$$\text{euclidean similarity inv-pow}(\mathbf{a}, \mathbf{b}) = \frac{1}{1 + \frac{d(\mathbf{a}, \mathbf{b})^2}{\alpha}} \quad (5.5)$$

where $d(\mathbf{a}, \mathbf{b})$ is the Euclidean distance between points \mathbf{a} and \mathbf{b} as defined in eq. 5.6 and α is a multiplicative factor that can be arbitrarily set at network definition when using one of the

inverted *Euclidean distance* functions. In the final Joint Embedding model implementation α is set to 1.0.

$$d(\mathbf{a}, \mathbf{b}) = \sqrt{\sum_{i=1}^D (a_i - b_i)^2} \quad (5.6)$$

The tags predicted probabilities outputted by the *Prediction Layer* are then used to compute the respective losses with respect to the ground truth tags $y_{t,i}, \forall i \in 1,..,11$, exactly as described in paper [3], which are combined together to form the final loss $L(\mathbf{x}, y)$.

Net Definition

Algorithm 17 Joint Embedding Net class, Init

```

1: class NET
2:   function INIT(self, use_malware, use_counts, use_tags, n_tags, feature_dimension, embedding_dimension, embeddingmax_embedding_norm, layer_sizes, dropout_p, activation_function)
3:     self.n_tags  $\leftarrow$  n_tags
4:     self.embedding_dimension  $\leftarrow$  embedding_dimension
5:     layers  $\leftarrow$  []
6:
7:     if layer_sizes is None then
8:       layer_sizes  $\leftarrow$  [512, 512, 128]
9:     end if
10:
11:    for all i, ls in enumerate(layer_sizes) do
12:      if i = 0 then
13:        layers.append(nn.Linear(feature_dimension, ls))
14:      else
15:        layers.append(nn.Linear(layer_sizes[i - 1], ls))
16:      end if
17:
18:      layers.append(nn.LayerNorm(ls))
19:      layers.append(activation_function())
20:      layers.append(nn.Dropout(dropout_p))
21:    end for
22:
23:    self.model_base  $\leftarrow$  nn.Sequential(*tuple(layers))
24:    self.sigmoid  $\leftarrow$  nn.Sigmoid()
25:
26:    self.pe_embedding  $\leftarrow$  nn.Sequential(nn.Linear(layer_sizes[-1],
27:                                              self.embedding_dimension))
28:    self.tags_embedding  $\leftarrow$  nn.Embedding(self.n_tags, self.embedding_dimension,
29:                                         max_norm=max_embedding_norm)
30:  end function

```

Algorithm 17 (*Joint Embedding* model init function) shows how the *Joint Embedding* model was implemented in *Python/Pytorch* code. More specifically, it describes how the model network is initialized/defined: first a **base topology** is defined similarly to the one used in the **ALOHA** model implementation previously described (a series of linear layers interleaved by layer normalization, an activation function and dropout). The **PE embedding** sub-network is, on the other hand, defined as a simple linear layer with output size equal to the chosen *embedding_dimension*. Finally, the **Tags embedding** is defined as an embedding matrix of size (*n_tags*, *embedding_dimension*), constraining the maximum norm of each tag embedding to *max_embedding_norm* (set to 1 throughout the experiments).

Algorithm 18 Joint Embedding Net class, Forward (with *Dot Product*)

```

31:   function FORWARD(self, data)
32:     rv  $\leftarrow \{\}$ 
33:     base_out  $\leftarrow \text{self.model\_base}(\text{data})$ 
34:
35:     pe_embedding  $\leftarrow \text{self.pe_embedding}(\text{base\_out})$ 
36:     tags_embedding  $\leftarrow \text{self.tags_embedding}(\text{Dataset.encoded\_tags})$ 
37:
38:     similarity_scores  $\leftarrow \text{torch.matmul}(\text{pe\_embedding}, \text{tags\_embedding.T})$ 
39:     rv[‘similarity’]  $\leftarrow \text{similarity\_scores}$ 
40:     rv[‘probability’]  $\leftarrow \text{self.sigmoid}(\text{similarity\_scores})$ 
41:     return rv
42:   end function

```

Algorithm 19 Joint Embedding Net class, Forward (with *Cosine similarity*)

```

43:   function FORWARD(self, data)
44:     ...
45:     similarity_scores  $\leftarrow \text{torch.div}(\text{torch.add}(\text{Cosine\_similarity}(\text{pe\_embedding},$ 
46:                               tags_embedding, dim=2), 1.0), 2.0)
47:     rv[‘similarity’]  $\leftarrow \text{similarity\_scores}$ 
48:     rv[‘probability’]  $\leftarrow \text{similarity\_scores}$ 
49:     return rv
50:   end function

```

Algorithm 20 Joint Embedding Net class, Forward (with *Inverted Pairwise Euclidean Distance*)

```

51:   function FORWARD(self, data)
52:     ...
53:     distances  $\leftarrow \text{torch.cdist}(\text{pe\_embedding}, \text{tags\_embedding}, \text{p}=2.0)$ 
54:     similarity_scores  $\leftarrow \text{distance\_to\_similarity}(\text{distances}, \text{a}=\text{sim\_function\_a},$ 
55:                               function=sim_function)
56:     rv[‘similarity’]  $\leftarrow \text{similarity\_scores}$ 
57:     rv[‘probability’]  $\leftarrow \text{similarity\_scores}$ 
58:     return rv
59:   end function

```

Algorithm 21 Joint Embedding Net class, Get_Embedding

```

60:   function GET_EMBEDDING(self, data)
61:     rv  $\leftarrow \{\}$ 
62:     base_out  $\leftarrow \text{self.model\_base}(\text{data})$ 
63:     pe_embedding  $\leftarrow \text{self.pe_embedding}(\text{base\_out})$ 
64:     rv[‘embedding’]  $\leftarrow \text{pe\_embedding}$ 
65:     return rv
66:   end function

```

Algorithm 22 Joint Embedding Net class, Get_Similarity (with *Dot Product*)

```

67:   function GET_SIMILARITY(self, first_embedding, second_embedding)
68:     similarity_scores  $\leftarrow \text{torch.matmul}(\text{first\_embedding}, \text{second\_embedding.T})$ 
69:     return {‘similarity’: similarity_scores, ‘probability’: self.sigmoid(similarity_scores)}
70:   end function

```

Alg. 18 (**Joint Embedding model forward function**), instead, describes how the input data (the PE files' numerical feature vectors) is forwarded through the network to produce the final tag predictions. In particular, the input data is first fed into the model base topology to produce an intermediate representation *base_out*. This transient output is then used as input

Algorithm 23 Joint Embedding Net class, Get_Similarity (with *Cosine similarity*)

```

71:   function GET_SIMILARITY(self, first_embedding, second_embedding)
72:     similarity_scores  $\leftarrow \text{torch.div}(\text{torch.add}(\text{F.cosine_similarity}(\text{first\_embedding},
73:                               second_embedding), 1.0), 2.0)
74:     return {’similarity’: similarity_scores, ’probability’: similarity_scores}
75:   end function$ 
```

Algorithm 24 Joint Embedding Net class, Get_Similarity (with *Inverted Pairwise Euclidean Distance*)

```

76:   function GET_SIMILARITY(self, first_embedding, second_embedding)
77:     distances  $\leftarrow \text{torch.cdist}(\text{first\_embedding}, \text{second\_embedding}, \text{p}=2.0)$ 
78:     similarity_scores  $\leftarrow \text{distance\_to\_similarity}(\text{distances}, \text{a}=\text{self.embedding\_dimension},
79:                               function=sim_function)
80:     return {’similarity’: similarity_scores, ’probability’: similarity_scores}
81:   end function$ 
```

Algorithm 25 Joint Embedding Net class, Compute Loss

```

82:   function COMPUTE_LOSS(predictions, labels, loss_wts)
83:     loss_dict  $\leftarrow \{ \text{’total’}: 0. \}$ 
84:
85:     similarity_loss  $\leftarrow \text{Binary\_cross\_entropy}(\text{predictions}[’probability’],
86:                               labels[’tags’],
87:                               reduction=’none’}).\text{sum}(\text{dim}=1).\text{mean}(\text{dim}=0)$ 
88:
89:     loss_dict[’jointEmbedding’]  $\leftarrow \text{similarity\_loss}$ 
90:     loss_dict[’total’]  $\leftarrow \text{similarity\_loss}$ 
91:
92:     return loss_dict
93:   end function
94: end class
```

Algorithm 26 Distance-to-Similarity function

```

1: function DISTANCE_TO_SIMILARITY(distance, a, function)
2:   if function = ’exp’ then
3:     similarity  $\leftarrow \text{torch.exp}(\text{torch.div}(\text{distances}, -a))$ 
4:   else if function = ’inv’ then
5:     similarity  $\leftarrow \text{torch.pow}(\text{torch.add}(\text{torch.div}(\text{distances}, a), 1.0), -1.0)$ 
6:   else if function = ’inv_pow’ then
7:     similarity  $\leftarrow \text{torch.pow}(\text{torch.add}(\text{torch.div}(\text{torch.pow}(\text{distances}, 2.0), a),
8:                               1.0), -1.0)$ 
9:   else
10:    raise Exception
11:   end if
12:
13:   return similarity
14: end function
```

for the ***PE Embedding*** topology to produce the samples’ representation in the joint embedding space (*pe_embedding*). On the other hand, the tags embeddings (one per tag) are extracted from the *tags_embedding* matrix. The **similarity scores** between the samples’ and tags’ embeddings is computed via a simple ***dot product*** (implemented through a matrix multiplication) between the embeddings. Finally the tags probabilities are obtained by mapping the similarity scores to the range [0,1] through the *sigmoid* function.

In alg. 19 (**Joint Embedding** model forward function with *cosine similarity*) is presented an alternative to the aforementioned **Model forward function** that uses the *Cosine similarity* instead of the *dot product* to compute the similarity scores. For the most part this version is equal to the previous one but for how the similarity scores are computed. In fact, in this case they correspond to the *Cosine similarity* between the *pe_embedding* and *tags_embedding*, which is a value in the range $[-1, 1]$, mapped to values in the $[0, 1]$ range. Moreover, in this case, since the similarity scores are already in the $[0, 1]$ range, the tags probabilities coincide with the similarity scores.

A further alternative to the **Model forward function** which uses one of the *Inverted Pairwise Euclidean Distances* proposed in sec. 5.1.2 instead of the *dot product*, is presented in alg. 20 (**Joint Embedding** model forward function with *Inverted Pairwise Euclidean Distance*). Again, this alternative is pretty much equal to the original forward function (alg. 18). In this case, however, the similarity scores are produced by inverting, through the use of the specially crafted *distance_to_similarity* function presented in alg. 26 (which simply applies the selected distance inversion formula among the ones mentioned in sec. 5.1.2), the euclidean distances between the samples' and the tags' embeddings in the joint latent space.

Alg. 21 (**JointEmbedding** model get embedding function) presents the *get_embedding* function which is used to extract the PE embedding representation from input data. This is particularly useful when performing different kinds of model performance evaluations together with the *get_similarity function*, which instead returns the similarity between two different input embeddings. Of course, depending on the similarity function selected at model definition the *get_similarity* function has to be different. In particular, if the *dot product* is selected the function, shown in algorithm 22 (**JointEmbedding** model get similarity function with **Dot Product**), simply computes the similarity scores as the dot product between the two input embeddings and then defines the probabilities by further applying the sigmoid activation function on those scores. On the other hand, if the *cosine similarity* is selected, the function (23 - **JointEmbedding** model get similarity function with *Cosine Similarity*) assigns to both similarity scores and resulting probabilities the normalized (to the $[0, 1]$ range) cosine similarity between the two embeddings. Finally, in case the **Inverted Pairwise Euclidean Distance** is selected, the function (24 - **JointEmbedding** model get similarity function with *Inverted Pairwise Euclidean Distance*) first computes the euclidean distances between the two input embeddings and then assigns to both similarity scores and resulting probabilities the inversion of such distances computed through the *distance_to_similarity* function.

Finally, in algorithm 25 is described how the loss which guides the training procedure is computed. In particular, it corresponds to the **binary cross entropy loss** between the predicted tag probabilities and the actual tag labels, as depicted by the formulas mentioned in par. 3.6.2.

5.2 Experiments

Both the models were trained (and validated) each for 10 *epochs* on the first 6M training and 1.153.846 validation samples of the pre-processed Sorel20M dataset (loaded with generator *alt3*) using different number of layers, number of nodes per layer (layer sizes), activation functions and dropout value for the model base topology, different optimizers (**Adam** or **SGD**), learning rates (LR), momentum, weight decay and loss weights. Moreover, in the case of the **Joint Embedding** model also different similarity score functions such as the cosine similarity and the inverse of the Euclidean distance were tested instead of the dot product. The resulting models were then evaluated on the first 1.846.154 test samples of Sorel20M dataset looking for their tag and/or malicious/benign label (present only in the ALOHA model implementation) prediction *TPR*, *accuracy*, *recall*, *precision*, *f1* score, *jaccard similarity* and *mean-per-sample-accuracy* at a FPR of 1%, as well as the resulting tag and/or malicious/benign *ROC* curves and *AUC* scores.

The final **ALOHA** architecture, which had a base topology consisting of 3 layers of sizes 512, 512 and 128 respectively, with dropout probability $p_d = 5\%$ and **ELU** activation function, in particular, was trained with a learning rate $LR = 10^{-3}$, **Adam** optimizer, no momentum nor weight decay and loss weights set to 1.0 for the malicious/benign head, 0.1 for the count head and 0.1 for the tag head (as described in the original paper [2]).

On the other hand, the base topology of the final **Joint Embedding** architecture consisted of 3 layers of sizes 512, 512, and 128 respectively, with dropout probability $p_d = 5\%$ and **ELU** activation function. The overall model was trained with a learning rate $LR = 10^{-3}$, **Adam** optimizer, no momentum nor weight decay (the loss weights are not used since there is a single output head). Moreover, the final model uses the dot product as similarity score between tags' and samples' embeddings.

The **ALOHA** and **Joint Embedding** model implementations results are reported in chapter 7.

5.3 Training and Evaluation algorithms

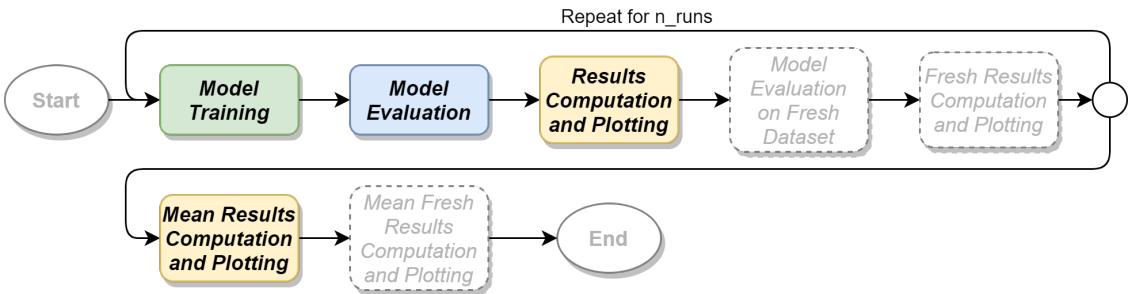


Figure 5.3: Workflow Model Training and Evaluation steps

Both the **ALOHA** and the **Joint Embedding** model implementations (and also the **Proposed Model** that will be defined in next chapter) are trained and evaluated following the procedures depicted by algorithms 27 (**Train network function**) and 28 (**Evaluate network function**).

5.3.1 Training

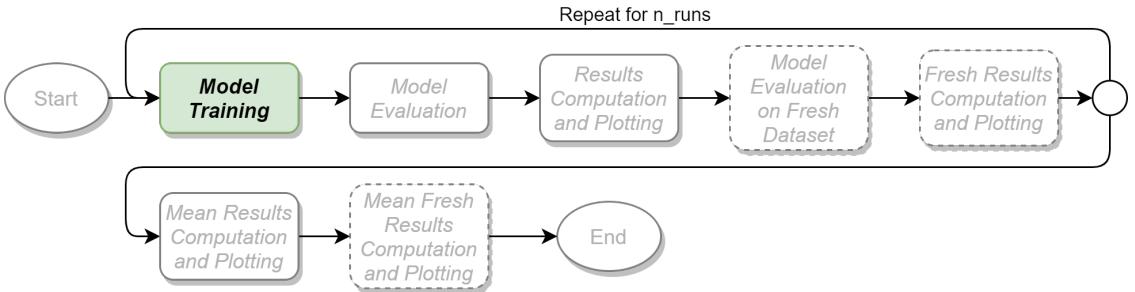


Figure 5.4: Workflow Model Training step

In particular, as shown in alg. 27, the training function first imports the chosen Net and Dataset generator (dataloader) implementations and some run variables situated in a *config* file, depending on the network and generator types selected by the user. Next, the model itself is defined calling the *init* member function of the previously imported *Net* class with all needed parameters (such as the number of tags, the input feature dimension, the layer sizes, the dropout probability, etc.). Then, the selected optimizer (**Adam** or **SGD**) is instantiated given the model parameters, chosen learning rate and weight decay. This is followed by the training and validation generators definition. Moreover, if the model has already been trained in a previous run and its state parameters checkpoints are available, both the model and optimizer states are restored. Next, the model is trained for '*epochs*' epochs. Each epoch the model is first set in training mode, then for all the mini-batches randomly sampled from the training generator the input features are forwarded through the network to get the output predictions after having re-set the optimizer previously computed gradients. The loss for the current mini batch is then computed and logged.

Algorithm 27 Train Network function

```
1: function TRAIN_NETWORK(ds_path, net_type, gen_type, batch_size, epochs,
   use_malicious_labels, use_count_labels, use_tag_labels, feature_dimension, workers, ...)
2:   Net, Dataset, get_generator, device, run_params  $\leftarrow$  import_modules(net_type, gen_type)
3:
4:   model  $\leftarrow$  Net(use_malicious_labels, use_count_labels, use_tag_labels,
5:     len(Dataset.tags), feature_dimension, run_params[‘layer_sizes’],
6:     run_params[‘dropout_p’], run_params[‘activation_function’])
7:   opt  $\leftarrow$  run_params[‘optimizer’]({model.parameters}, run_params[‘lr’],
8:     run_params[‘weight_decay’])
9:
10:  generator  $\leftarrow$  get_generator(ds_path, batch_size, ‘train’, workers, training_n_samples,
11:    use_malicious_labels, use_count_labels, use_tag_labels)
12:  val_generator  $\leftarrow$  get_generator(ds_path, batch_size, ‘validation’, workers,
13:    validation_n_samples, use_malicious_labels,
14:    use_count_labels, use_tag_labels)
15:  ...
16:  model.load(model_checkpoint_path)
17:  opt  $\leftarrow$  load_opt_state(opt, opt_checkpoint_path)
18:  for all epoch  $\in$  range(epochs) do
19:    model.train()
20:    for all i, (features, labels)  $\in$  enumerate(generator) do
21:      opt.zero_grad()
22:      out  $\leftarrow$  model(features)
23:      loss_dict  $\leftarrow$  model.compute_loss(out, labels, run_params[‘loss_wts’])
24:      loss  $\leftarrow$  loss_dict[‘total’]
25:      log(loss_dict)
26:      loss.backward()
27:      opt.step()
28:    end for
29:
30:    model.eval()
31:    for all i, (features, labels)  $\in$  enumerate(val_generator) do
32:      torch.no_gradout  $\leftarrow$  model(features)
34:      loss_dict  $\leftarrow$  model.compute_loss(out, labels)
35:      log(loss_dict)
36:    end for
37:
38:    model.save(epoch)
39:    save_opt_state(opt, epoch)
40:  end for
41: end function
```

Finally the model parameters are update by computing the loss gradients and back-propagating them through the network. Moreover, after each epoch training the model is also evaluated on samples from the validation dataset provided by the validation generator. In particular, the model is first set in *evaluation* mode, then for all the mini-batches of data from the validation generator the input features are passed through the net to produce the output predictions which are in turn used to compute the loss term. The aggregate loss, in this case, is simply logged (no gradients are computed). Finally, at the end of each epoch the model and optimizer state parameters are saved to file.

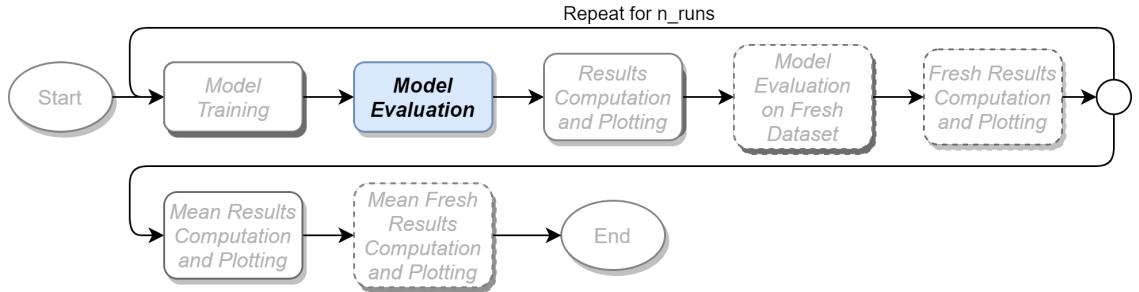


Figure 5.5: Workflow Model Evaluation step

Algorithm 28 Evaluate Network function

```

1: function TRAIN_NETWORK(ds_path, checkpoint_file, net_type, gen_type, batch_size, evaluate_malware, evaluate_count, evaluate_tags, ...)
2:   Net, Dataset, get_generator, device, run_params  $\leftarrow$  import_modules(net_type, gen_type)
3:
4:   model  $\leftarrow$  Net(evaluate_malware, evaluate_count, evaluate_tags,
5:                      len(Dataset.tags), feature_dimension, run_params[‘layer_sizes’],
6:                      run_params[‘dropout_p’], run_params[‘activation_function’])
7:
8:   model.load_state_dict(checkpoint_file)
9:   model.eval()
10:
11:  generator  $\leftarrow$  get_generator(ds_path, batch_size, ‘test’, workers, test_n_samples,
12:                           use_malicious_labels, use_count_labels,
13:                           use_tag_labels, return_shas=True)
14: ...
15: for all shas, features, labels  $\in$  generator do
16:   predictions  $\leftarrow$  model(features)
17:   save_to_file(labels, predictions)
18: end for
19: end function

```

5.3.2 Evaluation

The evaluation function, presented in 28, begins in a similar way as the training function by importing the Net and Dataset generator along with some run parameters, by calling the model constructor function (*init*) with all the needed parameters and by loading the model checkpoint chosen by the user. This function, however, continues by setting the model in *evaluation* mode and defining the generator for the test dataset. Then, for all the mini-batches of data from the test set the samples input features are passed to the network and the corresponding predictions are produced and saved to file.

5.3.3 Results Computation and plotting

The actual binary cross entropy statistics (such as *TPR*, *Accuracy*, *Recall*, etc.. at different FPRs) and the resulting AUC-ROC score and ROC curves are computed by a distinct set of functions not presented in this document since they simply implement and apply the corresponding mathematical formulas and save the ROC curves to file.

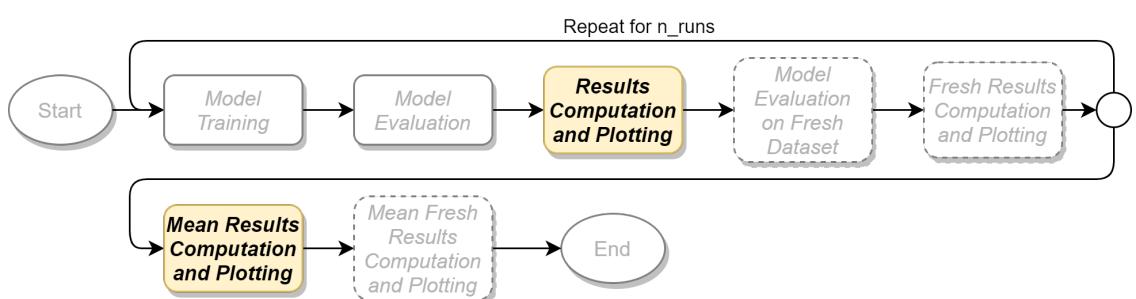


Figure 5.6: Workflow Results Computation and Plotting step

Chapter 6

Proposed Method

ALOHA model is based on the idea of enhancing the model performance and generalization on the task of *Malicious/Benign* label prediction with respect to a more simple single-task model by exploiting additional tasks (*Vendor Count* prediction, *SMART Tags* prediction, etc.) which are jointly optimized in a Multi-objective/multi-task learning fashion. Multi-objective learning has in fact proven to be beneficial for the model generalization and overall has a regularization effect providing better training curves and arguably better results.

On the other hand, the **Joint Embedding** model was designed with the idea of constructing an embedding space in which both the samples and tag labels can be mapped to, such that each sample ends up close to the tags it is associated to (and possibly far from the other unrelated tags). The *Joint Embedding* model is trained by optimizing its parameters on the single task of tag prediction by means of tags' and samples' embeddings similarity, with no additional tasks.

The natural follow-up to those models is a fusion between their two base ideas ultimately creating a model with the main objective of creating a joint embedding space, as done by the **Joint Embedding** model, while exploiting Multi-objective learning, as done by the **ALOHA** model. This model, which is called **Proposed Model** throughout this document, could potentially perform better in the SMART tag prediction task than the **Joint Embedding** model implementation since during the learning procedure its parameters' updates are 'guided' by the inclusion of additional targets. Moreover, it could also perform better than the **ALOHA** model implementation in the *Malicious/Benign* label prediction task since one of its additional tasks - the SMART tag prediction through similarity embedding - is harder than the original one - multi-label binary classification with one head for each tag. Harder complementary tasks are often used when exploiting *multi-task* learning because they help in finding a more informed latent representation of samples which in turn leads to better performances during testing/deployment.

Moreover, the fact that the learned embedding should pull samples close to their associated tags (and far from the others) also means that samples which are similar or exhibit similar capabilities (sharing the same tags) should be close to each other in the latent space. In turn, the learned representation of PE files in the embedding space, along with the distance function used (Euclidean distance) in this space, enables the computation of file-to-file similarities. This makes it possible to measure how similar two malware samples are in terms of their capabilities. The **Joint Embedding** and **Proposed Model** approaches are therefore interesting from the perspective of predicting the specific malware family of PE files based on their distance with respect to known malware samples. Notice, however, that both the **Joint Embedding** and the **Proposed Model** network are trained with a different objective - that of predicting the SMART tags associated to each malware sample (and also the *malicious/benign* label and *vendor count* for the proposed model). It is not uncommon in the ML field to train a model on a specific task while then testing (or deploying) it on a different, but most probably related, task with little degradation in model performance. In particular this technique is generally referred to *Transfer Learning*. Moreover, the models are trained on a dataset of malware samples which intrinsically belong to different malware families (although the effective family labels are not provided in the dataset), therefore the information of each sample's family is implicitly present.

Therefore the **Proposed Model** and the **Joint Embedding** model implementation performances are not only evaluated on the tasks they were trained for (*Malicious/Benign* label and *Vendor Count* prediction tasks and/or *SMART tag* prediction task), but also on the **Malware Family prediction** task and on the **Malware Family Ranking** task. Again, for the reasons mentioned above, the **Proposed Model** should perform better than the **Joint Embedding** model implementation also in these last tasks.

Then, if the **Proposed Model** performs well enough on the malware family *prediction* and *ranking* tasks, it could be possible to use its learned embedding space to define implicit signatures - the corresponding representation in this space, a.k.a. the embedding - of malware samples belonging to a specific malware family. These signatures could then be used to craft Yara Rules - with the use of a specially crafted custom Yara module that dynamically extracts the embedding representation for new PE files and compares them to the known signatures - to detect similar samples as belonging to the specific family represented by each rule. Yara Rules created in this way could potentially inherit the good properties of ML methods such as being able to detect not only known malware strains but also novel variants, being less susceptible to obfuscation attempts, etc. while being compatible with the highly expressive approach that is Yara.

6.1 Implementation

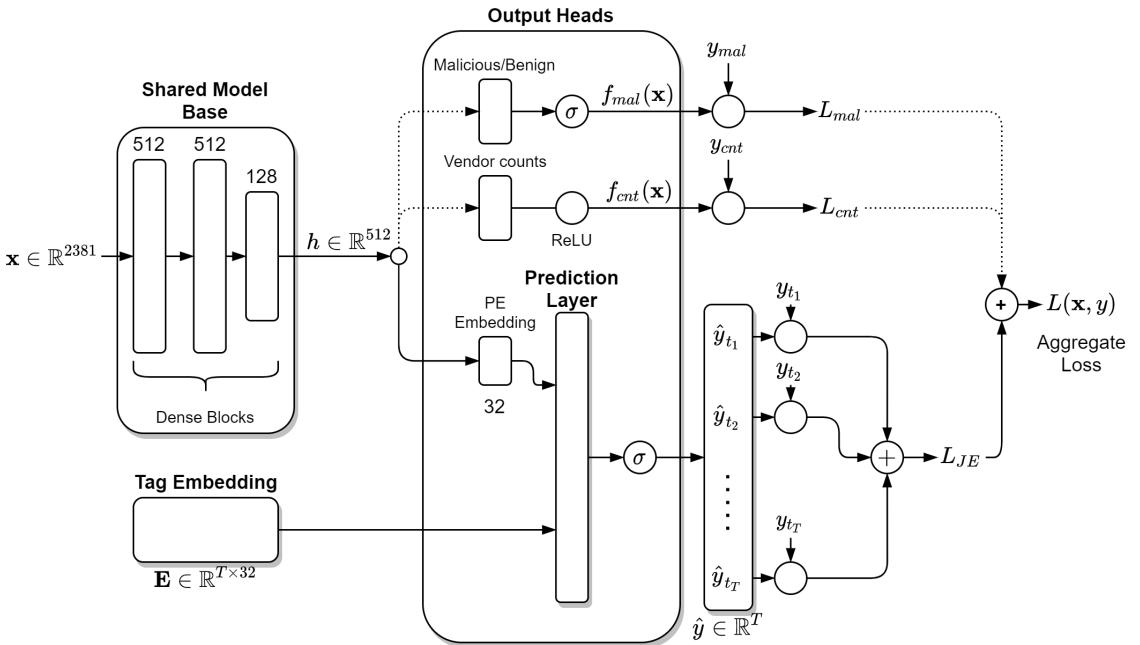


Figure 6.1: Proposed Model architecture

The proposed model implementation (fig. 6.1) consists of 3 main topologies: the *Shared Model Base* sub-network, the *Tag Embedding* and the multiple parallel *Output Heads*.

The model *shared base* topology is composed of 3 sequential dense blocks of output sizes 512, 512 and 128 respectively. These dense blocks have the same form as those used in the *ALOHA* and *Joint Embedding* models implementations described above (in 5.1.1 and 5.1.2 respectively): in particular a linear layer with dropout is used, followed by layer normalization and ELU activation function. Again at network definition it is possible to dynamically set the parameters for this base topology such as the exact number of layers, the linear layer's sizes, the dropout probability p_d and the non-linear activation function used. This model *shared base* topology transforms a given input sample feature vector \mathbf{x} (of size 2381) into its intermediate representation h of size 512 which is then used to the parallel output heads.

The *Tag Embedding* matrix $\mathbf{E} \in \mathbb{R}^{T \times 32}$, which maps the tags to their corresponding representation in the Joint Embedding space, is the same as the one used in the *Joint Embedding* model implementation (5.1.2) and is therefore also equal to the one used in the original implementation [3]. This matrix is used as second input for the *Prediction Layer* in the tag output head.

Similarly to the *ALOHA* model implementation (5.1.1), this model uses multiple parallel *output heads*, each relative to a different task, whose values are jointly optimized during training by back-propagating the gradients computed from an aggregate loss. In particular, the tasks used in this case are: the *Malicious/Benign label* prediction (optional), the *Vendor Count* estimation (optional) and the *SMART Tag* prediction based on the similarity between samples' and tags' embeddings. The first two output heads (*Malicious/Benign label* prediction head and *Vendor Count* estimation head) are the same ones used in the *ALOHA* model implementation (5.1.1) and can optionally be turned on or off at network definition.

In particular, the *Malicious/Benign* label prediction head is composed of a single dense layer followed by the *sigmoid* non-linearity. The loss between the malware/benign output of the network $f_{mal}(x)$ and the ground truth label y_{mal} is computed using the binary cross-entropy loss function.

The *Vendor Count* prediction head, instead, consists of a single linear layer whose parameters are updated in order to optimize the negative log-likelihood loss, as described in 3.6.1 (and [2]), between the predicted and the ground truth counts.

The *Tag* prediction head, on the other hand, is always used and consists of the same components used in the *Joint Embedding* model implementation (5.1.2): the linear layer mapping a sample intermediate representation h to its *PE Embedding* in the **Joint Embedding space** and the *Prediction Layer*, which produces a similarity score matrix between samples' and tags' embeddings. The resulting matrix is then run through a *sigmoid* non-linearity (by value) to produce the tags probabilities for sample x . The similarity function used in this model can be chosen to be one of the following at network definition: *dot product* (5.1), *cosine similarity* (5.2) or one of the *inverse Euclidean distances* (*exp* (5.3), *inv* (5.4) or *inv_pow* (5.5)). The *Tag prediction head* loss is computed as the sum over all tags of the individual cross entropy losses between the predicted tags and the ground truth ones, as described at 3.6.2 (and [3]).

Finally the individual head losses are aggregated together by computing their weighted sum. Different weights can be assigned to the distinct heads in order to tune their importance in the final loss term in turn affecting how the model's parameters are updated during back-propagation.

Net Definition

As presented by algorithm 29 (*Proposed Model init function*), the **Proposed Model** is initialized first by defining the *model shared topology* (*model_base*) as a sequence of blocks composed each of a linear layer, layer normalization, an activation function and dropout, exactly as previously done for the *ALOHA* and *Joint Embedding* model implementations. Next, similarly to the *Joint Embedding* implementation, the *pe_embedding* topology and the *tags_embedding* matrix are defined: the first as a single linear layer with output size equal to the *embedding_dimension* (= 32), and the latter as an embedding matrix of size (*n_tags*, *embedding_dimension*). The most important novelty compared to the *Joint Embedding* model is the presence of also the *Malicious/Benign* label and *Vendor Count* prediction heads from the *ALOHA* model implementation. In particular, the *Malicious/Benign* (*malware_head*) is defined as a linear layer of output size 1 immediately followed by a *sigmoid* non linearity. The *Vendor Count* head (*count_head*), on the other hand, is set to be a simple linear layer.

The input data (PE files' numerical feature vectors) are forwarded through the model as shown in alg. 30 (*Proposed Model forward function (with dot product)*). In particular, the it first passes through the model base topology (*model_base*) to produce an intermediate representation *base_out* which is then forwarded to the various parallel output heads. In particular, both the *malware label prediction head* and the *vendor count prediction head*, if enabled, produce their predictions given the transient latent representation *base_out* as input. On the other hand, for the *tags prediction* topology, first the intermediate representation *base_out* is fed into the *pe_embedding* sub-net producing the latent representation of the input data samples in the joint embedding

Algorithm 29 Proposed Model Net class, Init

```

1: class NET
2:   function INIT(self, use_malware, use_counts, use_tags, n_tags, feature_dimension, embedding_dimension, embeddingmax_embedding_norm, layer_sizes, dropout_p, activation_function)
3:     self.use_malware  $\leftarrow$  use_malware
4:     self.use_counts  $\leftarrow$  use_counts
5:     self.n_tags  $\leftarrow$  n_tags
6:     self.embedding_dimension  $\leftarrow$  embedding_dimension
7:     layers  $\leftarrow$  []
8:
9:     if layer_sizes is None then
10:      layer_sizes  $\leftarrow$  [512, 512, 128]
11:    end if
12:
13:    for all i, ls in enumerate(layer_sizes) do
14:      if i = 0 then
15:        layers.append(nn.Linear(feature_dimension, ls))
16:      else
17:        layers.append(nn.Linear(layer_sizes[i - 1], ls))
18:      end if
19:
20:      layers.append(nn.LayerNorm(ls))
21:      layers.append(activation_function())
22:      layers.append(nn.Dropout(dropout_p))
23:    end for
24:
25:    self.model_base  $\leftarrow$  nn.Sequential(*tuple(layers))
26:    self.sigmoid  $\leftarrow$  nn.Sigmoid()
27:
28:    self.pe_embedding  $\leftarrow$  nn.Sequential(nn.Linear(layer_sizes[-1],
29:                                              self.embedding_dimension))
30:    self.malware_head  $\leftarrow$  nn.Sequential(nn.Linear(layer_sizes[-1], 1),
31:                                         nn.Sigmoid())
32:    self.count_head  $\leftarrow$  nn.Sequential(nn.Linear(layer_sizes[-1], 1), nn.ReLU())
33:    self.tags_embedding  $\leftarrow$  nn.Embedding(self.n_tags, self.embedding_dimension,
34:                                         max_norm=max_embedding_norm)
35:  end function

```

space (PE embedding). Next, the similarity scores are computed by applying the dot product (via a simple matrix multiplication) between the PE embedding and the tags embedding matrix (*tags_embedding*). Finally, the tags probability predictions are the result of applying the *sigmoid* activation function on the previously computed similarity scores.

If, however, the user chooses to try a different similarity function other than the *dot product* two alternatives are available: the *cosine similarity* or the *inverted pairwise euclidean distance*. In particular both the *Proposed Model forward function (with cosine similarity)*, shown in alg. 31, and the *Proposed Model forward function (with inverted pairwise Euclidean distance)*, shown in alg. 32, are for the most part identical to the original forward function but for the computation of the similarity scores. These, in fact, correspond in one case to the cosine similarities between the PE files' and tags' embeddings, which are values in the range $[-1,1]$, normalized to the $[0,1]$ interval, and in the other to the euclidean distances between the two embeddings, inverted through the *distance_to_similarity* function already mentioned in sec. 5.1.2. Again, similarly to what was done in the *Joint Embedding* model implementation, in both cases the final tags probabilities coincide with the just computed similarity scores.

The *Proposed Model* implementation also uses exactly the same *get_embedding* and *get_similarity* functions as the ones used in the *Joint Embedding* model implementation, shown in alg. 21.

Algorithm 30 Proposed Model Net class, Forward (with *Dot Product*)

```

36:   function FORWARD(self, data)
37:     rv  $\leftarrow \{\}$ 
38:     base_out  $\leftarrow \text{self.model\_base}(\text{data})$ 
39:
40:     if self.use_malware then
41:       rv[‘malware’]  $\leftarrow \text{self.malware\_head}(\text{base\_out})$ 
42:     end if
43:
44:     if self.use_counts then
45:       rv[‘count’]  $\leftarrow \text{self.count\_head}(\text{base\_out})$ 
46:     end if
47:
48:     pe_embedding  $\leftarrow \text{self.pe\_embedding}(\text{base\_out})$ 
49:     tags_embedding  $\leftarrow \text{self.tags\_embedding}((\text{Dataset.encoded\_tags}))$ 
50:
51:     rv[‘similarity’]  $\leftarrow \text{torch.matmul}(\text{pe\_embedding}, \text{tags\_embedding.T})$ 
52:     rv[‘probability’]  $\leftarrow \text{self.sigmoid}(\text{similarity\_scores})$ 
53:
54:     return rv
55:   end function

```

Algorithm 31 Proposed Model Net class, Forward (with *Cosine similarity*)

```

56:   function FORWARD(self, data)
57:     ...
58:     similarity_scores  $\leftarrow \text{torch.div}(\text{torch.add}(\text{Cosine\_similarity}(\text{pe\_embedding},$ 
59:                               tags_embedding, dim=2), 1.0), 2.0)
60:     rv[‘similarity’]  $\leftarrow \text{similarity\_scores}$ 
61:     rv[‘probability’]  $\leftarrow \text{similarity\_scores}$ 
62:     return rv
63:   end function

```

Algorithm 32 Proposed Model Net class, Forward (with *Inverted Pairwise Euclidean Distance*)

```

64:   function FORWARD(self, data)
65:     ...
66:     distances  $\leftarrow \text{torch.cdist}(\text{pe\_embedding}, \text{tags\_embedding}, \text{p}=2.0)$ 
67:     similarity_scores  $\leftarrow \text{distance\_to\_similarity}(\text{distances}, \text{a}=\text{sim\_function\_a},$ 
68:                               function=sim_function)
69:     rv[‘similarity’]  $\leftarrow \text{similarity\_scores}$ 
70:     rv[‘probability’]  $\leftarrow \text{similarity\_scores}$ 
71:     return rv
72:   end function

```

Again, these functions are particularly useful when performing evaluations on the learnt embedding.

Finally, algorithm 33 (***Proposed Model* compute loss function**) describes how the ***Proposed Model*** loss computation is implemented in *Python/Pytorch* code. In particular the final total loss corresponds to the weighted sum of the losses from each enabled output head. The weights used in this computation can be defined by the user at network definition. Both the *Malicious/Benign* label prediction loss and the *Vendor Count* prediction loss computations are taken from the ***ALOHA*** model implementation and are therefore implemented as the **binary cross entropy loss** between the predicted malware label and the respective ground truth label and as the **Negative Log Likelihood with Poisson distribution** between the predicted count and the ground truth value, respectively. On the other hand, the *Tags* prediction loss computation is taken from the ***Joint Embedding*** model implementation and is therefore the per-sample

Algorithm 33 Proposed Model Net class, Compute Loss

```

73:   function COMPUTE LOSS(predictions, labels, loss_wts)
74:     if loss_wts is None then
75:       loss_wts  $\leftarrow \{\text{'malware': } 1.0, \text{'count': } 0.1, \text{'tags': } 1.0\}$ 
76:     end if
77:     loss_dict  $\leftarrow \{\text{'total': } 0.\}$ 
78:
79:     if 'malware'  $\in$  labels then
80:       malware_loss  $\leftarrow$  Binary_cross_entropy(predictions[‘malware’], labels[‘malware’])
81:       if 'malware'  $\in$  loss_wts then
82:         weight  $\leftarrow$  loss_wts[‘malware’]
83:       else
84:         weight  $\leftarrow 1.0$ 
85:       end if
86:
87:       loss_dict[‘malware’]  $\leftarrow$  malware_loss
88:       loss_dict[‘total’]  $\leftarrow$  loss_dict[‘total’] + malware_loss  $\times$  weight
89:     end if
90:
91:     if 'count'  $\in$  labels then
92:       count_loss  $\leftarrow$  PoissonNLLLoss(predictions[‘count’], labels[‘count’])
93:       if 'count'  $\in$  loss_wts then
94:         weight  $\leftarrow$  loss_wts[‘count’]
95:       else
96:         weight  $\leftarrow 1.0$ 
97:       end if
98:
99:       loss_dict[‘count’]  $\leftarrow$  count_loss
100:      loss_dict[‘total’]  $\leftarrow$  loss_dict[‘total’] + count_loss  $\times$  weight
101:    end if
102:
103:    if 'tags'  $\in$  labels then
104:      similarity_loss  $\leftarrow$  Binary_cross_entropy(predictions[‘probability’], labels[‘tags’],
105:                                              reduction=‘none’).sum(dim=1).mean(dim=0)
106:      if 'tags'  $\in$  loss_wts then
107:        weight  $\leftarrow$  loss_wts[‘tags’]
108:      else
109:        weight  $\leftarrow 1.0$ 
110:      end if
111:
112:      loss_dict[‘jointEmbedding’]  $\leftarrow$  similarity_loss
113:      loss_dict[‘total’]  $\leftarrow$  loss_dict[‘total’] + similarity_loss  $\times$  weight
114:    end if
115:
116:    return loss_dict
117:  end function
118: end class

```

mean of all tags losses, each of which is computed as the **binary cross entropy** loss between the predicted tag probability and the ground truth tag.

6.2 Experiments

The proposed model was trained multiple times for 10 *epochs* on the first 6M training and 1.153.846 validation samples of the pre-processed Sorel20M dataset (loaded with generator *alt3*)

with both the optional output heads (*Malicious/Benign* label prediction and *Vendor Count* prediction) on, changing the network hyper-parameters like the number of layers, layers dimensions (number of hidden units per layer), activation functions and dropout value for the base topology. Other hyper-parameters that were iteratively changed were the used optimizer (**Adam** or **SGD**), learning rate (LR), momentum, weight decay, loss weights and the similarity score function used (chosen between the *dot product*, the *cosine similarity* and one of the *inverted Euclidean Distances*: *exp*, *inv* or *inv-pow*).

Then, for the evaluation of the resulting model the first 1.846.154 test samples of the Sorel20M pre-processed dataset were used to asses the model *TPR*, *accuracy*, *recall*, *precision*, *f1 score*, *jaccard similarity*, and *mean-per-sample accuracy* at different *FPRs* of interest (particularly at $FPR = 1\%$) for the tags and malicious/benign label prediction task, as well as the corresponding ROC curves and AUC scores.

Moreover, this model learned samples' representation (embedding), together with the *Joint Embedding* model implementation one, was also tested on the fresh dataset, which contains 1.000 malicious samples belonging to 10 different malware families to asses its performance in the malware family prediction and ranking tasks with respect to the mentioned previous method. The malware families represented by the fresh dataset were chosen among the most wide-spread PE malware families present in Italy at the time of writing (as reported by [53]) for which *Malware Bazaar* [54] provided at least 100 samples.

In particular, for the malware family prediction task the learned representation was tested, as previously described in 4.2.1, by selecting a number n (10) of anchor samples per family (for a total of $n \times n_{families} = 100$ anchor samples) and assigning to the other $n_{samples} - 100 = 900$ non-anchor samples, for each family, a value $p \in [0,1]$ based on their Euclidean distance to the most similar anchor sample of that family. Finally, binary prediction statistics like *TPR*, *Accuracy*, *Recall*, *Precision* and *F1 Score* are computed for each family at different *FPRs* (most importantly at $FPR = 1\%$). Moreover, also the corresponding ROC curves and AUC-ROC scores are computed.

On the other hand, for the malware family ranking task the learned representation was tested (as described in 4.2.1) by first selecting q (100) query samples from the fresh dataset and, for each of them, computing the ranking of the other $n_{samples} - 1 = 999$ ordering them by similarity. Then, given the resulting rankings, the **MAP** and **MRR** scores were calculated.

The final **Proposed Model** architecture consisted of 3 layers of sizes 512, 512, 128 respectively (keeping them consistent to the *ALOHA* and *Joint Embedding* model implementation in order to better compare the final results), with dropout probability $p_d = 5\%$, **ELU** activation function and the *dot product* as similarity score between tags' and samples' embeddings. The training was carried out with a learning rate $LR = 10^{-3}$, **Adam** optimizer, no momentum nor weight decay, and loss weights set to 1.0 for the malicious/benign head, 0.1 for the count head and 1.0 for the tag head.

The **Proposed Model** results are reported in chapter 7.

6.3 Training and Evaluation algorithms

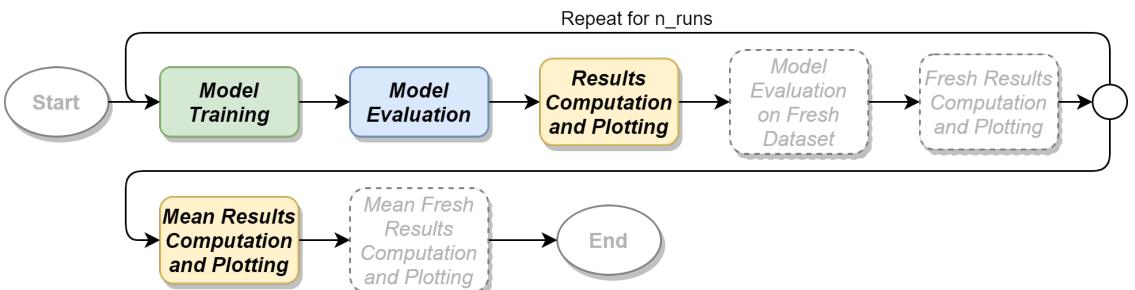


Figure 6.2: Workflow Model Training and Evaluation steps

As previously mentioned in section 5.3, the **Proposed Model** training and evaluation procedures are the same as the ones used for the **ALOHA** and **Joint Embedding** model implementations. In particular, the training procedure is presented in alg. 27, while the model evaluation is shown in alg. 28.

6.4 Evaluate Fresh algorithm

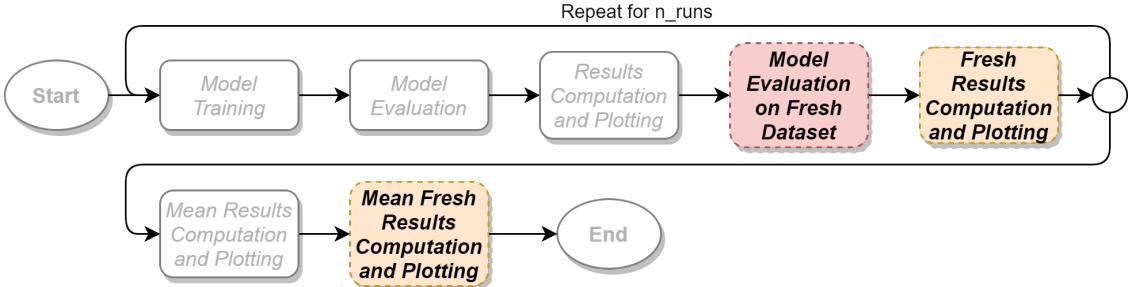


Figure 6.3: Workflow Model Evaluation on Fresh Dataset steps

Algorithm 34 Evaluate Fresh function

```

1: function EVALUATE_FRESH(ds_path, checkpoint_file, net_type, n_anchor_samples_per_family,
   hard, n_queries, batch_size)
2:   evaluate_fresh_predictions(ds_path, checkpoint_path, net_type,
   n_anchor_samples_per_family, hard, batch_size)
3:   evaluate_fresh_rankings(ds_path, checkpoint_path, net_type, n_queries, batch_size)
5: end function

```

As previously mentioned, both the **Proposed Model** and the **Joint Embedding** model implementations are also evaluated on the malware family prediction and ranking tasks. In particular the function presented in algorithm 34 (**Evaluate fresh function**) is used to produce both the evaluation results mentioned. More specifically, it calls both the following two functions in succession: *evaluate_fresh_predictions* and *evaluate_fresh_rankings*.

6.4.1 Model Evaluation on Malware Family Prediction/Ranking tasks

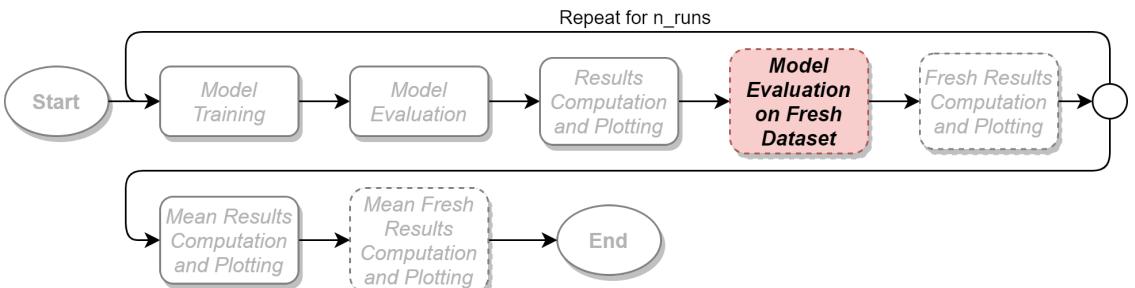


Figure 6.4: Workflow Model Evaluation on Fresh Dataset step

The *evaluate_fresh_predictions* function, shown in alg. 35, begins by importing the Net class corresponding to the chosen *net_type* and then instantiates it using the needed arguments (such as the *feature_dimension*, the *embedding_dimension*, the *layer_sizes*, etc.). Next, the function loads the model state checkpoint corresponding to a previous training run into the model current state, sets the model use modality to *evaluation* and instantiates the fresh dataset generator (dataloader). After these preliminary steps, in order to compute predict the family labels for

Algorithm 35 Evaluate Fresh Predictions function

```

1: function      EVALUATE_FRESH_PREDICTIONS(ds_path,           checkpoint_file,           net_type,
2:   n_anchor_samples_per_family, hard, batch_size)
3:   Net, similarity_type  $\leftarrow$  import_modules(net_type)
4:   model  $\leftarrow$  Net(use_malware=False, use_counts=False, n_tags=len(Dataset.tags),
5:                         feature_dimension=2381, embedding_dimension=32,
6:                         layer_sizes=run_params[‘layer_sizes’], dropout_p=run_params[‘dropout_p’],
7:                         activation_function=run_params[‘activation_function’])
8:   model.load_state_dict(checkpoint_path)
9:   model.eval()
10:  generator  $\leftarrow$  get_generator(ds_path, batch_size, return_shas=True, shuffle=True)
11:
12:  anchors  $\leftarrow$  {}
13:  for all shas, features, labels  $\in$  generator do
14:    embeddings  $\leftarrow$  model.get_embedding(features)[‘embedding’]
15:    for all n  $\in$  [0, n_families] do
16:      anchors.append(get_anchor_samples(shas, labels, embeddings, n,
17:                                         n_anchor_samples_per_family))
18:    end for
19:  end for
20:
21:  for all shas, features, labels  $\in$  generator do
22:    indices  $\leftarrow$  [i,  $\forall i, sha \in \text{enumerate}(shas)$  if sha not in anchors[‘shas’]]
23:    selected_shas  $\leftarrow$  [shas[i],  $\forall i \in indices$ ]
24:    selected_features  $\leftarrow$  features[indices]
25:    selected_labels  $\leftarrow$  labels[indices]
26:    embeddings  $\leftarrow$  model.get_embedding(selected_features)[‘embedding’]
27:    if similarity_type = ‘invertedEuclideanDistance’ then
28:      similarity_scores  $\leftarrow$  model.get_similarity(embeddings,
29:                                         anchors[‘embeddings’])[‘similarity’]
30:    else
31:      similarity_scores  $\leftarrow$  torch.div(torch.add(F.cosine_similarity(embeddings,
32:                                         anchors[‘embeddings’]), dim=2), 1.0), 2.0)
33:    end if
34:
35:    rv  $\leftarrow$  {}
36:    for all i  $\in$  [0, n_families] do
37:      rv[‘label_{i}’.format(label_to_sig(i))]  $\leftarrow$  [int(lab = i),  $\forall lab \in selected_labels$ ]
38:      if hard then
39:        rv[‘pred_{i}’.format(label_to_sig(i))]  $\leftarrow$  [int(torch.max(sims[[j,
40:                                          $\forall j \in [0, \text{len}(sims)]$  if anchors[‘labels’][j] == i]),  

41:                                         = torch.max(sims)),  $\forall sims \in similarity_scores$ ]
42:      else
43:        rv[‘pred_{i}’.format(label_to_sig(i))]  $\leftarrow$  [torch.max(sims[[j,
44:                                          $\forall j \in [0, \text{len}(sims)]$  if anchors[‘labels’][j] == i]),  

45:                                          $\forall sims \in similarity_scores$ ]
46:      end if
47:    end for
48:
49:    save_to_file(rv)
50:  end for
51: end function

```

the fresh dataset samples, a number of anchor samples is needed. Therefore, the function cycles through some of the mini-batches from the generator getting the *shas*, *labels* and *embeddings* for

Algorithm 36 Evaluate Fresh Rankings function

```

1: function EVALUATE_FRESH_RANKINGS(ds_path, checkpoint_file, net_type, n_queries,
2:   batch_size)
3:
4:   Net  $\leftarrow$  import_modules(net_type)
5:
6:
7:   model  $\leftarrow$  Net(use_malware=False, use_counts=False, n_tags=len(Dataset.tags),
8:     feature_dimension=2381, embedding_dimension=32,
9:     layer_sizes=run_params[‘layer_sizes’], dropout_p=run_params[‘dropout_p’],
10:    activation_function=run_params[‘activation_function’])
11:
12:   model.load_state_dict(checkpoint_path)
13:   model.eval()
14:   generator  $\leftarrow$  get_generator(ds_path, batch_size, return_shas=True, shuffle=True)
15:
16:
17:   queries  $\leftarrow$  None
18:   rank_per_query  $\leftarrow$  {}
19:   for all shas, features, labels  $\in$  generator do
20:     embeddings  $\leftarrow$  model.get_embedding(features)[‘embedding’]
21:
22:     if queries is None then
23:       queries  $\leftarrow$  {‘shas’: shas[n_queries], ‘embeddings’: embeddings[n_queries],
24:         ‘labels’: labels[n_queries]}
25:     end if
26:
27:     similarity_scores  $\leftarrow$  model.get_similarity(queries[‘embeddings’],
28:       embeddings)[‘similarity’]
29:
30:     for all i, s  $\in$  enumerate(queries[‘shas’]) do
31:       rank_per_query.update({
32:         i: {‘query_sha’: s,
33:           ‘ground_truth_label’: int(queries[‘labels’][i]),
34:           ‘ground_truth_family’: label_to_sig(int(queries[‘labels’][i])),
35:           ‘rank_shas’: shas[similarity_scores[i, j,  $\forall j \in (0, \text{len}(\text{similarity_scores}[i]))$ 
36:             if shas[j]  $\neq$  s].argsort()],
37:           ‘rank_labels’: [lab,  $\forall lab \in labels[similarity_scores[i, j,$ 
38:              $\forall j \in (0, \text{len}(\text{similarity_scores}[i])) \text{ if } shas[j] \neq s].argsort()]],
39:           ‘rank_families’: [label_to_sig(lab),  $\forall lab \in labels[similarity_scores[i, j,$ 
40:              $\forall j \in (0, \text{len}(\text{similarity_scores}[i])) \text{ if } shas[j] \neq s].argsort()]]}}}
41:       end for
42:     end for
43:
44:     rs  $\leftarrow$  binarize(rank_per_query)
45:     log(mean_reciprocal_rank(rs))
46:     log(mean_average_precision(rs))
47:
48:     queries_indexes  $\leftarrow$  {‘max_rr’: max_reciprocal_rank_index(rs),
49:       ‘min_rr’: min_reciprocal_rank_index(rs),
50:       ‘max_ap’: max_average_precision_index(rs),
51:       ‘min_ap’: min_average_precision_index(rs),
52:       ‘random’: random.choice([v  $\forall v \in (0, \text{len}(\text{rank_per_query.keys}()))$ 
53:         if v  $\notin$  [max_reciprocal_rank_index(rs), min_reciprocal_rank_index(rs),
54:           max_average_precision_index(rs), min_average_precision_index(rs)]]}}
55:
56:     ranks_to_save  $\leftarrow$  {key: rank_per_query[index]  $\forall key, index \in queries_indexes.items()$ }
57:     save_to_file(ranks_to_save)
58:   end function$$ 
```

`n_anchor_samples_per_family` samples for each of the `n_families` families. At this point the actual family predictions can be made for all the other **non-anchor** samples. The function, in fact, cycles through all the mini-batches from the fresh dataset generator (from the beginning) and filters out all the **anchor** samples in order to have the *shas*, *labels* and *features* of only **non-anchor** samples. Next, the **non-anchor** samples embeddings are obtained by passing the input features through the network. The similarity scores are computed by using the model's `get_similarity` function or by computing the normalized (to the [0,1] range) **Cosine Similarity** between the **non-anchor** samples' embeddings and the **anchor** samples' ones, depending on the similarity function used by the model. Finally, the similarities just computed are used to define the per-sample **hard** or **soft** family predictions. In particular, for each family, the resulting ground truth label vector (with one value per non-anchor sample) values are each set to 1.0 if the current **non-anchor** sample's label is equal to the current family label, 0.0 otherwise. Then, if the **hard** prediction mode is enabled, then, for each family, the resulting prediction vector (again with one value per sample) values are set to 1.0 if the maximum similarity between the current **non-anchor** sample and the **anchor** samples from the current family is equal to the overall maximum similarity between the current **non-anchor** sample and all **anchor** samples, to 0.0 otherwise. This is done in order to account for the case in which two malware families' anchor samples have the same minimum distance to the current **non-anchor** sample, in such instance this function assigns both families to the same **non-anchor** sample. On the other hand, if the **soft** prediction mode is enabled, then, for each family, the resulting prediction vector values are set to the maximum similarity between the current **non-anchor** sample and the **anchor** sample from the current family. Therefore, in this case, the prediction values can be values in the range [0,1] instead of only the values 0 and 1. In any case, the function ends by saving the malware family predictions to file.

The `evaluate_fresh_rankings` function, presented in algorithm 36, again begins by importing the selected *Net* class and instantiating the model with appropriate parameters; then it continues by loading the model state parameters from a checkpoint of a previous training run, setting the model use modality to *evaluation* and defining the fresh dataset generator (dataloader). Next, the function cycles through all the mini-batches of data from the fresh dataset generator and gets the samples embeddings by feeding the model with their numerical feature vectors. Then, the function selects the first `n_samples` samples (their *shas*, *labels* and *embeddings*) and uses them as **query** samples in order to compute family rankings. In particular, as first step for generating the family rankings, the function computes the similarity scores between the **query** samples' embeddings and all the mini-batch embeddings (which may contain also the **query** sample latent representation) by using the model's `get_similarity` function. Next, for each **query** sample the function generated one ranking by ordering the samples' *shas*, *family labels* and *family names* by their similarity with the current **query** sample (excluding the current **query** sample from the ranking). Finally, after having *binarized* the ranks labels (essentially by assigning a 1.0 to all samples in the ranking that are of the same family as the current **query** sample, 0.0 to the rest), the function computes the **MRR** and **MAP** scores and saves to file a bunch of interesting rankings, namely the ones which produces the maximum RR (reciprocal rank), the maximum AP (average precision), the minimum RR and minimum AP, plus a random ranking chosen between all the other ones.

6.4.2 Fresh Results Computation and Plotting

The actual binary cross entropy statistics for the malware family prediction task (such as *TPR*, *Accuracy*, *Recall*, etc.. at different *FPRs*) and the resulting *AUC-ROC* scores and *ROC* curves are computed by a distinct set of functions similar to the one mentioned in section 5.3 not presented in this document.

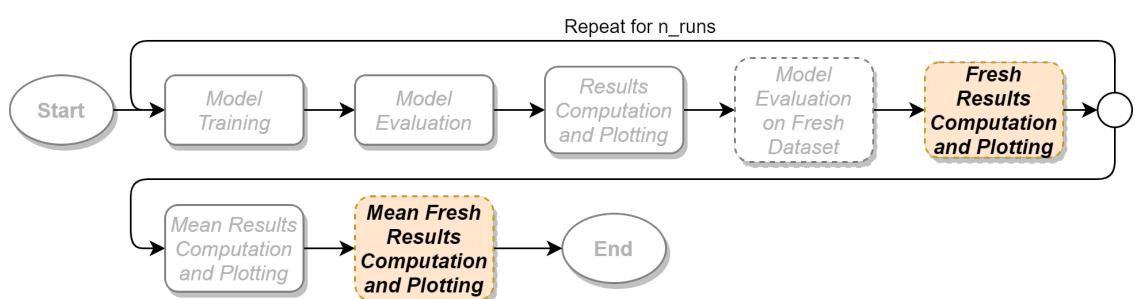


Figure 6.5: Workflow Fresh Results Computation and Plotting step

Chapter 7

Results

The following results are ‘fake’. They were taken from a model trained on just 8.192 training samples and are here presented only to show the tables structures.

In this chapter the evaluation results of the following model implementations will be presented and compared:

- **ALOHA (M/B only)**: *ALOHA* model implementation described in section 5.1.1, with only the *Malicious/Benign* label prediction head enabled. This model thus represents a simple single task FNN designed exclusively for the malicious label prediction task.
- **ALOHA**: final *ALOHA* model implementation described in section 5.1.1, with all the additional output heads enabled.
- **Joint Embedding**: final *Joint Embedding* model implementation described in section 5.1.2.
- **Proposed Model**: final *Proposed Model* implementation described in section 6.1.

In the first section (7.1) of this chapter, the models’ results concerning the **Malicious/Benign** label prediction task are presented. In particular for each model are shown the corresponding **AUC-ROC** (Area Under ROC Curve) score and the various binary cross entropy loss statistics (**TPR** (True Positive Rate), **Accuracy**, **Recall**, **Precision** and **F1 Score**) at different **FPRs** (False Positive Rates). Moreover, for the **TPR** statistic, also the **Proposed Model** Error and Standard Deviation Reductions with respect to the other (previous) models are presented, in percentage. Next, the **ROC** (Receiving Operating Characteristic) curves of the different models for the malicious label prediction task are presented. Finally, it is presented a brief summary of the different statistics for the different models at **FPR** = 1%.

In the next section (7.2), the models’ evaluation results for the SMART tagging prediction task are presented. More specifically, a table shows, for each predicted tag, the **AUC-ROC** scores relative to the different models. Another table presents the tag prediction per-tag binary cross entropy statistics at **FPR** = 1%. Finally, for each model it is presented a single graph showing the **ROC** curves of all predicted tags.

A further section (7.3) is used for showing the models’ *Mean per-sample scores* corresponding to the different tags for each model. In particular, here are presented the **Jaccard Similarity** and **Mean per-Sample Accuracy** scores at different **FPRs** achieved by the models.

A final section (7.4), instead, contains some tables showing the different models’ performance evaluations at the **Malware Family Prediction and Ranking** tasks. More specifically, a table shows the rankings’ **MRR** (Mean Reciprocal Rank) and **MAP** (Mean Average Precision) for the **Joint Embedding** and **Proposed Model** implementations. The **ALOHA** implementation, in fact, does not generate the joint embedding space needed for the ranking task. Next, two more tables show the models’ malware family prediction **AUC-ROC** scores and binary cross entropy statistics at **FPR** = 1%, similarly to what was done for the SMART tagging task. Finally,

the actual family prediction ***ROC*** curves for the 2 mentioned model implementations (**Joint Embedding** and **Proposed Model**) are presented, along with the 4 fresh dataset samples' rankings which produced the maximum and minimum AP (Average Precision) and RR (Reciprocal Rank) scores.

7.1 Malware Label results

Malware Label	ALOHA (M/B only)	ALOHA	Joint Embedding	Proposed Model
AUC-ROC	0.995±0.000	0.994±0.001	-	0.995±0.001

Table 7.1: AUC-ROC (Area Under Curve) of the different models for the **Malware Label** prediction task. Results were aggregated over **3** training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**.

Malware Label	FPR				
	10 ⁻⁵	10 ⁻⁴	10 ⁻³	10 ⁻²	10 ⁻¹
TPR					
ALOHA (M/B only)	0.281±0.093	0.671±0.020	0.864±0.014	0.950±0.003	0.989±0.000
ALOHA	0.346±0.014	0.672±0.023	0.855±0.013	0.952±0.003	0.987±0.003
Joint Embedding	-	-	-	-	-
Proposed Model	0.514±0.037	0.771±0.005	0.892±0.015	0.959±0.008	0.990±0.000
Error Reduction wrt ALOHA (M/B only)	32.4%	30.4%	20.6%	18.0%	9.1%
Error Reduction wrt ALOHA	25.7%	30.2%	25.5%	14.6%	23.1%
Error Reduction wrt Joint Embedding	-	-	-	-	-
Std Reduction wrt ALOHA (M/B only)	60.2%	75.0%	-7.1%	-166.7%	0.0%
Std Reduction wrt ALOHA	-164.3%	78.3%	-15.4%	-166.7%	100.0%
Std Reduction wrt Joint Embedding	-	-	-	-	-
Accuracy					
ALOHA (M/B only)	0.721±0.037	0.873±0.008	0.947±0.005	0.975±0.001	0.934±0.000
ALOHA	0.747±0.005	0.873±0.009	0.943±0.005	0.975±0.001	0.934±0.001
Joint Embedding	-	-	-	-	-
Proposed Model	0.812±0.014	0.911±0.002	0.957±0.006	0.978±0.003	0.935±0.000
Recall					
ALOHA (M/B only)	0.279±0.095	0.671±0.020	0.864±0.014	0.950±0.003	0.989±0.000
ALOHA	0.346±0.014	0.672±0.023	0.855±0.013	0.952±0.003	0.987±0.003

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Malware Label	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
Joint Embedding	-	-	-	-	-
Proposed Model	0.514 ± 0.037	0.771 ± 0.005	0.892 ± 0.015	0.959 ± 0.008	0.990 ± 0.000
Precision					
ALOHA (M/B only)	1.000 ± 0.000	1.000 ± 0.000	0.998 ± 0.000	0.984 ± 0.000	0.862 ± 0.000
ALOHA	1.000 ± 0.000	1.000 ± 0.000	0.998 ± 0.000	0.984 ± 0.000	0.862 ± 0.000
Joint Embedding	-	-	-	-	-
Proposed Model	1.000 ± 0.000	1.000 ± 0.000	0.998 ± 0.000	0.984 ± 0.000	0.862 ± 0.000
F1 Score					
ALOHA (M/B only)	0.427 ± 0.122	0.803 ± 0.014	0.926 ± 0.008	0.967 ± 0.002	0.921 ± 0.000
ALOHA	0.514 ± 0.016	0.804 ± 0.016	0.921 ± 0.007	0.967 ± 0.001	0.920 ± 0.002
Joint Embedding	-	-	-	-	-
Proposed Model	0.678 ± 0.032	0.870 ± 0.003	0.942 ± 0.008	0.971 ± 0.004	0.922 ± 0.000

Table 7.2: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Malware Label** prediction task at different **FPRs (False Positive Rates)**. Results were aggregated over **3** training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

Results

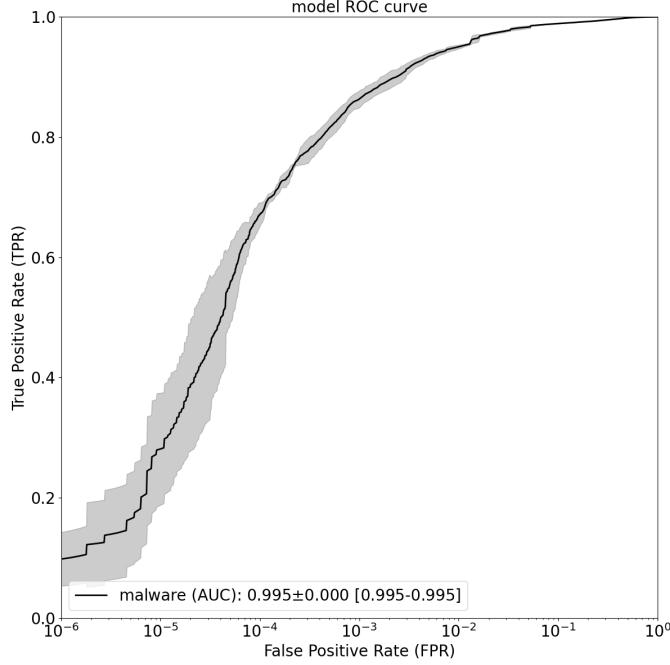


Figure 7.1: ROC curve and AUC statistics of **ALOHA (M/B only)** model for the **Malware Label**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

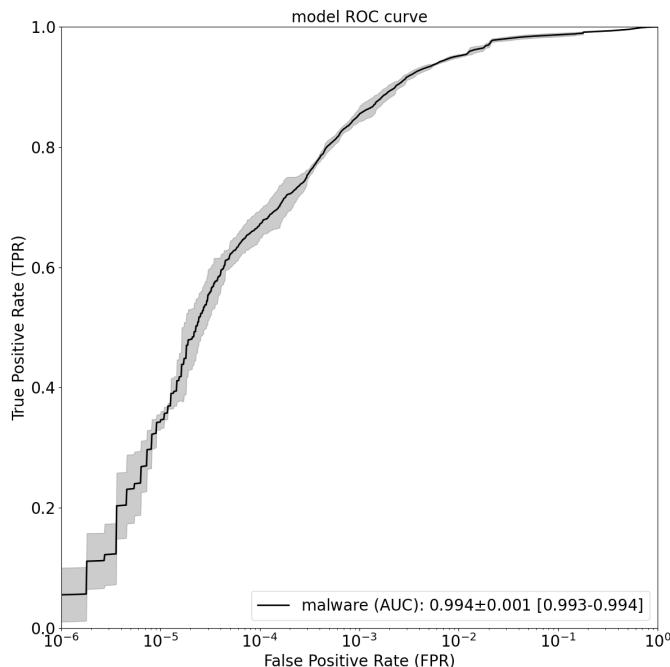


Figure 7.2: ROC curve and AUC statistics of **ALOHA** model for the **Malware Label**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

Results

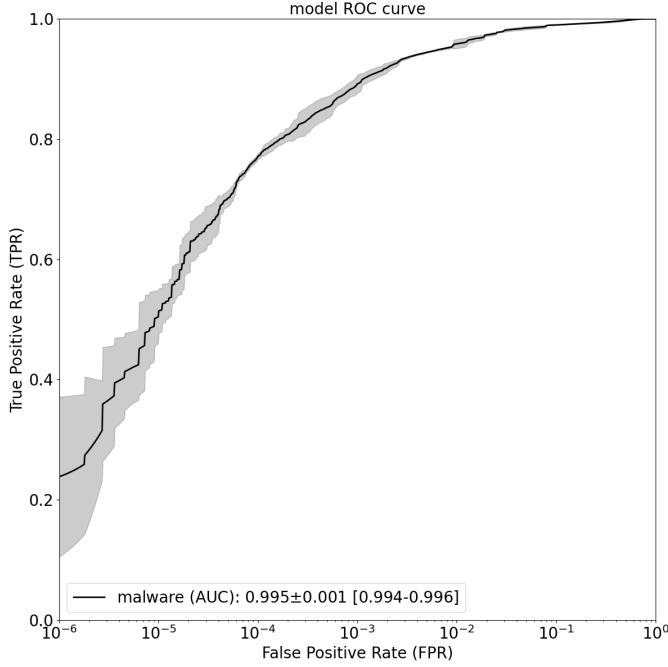


Figure 7.3: ROC curve and AUC statistics of **Proposed Model** for the **Malware Label**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

7.1.1 Summary

Malware Label (at FPR = 1%)					
Model	TPR	Accuracy	Precision	Recall	F1 score
ALOHA (M/B only)	0.950±0.003	0.975±0.001	0.984±0.000	0.950±0.003	0.967±0.002
ALOHA	0.952±0.003	0.975±0.001	0.984±0.000	0.952±0.003	0.967±0.001
Joint Embedding	-	-	-	-	-
Proposed Model	0.959±0.008	0.978±0.003	0.984±0.000	0.959±0.008	0.971±0.004

Table 7.3: Summary of the mean and standard deviation results of the different models for the **Malware Label** prediction task at **FPR** = 1%. Results were aggregated over **3** training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**.

7.2 Per-tag results

AUC-ROC	ALOHA (M/B only)	ALOHA	Joint Embedding	Proposed Model
Adware Tag	-	0.964±0.001	0.972±0.002	0.973±0.001
Crypto-miner Tag	-	0.980±0.003	0.986±0.003	0.990±0.003
Downloader Tag	-	0.968±0.007	0.974±0.002	0.980±0.002
Dropper Tag	-	0.968±0.005	0.972±0.001	0.977±0.001
File-infecter Tag	-	0.985±0.000	0.986±0.000	0.987±0.002
Flooder Tag	-	0.983±0.002	0.988±0.000	0.988±0.001

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AUC-ROC	ALOHA (M/B only)	ALOHA	Joint Embedding	Proposed Model
Installer Tag	-	0.967±0.007	0.975±0.001	0.975±0.001
Packed Tag	-	0.978±0.003	0.980±0.001	0.982±0.001
Ransomware Tag	-	0.977±0.003	0.983±0.003	0.987±0.001
Spyware Tag	-	0.959±0.005	0.971±0.001	0.968±0.001
Worm Tag	-	0.966±0.010	0.928±0.014	0.940±0.014

Table 7.4: Mean and standard deviation AUC-ROC (Area Under Curve) of the different models for the prediction of the different Tags. Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**.

	TPR	Accuracy	Precision	Recall	F1 score
Adware Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.682±0.010	0.972±0.001	0.806±0.002	0.682±0.010	0.739±0.007
Joint Embedding	0.698±0.008	0.973±0.000	0.810±0.002	0.698±0.008	0.750±0.005
Proposed Model	0.721±0.014	0.975±0.001	0.815±0.003	0.721±0.014	0.765±0.009
Crypto-miner Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.773±0.148	0.987±0.002	0.515±0.052	0.773±0.148	0.616±0.086
Joint Embedding	0.896±0.008	0.989±0.000	0.556±0.002	0.896±0.008	0.687±0.004
Proposed Model	0.905±0.006	0.989±0.000	0.559±0.002	0.905±0.006	0.691±0.003
Downloader Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.632±0.034	0.960±0.003	0.852±0.007	0.632±0.034	0.725±0.025
Joint Embedding	0.635±0.031	0.960±0.003	0.853±0.006	0.635±0.031	0.728±0.023
Proposed Model	0.667±0.012	0.963±0.001	0.859±0.002	0.667±0.012	0.751±0.009
Dropper Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.673±0.037	0.950±0.005	0.908±0.005	0.673±0.037	0.772±0.026
Joint Embedding	0.702±0.031	0.953±0.004	0.911±0.004	0.702±0.031	0.793±0.021
Proposed Model	0.690±0.016	0.952±0.002	0.910±0.002	0.690±0.016	0.785±0.011
File-infector Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.875±0.003	0.971±0.001	0.944±0.000	0.875±0.003	0.908±0.002

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	TPR	Accuracy	Precision	Recall	F1 score
Joint Embedding	0.867±0.015	0.970±0.002	0.943±0.001	0.867±0.015	0.903±0.009
Proposed Model	0.883±0.014	0.973±0.002	0.945±0.001	0.883±0.014	0.913±0.008
Flooder Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.876±0.016	0.990±0.000	0.137±0.002	0.876±0.016	0.237±0.004
Joint Embedding	0.916±0.003	0.990±0.000	0.142±0.000	0.916±0.003	0.246±0.001
Proposed Model	0.906±0.007	0.990±0.000	0.141±0.001	0.906±0.007	0.244±0.002
Installer Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.717±0.054	0.985±0.001	0.564±0.019	0.717±0.054	0.631±0.033
Joint Embedding	0.777±0.010	0.986±0.000	0.584±0.003	0.777±0.010	0.667±0.005
Proposed Model	0.759±0.042	0.986±0.001	0.578±0.014	0.759±0.042	0.656±0.024
Packed Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.715±0.003	0.952±0.000	0.919±0.000	0.715±0.003	0.804±0.002
Joint Embedding	0.764±0.014	0.959±0.002	0.924±0.001	0.764±0.014	0.836±0.009
Proposed Model	0.735±0.015	0.955±0.002	0.921±0.002	0.735±0.015	0.817±0.010
Ransomware Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.850±0.002	0.983±0.000	0.824±0.000	0.850±0.002	0.837±0.001
Joint Embedding	0.856±0.003	0.983±0.000	0.825±0.000	0.856±0.003	0.840±0.002
Proposed Model	0.856±0.001	0.983±0.000	0.825±0.000	0.856±0.001	0.841±0.001
Spyware Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.582±0.031	0.946±0.003	0.874±0.006	0.582±0.031	0.698±0.025
Joint Embedding	0.688±0.033	0.958±0.004	0.892±0.004	0.688±0.033	0.777±0.022
Proposed Model	0.660±0.003	0.955±0.000	0.888±0.000	0.660±0.003	0.757±0.002
Worm Tag (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.605±0.012	0.929±0.002	0.919±0.002	0.605±0.012	0.730±0.009
Joint Embedding	0.656±0.004	0.937±0.001	0.925±0.000	0.656±0.004	0.767±0.003
Proposed Model	0.657±0.007	0.937±0.001	0.925±0.001	0.657±0.007	0.768±0.005

Table 7.5: Summary of the mean and standard deviation results of the different models for the prediction of the different tags at **FPR** = 1%. Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**.

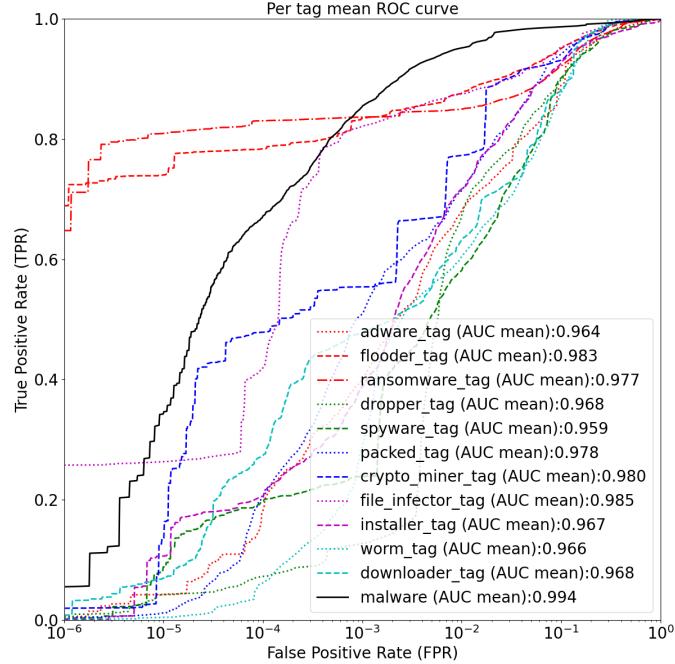


Figure 7.4: Mean ROC curve and AUC statistics of **ALOHA** model for the tags/labels. The line represents the *mean* TPR at a given FPR. Statistics were computed over 3 training runs, each with random parameter initialization.

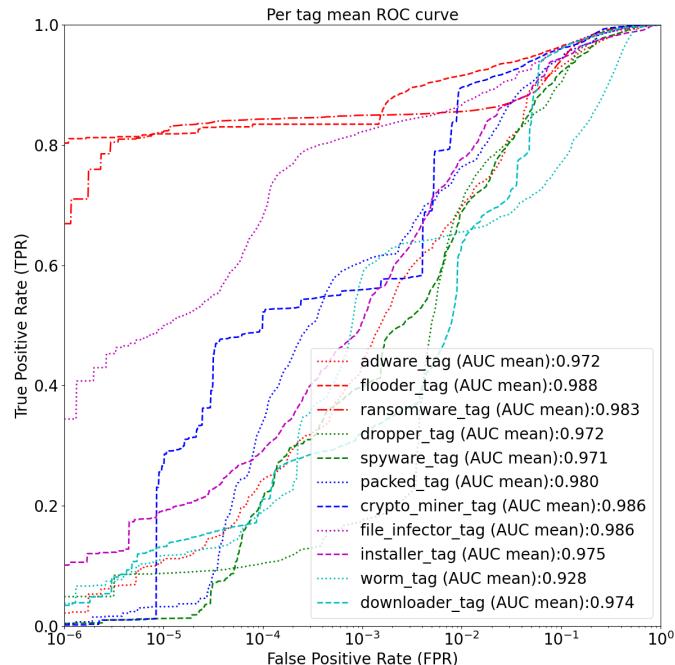


Figure 7.5: Mean ROC curve and AUC statistics of **Joint Embedding** model for all tags/labels. The line represents the *mean* TPR at a given FPR. Statistics were computed over 3 training runs, each with random parameter initialization.

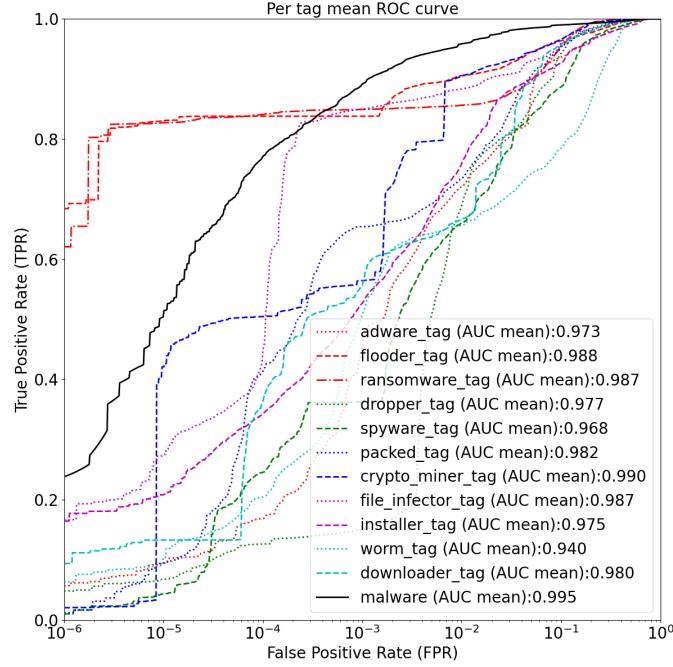


Figure 7.6: Mean ROC curve and AUC statistics of Proposed Model for the tags/labels. The line represents the *mean* TPR at a given FPR. Statistics were computed over **3** training runs, each with random parameter initialization.

7.3 Mean per-sample tagging scores

Mean per-sample tagging scores	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
Jaccard Similarity					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.674 ± 0.004	0.716 ± 0.013	0.787 ± 0.014	0.864 ± 0.003	0.793 ± 0.002
Joint Embedding	0.703 ± 0.016	0.751 ± 0.016	0.826 ± 0.020	0.880 ± 0.005	0.784 ± 0.003
Proposed Model	0.693 ± 0.017	0.748 ± 0.017	0.830 ± 0.007	0.877 ± 0.002	0.797 ± 0.005
Mean per-Sample Accuracy					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.633 ± 0.002	0.658 ± 0.014	0.716 ± 0.031	0.791 ± 0.002	0.699 ± 0.003
Joint Embedding	0.643 ± 0.005	0.680 ± 0.012	0.763 ± 0.021	0.806 ± 0.004	0.684 ± 0.006
Proposed Model	0.645 ± 0.005	0.682 ± 0.006	0.764 ± 0.013	0.806 ± 0.002	0.699 ± 0.004

Table 7.6: Mean and standard deviation of mean per-sample tagging results (*Jaccard simialrity* and *mean per-sample accuracy*) for the different models. Results were aggregated over **3** training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**.

7.4 Family Ranking Evaluation

	ALOHA (M/B only)	ALOHA	Joint Embedding	Proposed Model
MRR	-	-	0.252±0.029	0.255±0.024
MAP	-	-	0.106±0.001	0.106±0.001

Table 7.7: Mean and standard deviation MRR (Mean Reciprocal Rank) and MAP (Mean Average Precision) results for the different models on the Family Ranking task. Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**.

AUC-ROC	ALOHA (M/B only)	ALOHA	Joint Embedding	Proposed Model
Agenttesla	-	-	0.616±0.091	0.629±0.052
Avemariarat	-	-	0.495±0.046	0.510±0.029
Azorult	-	-	0.588±0.016	0.588±0.009
Formbook	-	-	0.533±0.026	0.513±0.035
Heodo	-	-	0.798±0.016	0.778±0.025
Loki	-	-	0.577±0.016	0.568±0.015
Masslogger	-	-	0.513±0.040	0.550±0.029
Netwire	-	-	0.503±0.008	0.486±0.024
Njrat	-	-	0.722±0.030	0.651±0.023
Remcorsat	-	-	0.545±0.077	0.531±0.055

Table 7.8: Mean and standard deviation AUC-ROC (Area Under Curve) of the different models for the prediction of the different families on fresh dataset samples. Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**.

	TPR	Accuracy	Precision	Recall	F1 score
Agenttesla (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-
Joint Embedding	0.014±0.003	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.015±0.003	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Avemariarat (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-

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	TPR	Accuracy	Precision	Recall	F1 score
Joint Embedding	0.010±0.002	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.010±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Azorult (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-
Joint Embedding	0.013±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.013±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Formbook (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-
Joint Embedding	0.011±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.010±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Heodo (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-
Joint Embedding	0.033±0.009	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.026±0.005	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Loki (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-
Joint Embedding	0.012±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.012±0.000	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Masslogger (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-
Joint Embedding	0.011±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.012±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Netwire (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-
Joint Embedding	0.010±0.000	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.009±0.001	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Njrat (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-

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	TPR	Accuracy	Precision	Recall	F1 score
Joint Embedding	0.022±0.004	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.017±0.002	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Remcorstat (at FPR = 1%)					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	-	-	-	-	-
Joint Embedding	0.013±0.005	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000
Proposed Model	0.011±0.002	0.900±0.000	1.000±0.000	0.000±0.000	0.000±0.000

Table 7.9: Summary of the mean and standard deviation results of the different models for the prediction of the different families on fresh dataset samples at **FPR = 1%**. Results were aggregated over **3** training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**.

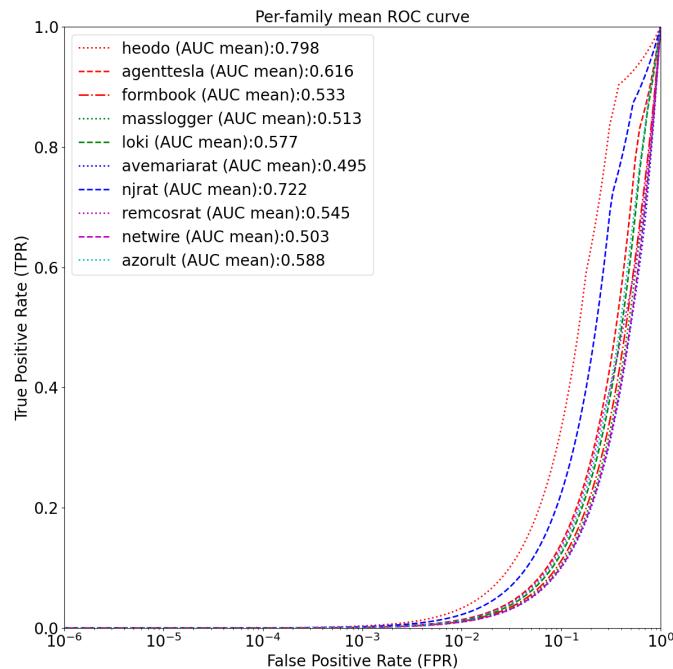


Figure 7.7: Mean ROC curve and AUC statistics of **Joint Embedding** model for the prediction of all families on fresh dataset samples. The line represents the *mean* TPR at a given FPR. Statistics were computed over **3** training runs, each with random parameter initialization.

Results

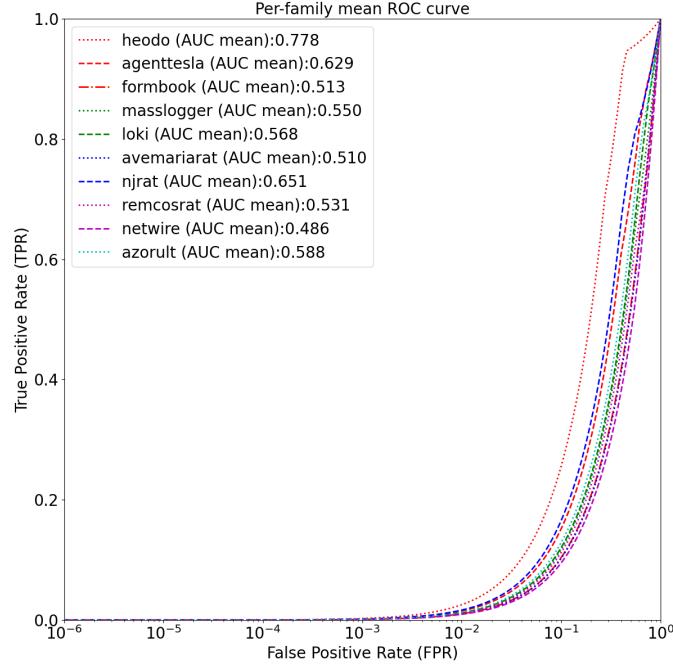


Figure 7.8: Mean ROC curve and AUC statistics of **Proposed Model** for the prediction of all families on fresh dataset samples. The line represents the *mean* TPR at a given FPR. Statistics were computed over **3** training runs, each with random parameter initialization.

7.4.1 Example rankings

Max AP	Joint Embedding			Proposed Model		
	Sha256	Label	Family	Sha256	Label	Family
Query	f8be22..	2	formbook	e8cac4..	7	remcosrat
0	3418e1..	7	remcosrat	351afe..	0	heodo
1	63a870..	3	masslogger	7bcd06..	1	agenttesla
2	271006..	9	azorult	e8cac4..	7	remcosrat
3	91069b..	1	agenttesla	72c2e8..	3	masslogger
4	0dcebc..	4	loki	057883..	3	masslogger
5	0ad577..	6	njrat	e1108e..	9	azorult
6	13b82d..	0	heodo	e482d2..	8	netwire
7	cb79b3..	0	heodo	46d544..	7	remcosrat
8	3b8dd3..	8	netwire	893e86..	5	avemariarat
9	e124d3..	0	heodo	8d55fa..	9	azorult

Table 7.10: **Joint Embedding** and **Proposed Model** example rankings (limited to the first 10 samples) having the maximum Average Precision (max AP). The elements matching the query sample are shown in **bold**. The next occurrence of family ‘formbook’ in the joint Embedding ranking is at positions **20**. On the other hand, the next occurrence of family ‘remcosrat’ in the Proposed Model ranking is at position **12**.

Max RR	Joint Embedding			Proposed Model		
	Sha256	Label	Family	Sha256	Label	Family
Query	82841c..	7	remcosrat	963553..	0	heodo
0	3418e1..	7	remcosrat	351afe..	0	heodo
1	63a870..	3	masslogger	7bcd06..	1	agenttesla
2	271006..	9	azorult	e8cac4..	7	remcosrat
3	91069b..	1	agenttesla	72c2e8..	3	masslogger
4	0dcebc..	4	loki	057883..	3	masslogger
5	33be12..	5	avemariarat	e1108e..	9	azorult
6	13b82d..	0	heodo	e482d2..	8	netwire
7	cb79b3..	0	heodo	46d544..	7	remcosrat
8	3b8dd3..	8	netwire	893e86..	5	avemariarat
9	e124d3..	0	heodo	8d55fa..	9	azorult

Table 7.11: **Joint Embedding** and **Proposed Model** example rankings (limited to the first 10 samples) having the maximum Reciprocal Rank (max RR). The elements matching the query sample are shown in **bold**. The next occurrence of family 'remcosrat' in the joint Embedding ranking is at positions **11**. On the other hand, the next occurrence of family 'heodo' in the Proposed Model ranking is at position **19**.

Min AP	Joint Embedding			Proposed Model		
	Sha256	Label	Family	Sha256	Label	Family
Query	9a825e..	5	avemariarat	a4a8d0..	4	loki
0	3418e1..	7	remcosrat	351afe..	0	heodo
1	63a870..	3	masslogger	7bcd06..	1	agenttesla
2	271006..	9	azorult	e8cac4..	7	remcosrat
3	91069b..	1	agenttesla	72c2e8..	3	masslogger
4	0dcebc..	4	loki	057883..	3	masslogger
5	0ad577..	6	njrat	e1108e..	9	azorult
6	13b82d..	0	heodo	e482d2..	8	netwire
7	cb79b3..	0	heodo	46d544..	7	remcosrat
8	3b8dd3..	8	netwire	893e86..	5	avemariarat
9	e124d3..	0	heodo	8d55fa..	9	azorult

Table 7.12: **Joint Embedding** and **Proposed Model** example rankings (limited to the first 10 samples) having the minimum Average Precision (min AP). The elements matching the query sample are shown in **bold**. The next occurrence of family 'avemariarat' in the joint Embedding ranking is at positions **28**. On the other hand, the next occurrence of family 'loki' in the Proposed Model ranking is at position **70**.

Min RR	Joint Embedding			Proposed Model		
	Sha256	Label	Family	Sha256	Label	Family
Query	b170f2..	5	avemariarat	a63d9a..	4	loki
0	3418e1..	7	remcosrat	351afe..	0	heodo
1	63a870..	3	masslogger	7bcd06..	1	agenttesla
2	271006..	9	azorult	e8cac4..	7	remcosrat
3	91069b..	1	agenttesla	72c2e8..	3	masslogger
4	0dcebc..	4	loki	057883..	3	masslogger
5	0ad577..	6	njrat	e1108e..	9	azorult
6	13b82d..	0	heodo	e482d2..	8	netwire
7	cb79b3..	0	heodo	46d544..	7	remcosrat
8	3b8dd3..	8	netwire	893e86..	5	avemariarat
9	e124d3..	0	heodo	8d55fa..	9	azorult

Table 7.13: **Joint Embedding** and **Proposed Model** example rankings (limited to the first 10 samples) having the minimum Reciprocal Rank (min RR). The elements matching the query sample are shown in **bold**. The next occurrence of family '**avemariarat**' in the joint Embedding ranking is at positions **28**. On the other hand, the next occurrence of family '**loki**' in the Proposed Model ranking is at position **70**.

Chapter 8

Conclusions

Qui si inseriscono brevi conclusioni sul lavoro svolto, senza ripetere inutilmente il sommario.

Si possono evidenziare i punti di forza e quelli di debolezza, nonché i possibili sviluppi futuri o attività da svolgere per migliorare i risultati.

Chapter 9

Appendix

9.1 More Detailed Per-Tag Results

The following results are 'fake'. They were taken from a model trained on just 8.192 training samples and are here presented only to show the tables structures.

Here are presented more detailed results regarding the different models' predictions in the SMART tag prediction task. More specifically, for each task are reported the *Binary cross entropy* statistics (**TPR**, **Accuracy**, **Recall**, etc.) at different **FPRs** for the models and the corresponding **ROC** curve plots (with mean and standard deviation).

9.1.1 Adware tag results

Adware Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.042 ± 0.055	0.181 ± 0.037	0.402 ± 0.040	0.682 ± 0.010	0.890 ± 0.027
Joint Embedding	0.110 ± 0.050	0.244 ± 0.015	0.433 ± 0.037	0.698 ± 0.008	0.931 ± 0.002
Proposed Model	0.095 ± 0.030	0.169 ± 0.021	0.456 ± 0.023	0.721 ± 0.014	0.937 ± 0.003
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	5.5%	-1.5%	9.0%	12.3%	42.7%
Error Reduction wrt Joint Embedding	-1.7%	-9.9%	4.1%	7.6%	8.7%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	45.5%	43.2%	42.5%	-40.0%	88.9%
Std Reduction wrt Joint Embedding	40.0%	-40.0%	37.8%	-75.0%	-50.0%
Accuracy					

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		FPR				
		10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
Adware Tag		-	-	-	-	-
ALOHA (M/B only)		-	-	-	-	-
ALOHA		0.945 ± 0.003	0.953 ± 0.002	0.965 ± 0.002	0.972 ± 0.001	0.899 ± 0.002
Joint Embedding		0.949 ± 0.003	0.956 ± 0.001	0.966 ± 0.002	0.973 ± 0.000	0.902 ± 0.000
Proposed Model		0.948 ± 0.002	0.952 ± 0.001	0.968 ± 0.001	0.975 ± 0.001	0.902 ± 0.000
Recall						
ALOHA (M/B only)		-	-	-	-	-
ALOHA		0.042 ± 0.055	0.181 ± 0.037	0.402 ± 0.040	0.682 ± 0.010	0.890 ± 0.027
Joint Embedding		0.110 ± 0.050	0.244 ± 0.015	0.433 ± 0.037	0.698 ± 0.008	0.931 ± 0.002
Proposed Model		0.095 ± 0.030	0.169 ± 0.021	0.456 ± 0.023	0.721 ± 0.014	0.937 ± 0.003
Precision						
ALOHA (M/B only)		-	-	-	-	-
ALOHA		0.964 ± 0.033	0.991 ± 0.002	0.961 ± 0.004	0.806 ± 0.002	0.352 ± 0.007
Joint Embedding		0.998 ± 0.001	0.993 ± 0.000	0.963 ± 0.003	0.810 ± 0.002	0.362 ± 0.000
Proposed Model		0.998 ± 0.001	0.990 ± 0.001	0.965 ± 0.002	0.815 ± 0.003	0.364 ± 0.001
F1 Score						
ALOHA (M/B only)		-	-	-	-	-
ALOHA		0.076 ± 0.097	0.305 ± 0.053	0.565 ± 0.039	0.739 ± 0.007	0.504 ± 0.011
Joint Embedding		0.195 ± 0.081	0.391 ± 0.019	0.597 ± 0.036	0.750 ± 0.005	0.522 ± 0.001
Proposed Model		0.172 ± 0.050	0.288 ± 0.030	0.619 ± 0.021	0.765 ± 0.009	0.524 ± 0.001

Table 9.1: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Adware Tag** prediction task at different **FPRs (False Positive Rates)**. Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

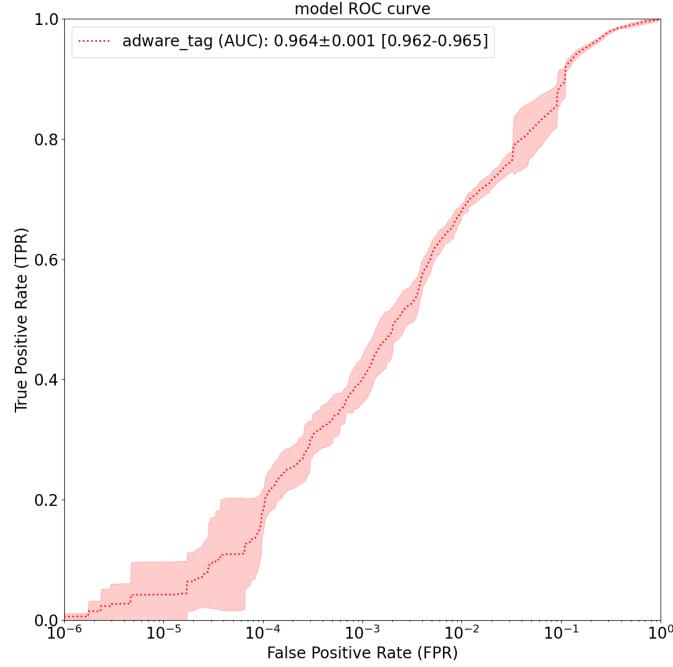


Figure 9.1: ROC curve and AUC statistics of **ALOHA** model for the **Adware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

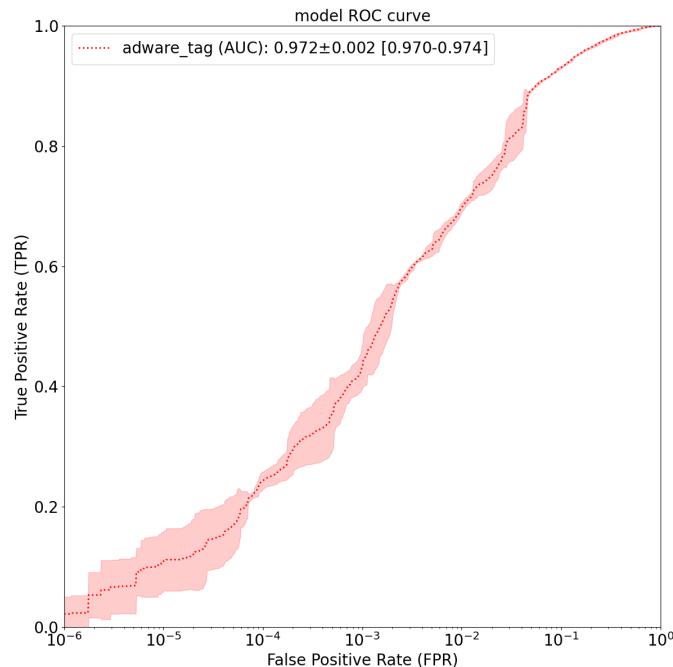


Figure 9.2: ROC curve and AUC statistics of **Joint Embedding** model for the **Adware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

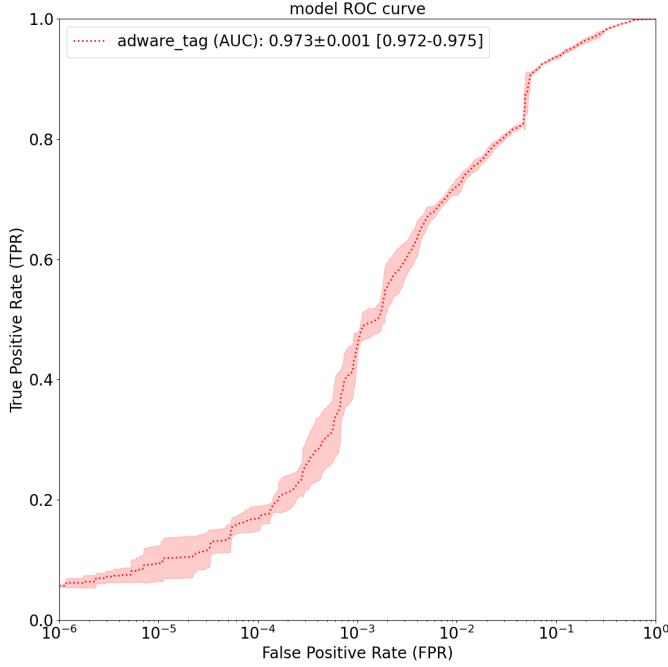


Figure 9.3: ROC curve and AUC statistics of **Proposed Model** for the **Adware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.2 Crypto-miner tag results

Crypto-miner Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.136 ± 0.048	0.479 ± 0.012	0.554 ± 0.014	0.773 ± 0.148	0.931 ± 0.016
Joint Embedding	0.266 ± 0.113	0.522 ± 0.015	0.560 ± 0.029	0.896 ± 0.008	0.961 ± 0.015
Proposed Model	0.421 ± 0.045	0.505 ± 0.005	0.565 ± 0.019	0.905 ± 0.006	0.966 ± 0.015
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	33.0%	5.0%	2.5%	58.1%	50.7%
Error Reduction wrt Joint Embedding	21.1%	-3.6%	1.1%	8.7%	12.8%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	6.3%	58.3%	-35.7%	95.9%	6.3%
Std Reduction wrt Joint Embedding	60.2%	66.7%	34.5%	25.0%	0.0%
Accuracy					
ALOHA (M/B only)	-	-	-	-	-

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Crypto-miner Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA	0.988±0.001	0.993±0.000	0.993±0.000	0.987±0.002	0.900±0.000
Joint Embedding	0.990±0.002	0.993±0.000	0.993±0.000	0.989±0.000	0.901±0.000
Proposed Model	0.992±0.001	0.993±0.000	0.993±0.000	0.989±0.000	0.901±0.000
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.136±0.048	0.479±0.012	0.554±0.014	0.773±0.148	0.931±0.016
Joint Embedding	0.266±0.113	0.522±0.015	0.560±0.029	0.896±0.008	0.961±0.015
Proposed Model	0.421±0.045	0.505±0.005	0.565±0.019	0.905±0.006	0.966±0.015
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.994±0.002	0.985±0.000	0.886±0.003	0.515±0.052	0.115±0.002
Joint Embedding	0.997±0.002	0.987±0.000	0.887±0.006	0.556±0.002	0.119±0.002
Proposed Model	0.998±0.000	0.986±0.000	0.888±0.003	0.559±0.002	0.119±0.002
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.237±0.075	0.644±0.011	0.682±0.011	0.616±0.086	0.205±0.003
Joint Embedding	0.407±0.152	0.683±0.013	0.686±0.024	0.687±0.004	0.211±0.003
Proposed Model	0.591±0.045	0.668±0.005	0.690±0.015	0.691±0.003	0.212±0.003

Table 9.2: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Crypto-miner Tag** prediction task at different **FPRs** (*False Positive Rates*). Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

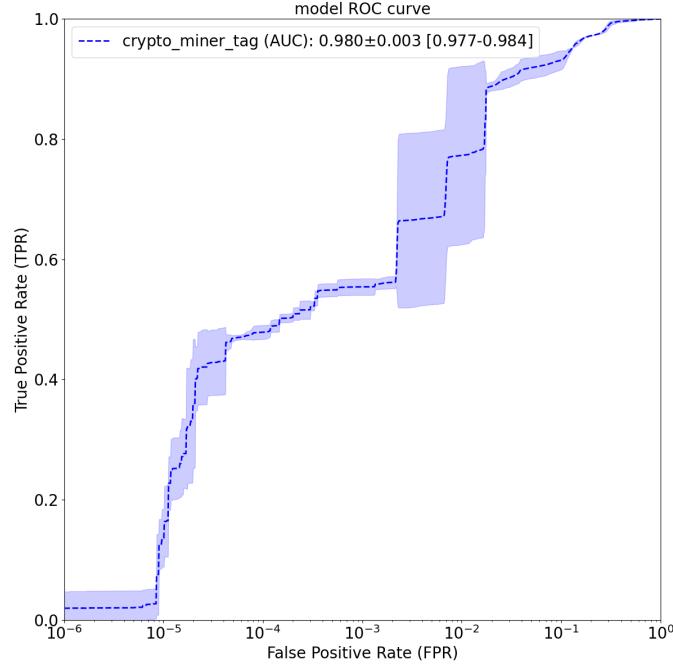


Figure 9.4: ROC curve and AUC statistics of **ALOHA** model for the **Crypto-miner Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

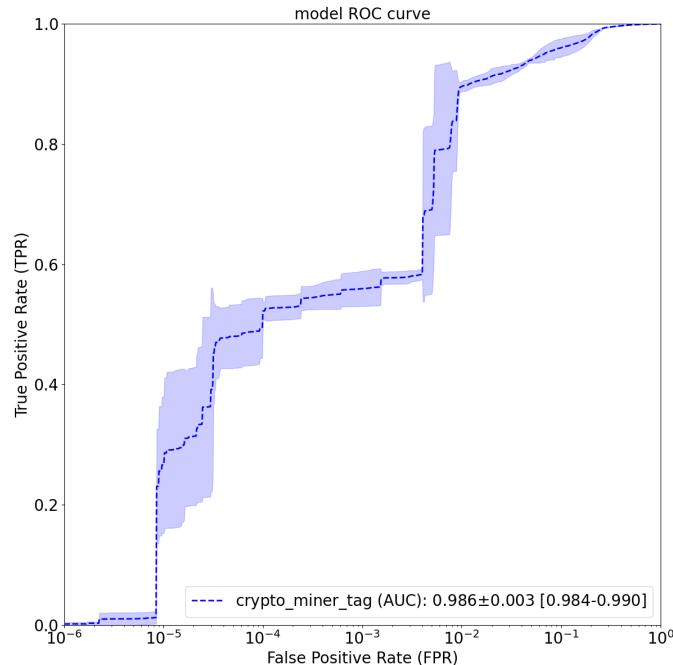


Figure 9.5: ROC curve and AUC statistics of **Joint Embedding** model for the **Crypto-miner Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

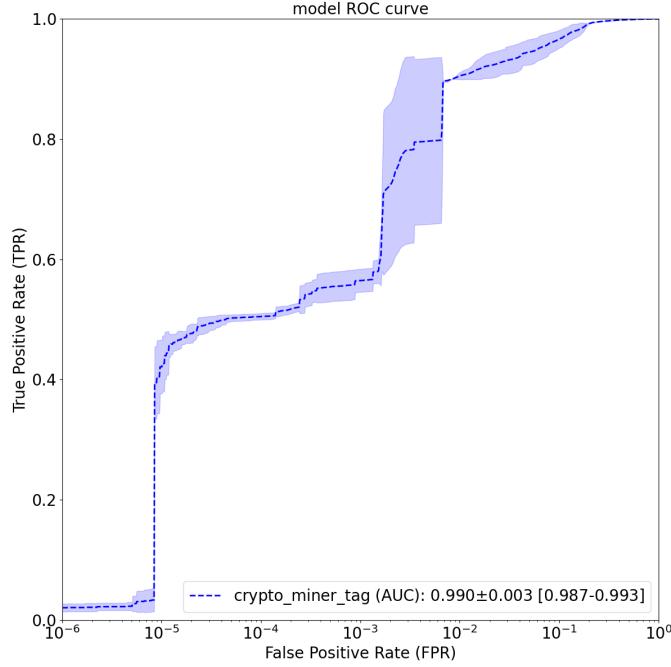


Figure 9.6: ROC curve and AUC statistics of **Proposed Model** for the **Crypto-miner Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.3 Downloader tag results

Downloader Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.070 ± 0.030	0.274 ± 0.170	0.478 ± 0.150	0.632 ± 0.034	0.883 ± 0.055
Joint Embedding	0.132 ± 0.087	0.211 ± 0.134	0.320 ± 0.224	0.635 ± 0.031	0.960 ± 0.003
Proposed Model	0.133 ± 0.101	0.364 ± 0.045	0.557 ± 0.044	0.667 ± 0.012	0.957 ± 0.006
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	6.8%	12.4%	15.1%	9.5%	63.2%
Error Reduction wrt Joint Embedding	0.1%	19.4%	34.9%	8.8%	-7.5%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	-236.7%	73.5%	70.7%	64.7%	89.1%
Std Reduction wrt Joint Embedding	-16.1%	66.4%	80.4%	61.3%	-100.0%
Accuracy					

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Downloader Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.922 ± 0.003	0.939 ± 0.014	0.955 ± 0.013	0.960 ± 0.003	0.899 ± 0.005
Joint Embedding	0.927 ± 0.007	0.934 ± 0.011	0.942 ± 0.019	0.960 ± 0.003	0.905 ± 0.000
Proposed Model	0.928 ± 0.008	0.947 ± 0.004	0.962 ± 0.004	0.963 ± 0.001	0.905 ± 0.001
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.070 ± 0.030	0.274 ± 0.170	0.478 ± 0.150	0.632 ± 0.034	0.883 ± 0.055
Joint Embedding	0.132 ± 0.087	0.211 ± 0.134	0.320 ± 0.224	0.635 ± 0.031	0.960 ± 0.003
Proposed Model	0.133 ± 0.101	0.364 ± 0.045	0.557 ± 0.044	0.667 ± 0.012	0.957 ± 0.006
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.998 ± 0.001	0.995 ± 0.003	0.975 ± 0.010	0.852 ± 0.007	0.446 ± 0.016
Joint Embedding	0.999 ± 0.001	0.990 ± 0.008	0.928 ± 0.065	0.853 ± 0.006	0.467 ± 0.001
Proposed Model	1.000 ± 0.000	0.997 ± 0.000	0.981 ± 0.001	0.859 ± 0.002	0.466 ± 0.002
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.129 ± 0.053	0.404 ± 0.195	0.627 ± 0.149	0.725 ± 0.025	0.592 ± 0.026
Joint Embedding	0.223 ± 0.132	0.328 ± 0.186	0.437 ± 0.263	0.728 ± 0.023	0.628 ± 0.001
Proposed Model	0.221 ± 0.158	0.532 ± 0.049	0.710 ± 0.036	0.751 ± 0.009	0.627 ± 0.003

Table 9.3: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Downloader Tag** prediction task at different **FPRs (False Positive Rates)**. Results were aggregated over **3** training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

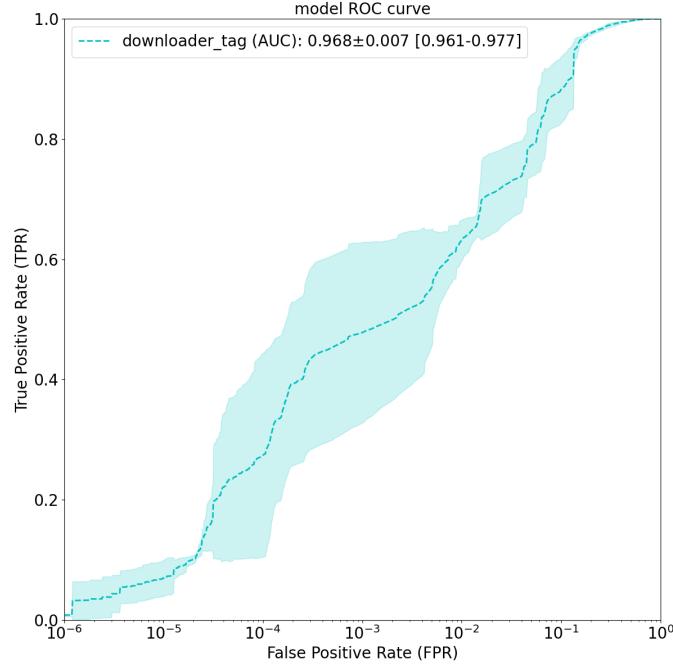


Figure 9.7: ROC curve and AUC statistics of **ALOHA** model for the **Downloader Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

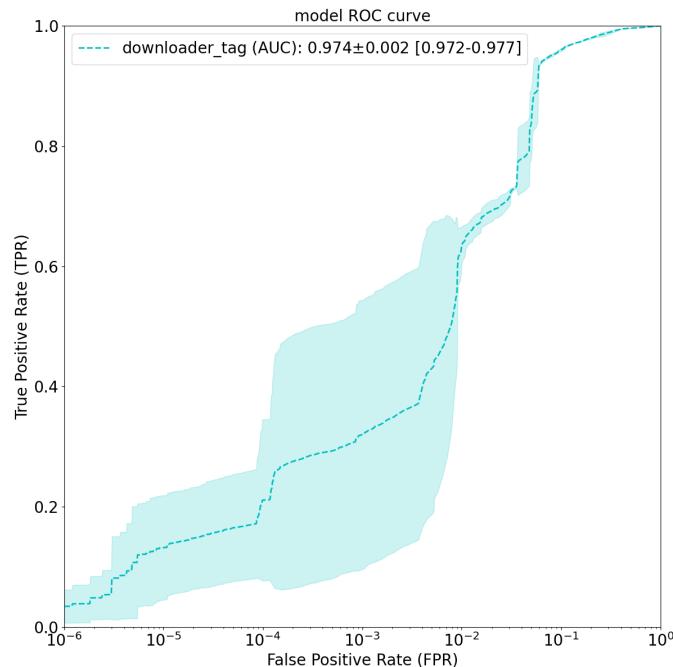


Figure 9.8: ROC curve and AUC statistics of **Joint Embedding** model for the **Downloader Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

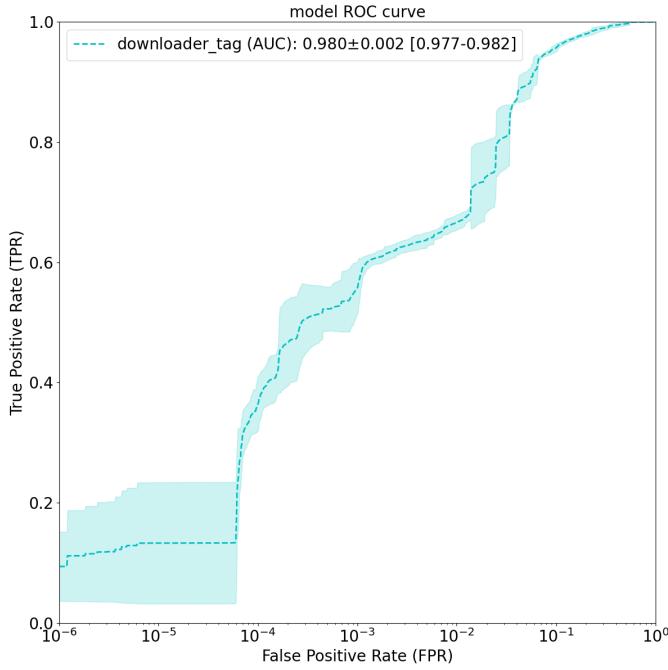


Figure 9.9: ROC curve and AUC statistics of **Proposed Model** for the **Downloader Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.4 Dropper tag results

Dropper Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.042 ± 0.022	0.073 ± 0.023	0.127 ± 0.053	0.673 ± 0.037	0.899 ± 0.017
Joint Embedding	0.088 ± 0.088	0.104 ± 0.089	0.176 ± 0.130	0.702 ± 0.031	0.904 ± 0.019
Proposed Model	0.068 ± 0.015	0.127 ± 0.047	0.151 ± 0.042	0.690 ± 0.016	0.943 ± 0.017
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	2.7%	5.8%	2.7%	5.2%	43.6%
Error Reduction wrt Joint Embedding	-2.2%	2.6%	-3.0%	-4.0%	40.6%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	31.8%	-104.3%	20.8%	56.8%	0.0%
Std Reduction wrt Joint Embedding	83.0%	47.2%	67.7%	48.4%	10.5%
Accuracy					

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Dropper Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.878 ± 0.003	0.882 ± 0.003	0.888 ± 0.007	0.950 ± 0.005	0.900 ± 0.002
Joint Embedding	0.884 ± 0.011	0.886 ± 0.011	0.894 ± 0.016	0.953 ± 0.004	0.901 ± 0.002
Proposed Model	0.881 ± 0.002	0.888 ± 0.006	0.891 ± 0.005	0.952 ± 0.002	0.905 ± 0.002
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.042 ± 0.022	0.073 ± 0.023	0.127 ± 0.054	0.673 ± 0.037	0.899 ± 0.017
Joint Embedding	0.088 ± 0.088	0.104 ± 0.089	0.174 ± 0.129	0.702 ± 0.031	0.904 ± 0.019
Proposed Model	0.066 ± 0.014	0.126 ± 0.047	0.150 ± 0.042	0.690 ± 0.016	0.943 ± 0.017
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.998 ± 0.001	0.989 ± 0.004	0.937 ± 0.031	0.908 ± 0.005	0.568 ± 0.005
Joint Embedding	0.991 ± 0.012	0.946 ± 0.069	0.880 ± 0.130	0.911 ± 0.004	0.570 ± 0.005
Proposed Model	0.999 ± 0.000	0.994 ± 0.002	0.954 ± 0.011	0.910 ± 0.002	0.580 ± 0.004
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.080 ± 0.041	0.134 ± 0.041	0.220 ± 0.085	0.772 ± 0.026	0.696 ± 0.009
Joint Embedding	0.149 ± 0.144	0.176 ± 0.146	0.274 ± 0.191	0.793 ± 0.021	0.699 ± 0.009
Proposed Model	0.124 ± 0.026	0.221 ± 0.072	0.258 ± 0.062	0.785 ± 0.011	0.718 ± 0.008

Table 9.4: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Dropper Tag** prediction task at different **FPRs (False Positive Rates)**. Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

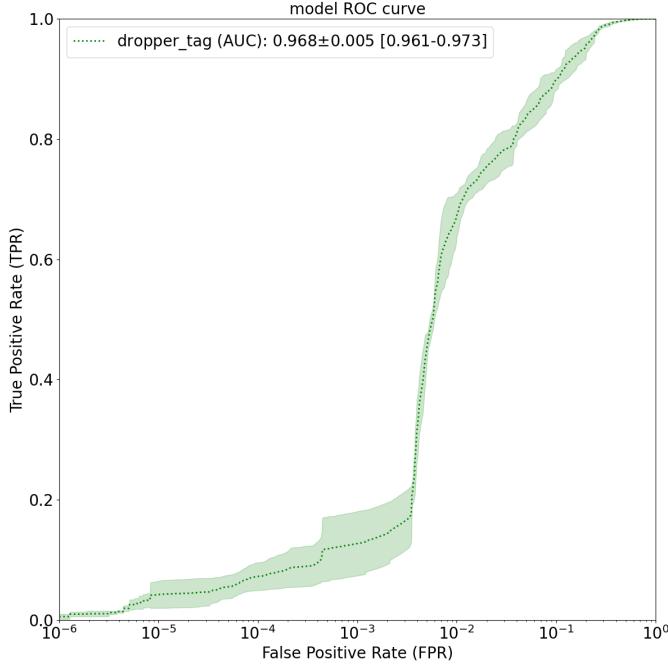


Figure 9.10: ROC curve and AUC statistics of **ALOHA** model for the **Dropper Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

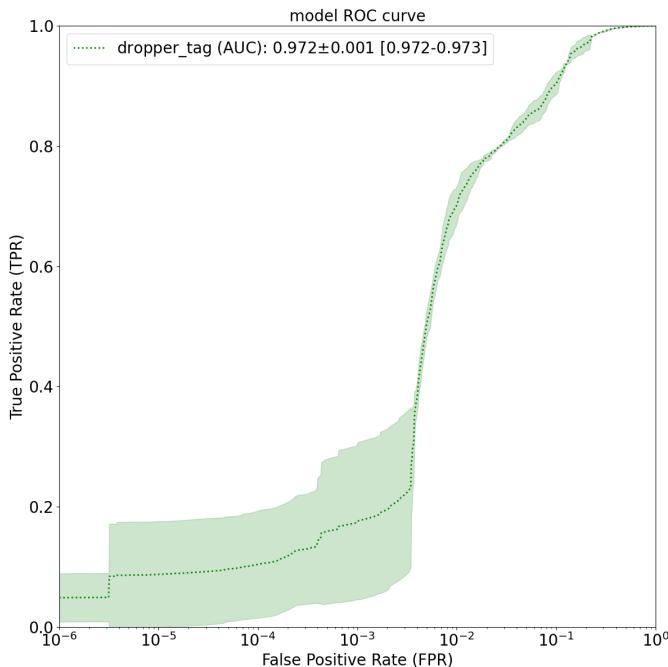


Figure 9.11: ROC curve and AUC statistics of **Joint Embedding** model for the **Dropper Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

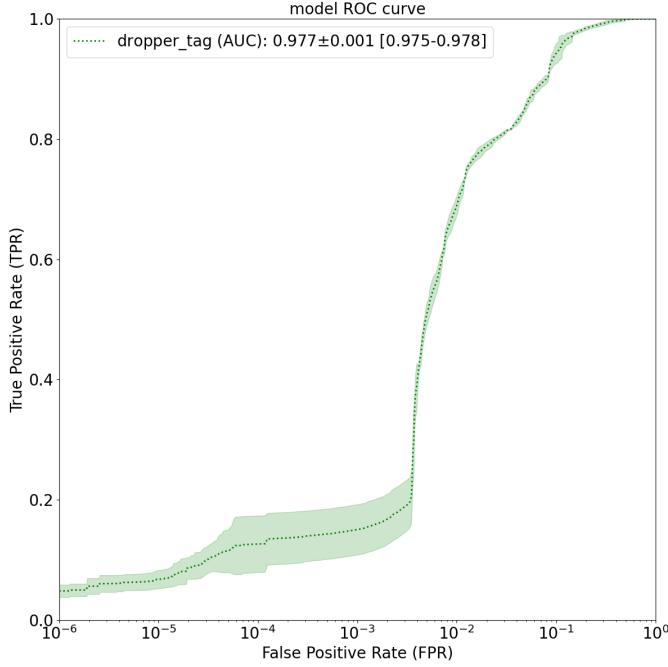


Figure 9.12: ROC curve and AUC statistics of **Proposed Model** for the **Dropper Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.5 File-infecto tag results

File-infecto Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.263 ± 0.110	0.419 ± 0.081	0.821 ± 0.013	0.875 ± 0.003	0.953 ± 0.001
Joint Embedding	0.500 ± 0.133	0.681 ± 0.084	0.822 ± 0.041	0.867 ± 0.015	0.962 ± 0.003
Proposed Model	0.275 ± 0.263	0.449 ± 0.246	0.850 ± 0.016	0.883 ± 0.014	0.959 ± 0.012
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	1.6%	5.2%	16.2%	6.4%	12.8%
Error Reduction wrt Joint Embedding	-45.0%	-72.7%	15.7%	12.0%	-7.9%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	-139.1%	-203.7%	-23.1%	-366.7%	-1100.0%
Std Reduction wrt Joint Embedding	-97.7%	-192.9%	61.0%	6.7%	-300.0%
Accuracy					
ALOHA (M/B only)	-	-	-	-	-

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File-infector Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA	0.881±0.018	0.906±0.013	0.970±0.002	0.971±0.001	0.908±0.000
Joint Embedding	0.919±0.021	0.948±0.014	0.970±0.007	0.970±0.002	0.910±0.001
Proposed Model	0.883±0.043	0.911±0.040	0.975±0.003	0.973±0.002	0.910±0.002
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.263±0.110	0.419±0.081	0.821±0.013	0.875±0.003	0.953±0.001
Joint Embedding	0.500±0.133	0.681±0.084	0.822±0.041	0.867±0.015	0.962±0.003
Proposed Model	0.275±0.264	0.448±0.247	0.850±0.016	0.883±0.014	0.959±0.012
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	1.000±0.000	0.999±0.000	0.994±0.000	0.944±0.000	0.647±0.000
Joint Embedding	1.000±0.000	0.999±0.000	0.994±0.000	0.943±0.001	0.650±0.001
Proposed Model	0.999±0.001	0.999±0.001	0.994±0.000	0.945±0.001	0.649±0.003
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.405±0.133	0.586±0.082	0.899±0.008	0.908±0.002	0.771±0.001
Joint Embedding	0.656±0.112	0.807±0.060	0.899±0.025	0.903±0.009	0.775±0.002
Proposed Model	0.369±0.302	0.582±0.218	0.916±0.009	0.913±0.008	0.774±0.006

Table 9.5: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **File-infector Tag** prediction task at different **FPRs** (*False Positive Rates*). Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

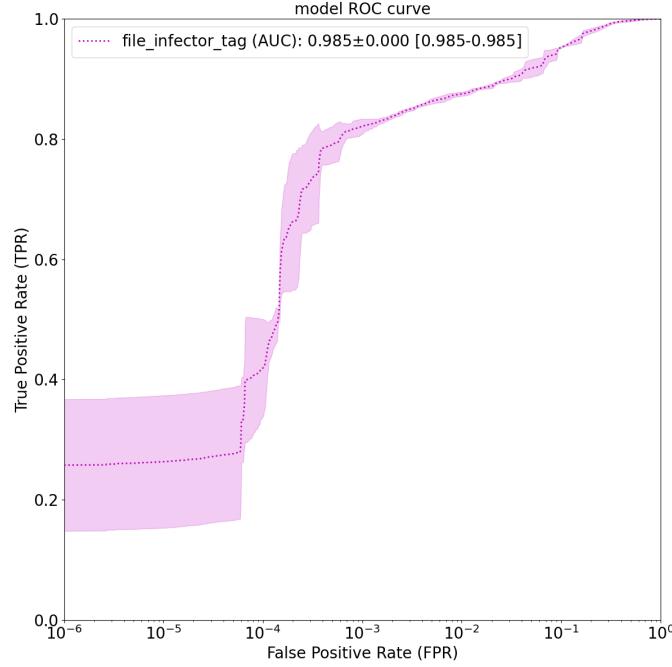


Figure 9.13: ROC curve and AUC statistics of **ALOHA** model for the **File-infector Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

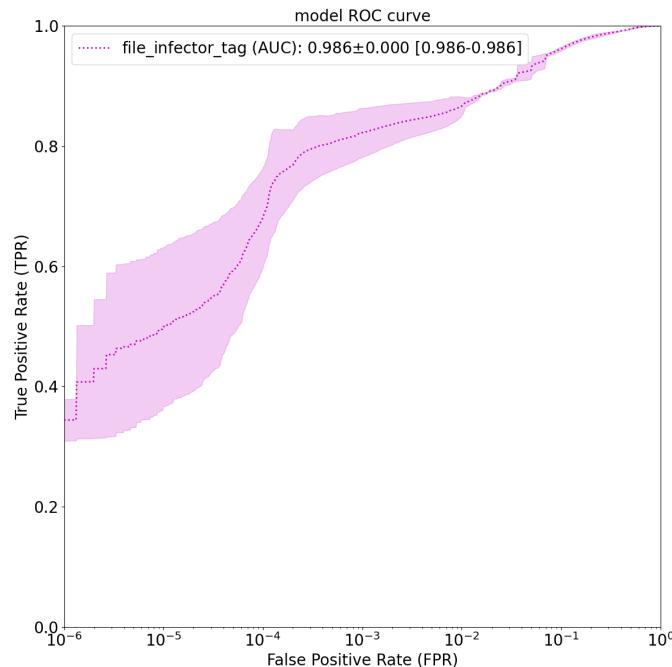


Figure 9.14: ROC curve and AUC statistics of **Joint Embedding** model for the **File-infector Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

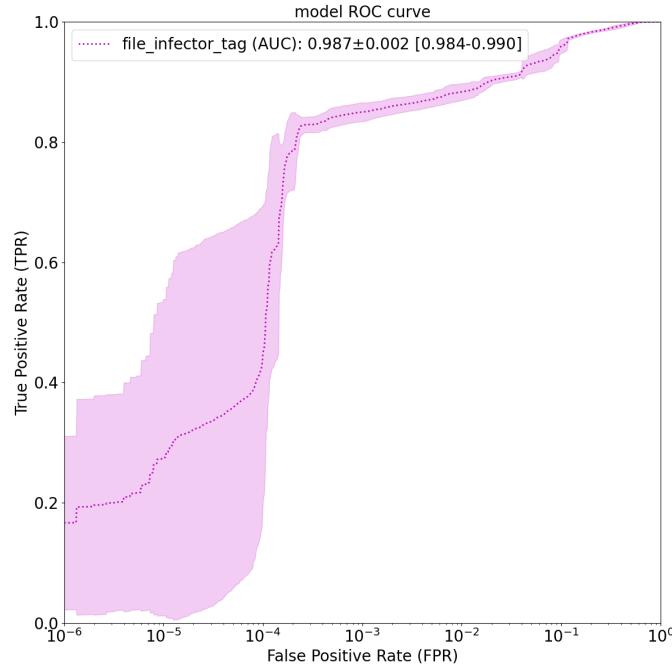


Figure 9.15: ROC curve and AUC statistics of Proposed Model for the File-infector Tag. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over 3 training runs, each with random parameter initialization.

9.1.6 Flooder tag results

Flooder Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.741 ± 0.102	0.784 ± 0.054	0.835 ± 0.009	0.876 ± 0.016	0.951 ± 0.009
Joint Embedding	0.819 ± 0.018	0.835 ± 0.013	0.835 ± 0.013	0.916 ± 0.003	0.964 ± 0.002
Proposed Model	0.831 ± 0.011	0.838 ± 0.004	0.838 ± 0.004	0.906 ± 0.007	0.964 ± 0.004
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	34.7%	25.0%	1.8%	24.2%	26.5%
Error Reduction wrt Joint Embedding	6.6%	1.8%	1.8%	-11.9%	0.0%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	89.2%	92.6%	55.6%	56.3%	55.6%
Std Reduction wrt Joint Embedding	38.9%	69.2%	69.2%	-133.3%	-100.0%
Accuracy					

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Flooder Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA (M/B only)	-	-	-	-	-
ALOHA	1.000 ± 0.000	1.000 ± 0.000	0.999 ± 0.000	0.990 ± 0.000	0.900 ± 0.000
Joint Embedding	1.000 ± 0.000	1.000 ± 0.000	0.999 ± 0.000	0.990 ± 0.000	0.900 ± 0.000
Proposed Model	1.000 ± 0.000	1.000 ± 0.000	0.999 ± 0.000	0.990 ± 0.000	0.900 ± 0.000
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.741 ± 0.102	0.784 ± 0.054	0.835 ± 0.009	0.876 ± 0.016	0.951 ± 0.009
Joint Embedding	0.819 ± 0.018	0.835 ± 0.013	0.835 ± 0.013	0.916 ± 0.003	0.964 ± 0.002
Proposed Model	0.831 ± 0.011	0.838 ± 0.004	0.838 ± 0.004	0.906 ± 0.007	0.964 ± 0.004
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.992 ± 0.001	0.934 ± 0.004	0.604 ± 0.004	0.137 ± 0.002	0.017 ± 0.000
Joint Embedding	0.996 ± 0.003	0.943 ± 0.006	0.601 ± 0.003	0.142 ± 0.000	0.017 ± 0.000
Proposed Model	0.994 ± 0.001	0.935 ± 0.004	0.603 ± 0.001	0.141 ± 0.001	0.017 ± 0.000
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.844 ± 0.070	0.851 ± 0.034	0.701 ± 0.006	0.237 ± 0.004	0.033 ± 0.000
Joint Embedding	0.898 ± 0.011	0.886 ± 0.009	0.699 ± 0.007	0.246 ± 0.001	0.034 ± 0.000
Proposed Model	0.905 ± 0.007	0.884 ± 0.002	0.701 ± 0.003	0.244 ± 0.002	0.034 ± 0.000

Table 9.6: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Flooder Tag** prediction task at different **FPRs (False Positive Rates)**. Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

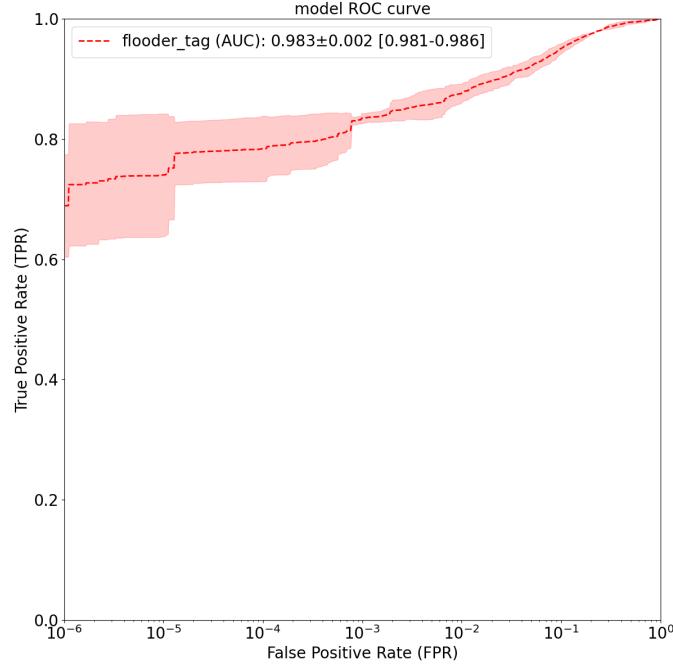


Figure 9.16: ROC curve and AUC statistics of **ALOHA** model for the **Flooder Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

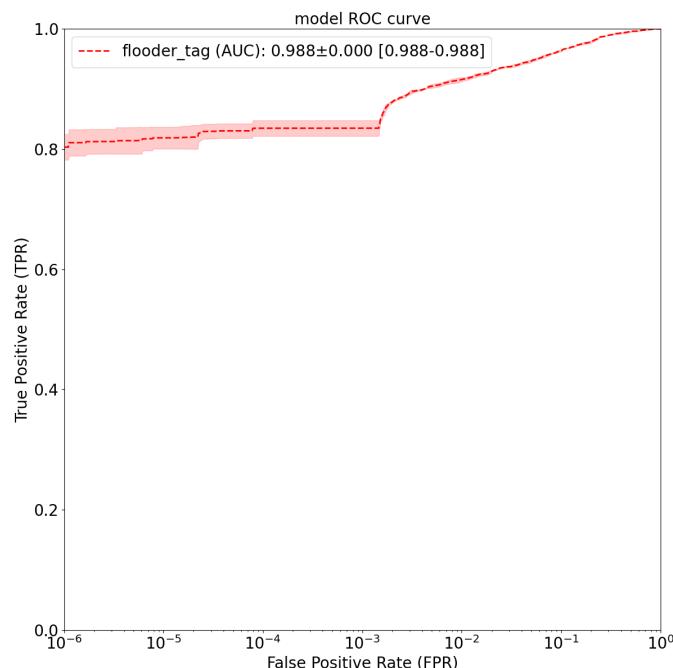


Figure 9.17: ROC curve and AUC statistics of **Joint Embedding** model for the **Flooder Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

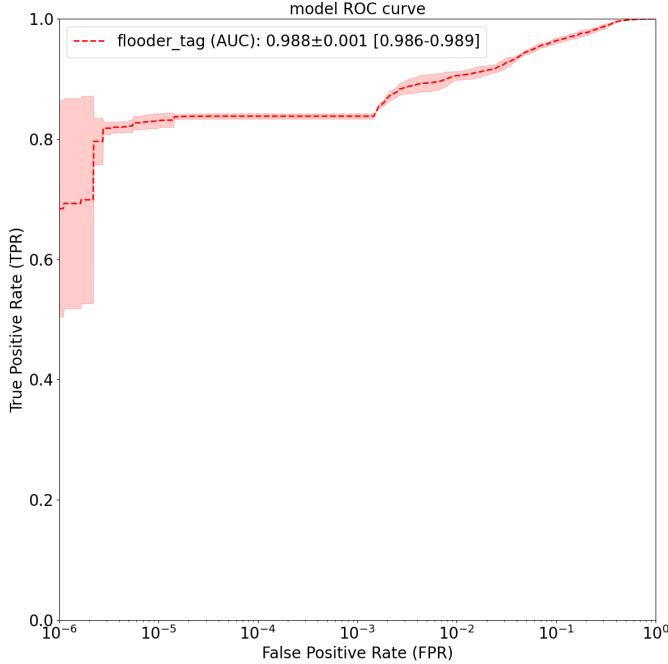


Figure 9.18: ROC curve and AUC statistics of **Proposed Model** for the **Flooder Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.7 Installer tag results

Installer Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.106 ± 0.073	0.206 ± 0.011	0.388 ± 0.059	0.717 ± 0.054	0.927 ± 0.021
Joint Embedding	0.191 ± 0.013	0.292 ± 0.026	0.506 ± 0.054	0.777 ± 0.010	0.944 ± 0.006
Proposed Model	0.208 ± 0.009	0.343 ± 0.033	0.527 ± 0.063	0.759 ± 0.042	0.926 ± 0.013
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	11.4%	17.3%	22.7%	14.8%	-1.4%
Error Reduction wrt Joint Embedding	2.1%	7.2%	4.3%	-8.1%	-32.1%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	87.7%	-200.0%	-6.8%	22.2%	38.1%
Std Reduction wrt Joint Embedding	30.8%	-26.9%	-16.7%	-320.0%	-116.7%
Accuracy					
ALOHA (M/B only)	-	-	-	-	-

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Installer Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA	0.984±0.001	0.986±0.000	0.988±0.001	0.985±0.001	0.900±0.000
Joint Embedding	0.986±0.000	0.987±0.000	0.990±0.001	0.986±0.000	0.901±0.000
Proposed Model	0.986±0.000	0.988±0.001	0.991±0.001	0.986±0.001	0.900±0.000
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.106±0.073	0.206±0.011	0.388±0.059	0.717±0.054	0.927±0.021
Joint Embedding	0.191±0.013	0.292±0.026	0.506±0.054	0.777±0.010	0.944±0.006
Proposed Model	0.208±0.009	0.343±0.033	0.527±0.063	0.759±0.042	0.926±0.013
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.955±0.059	0.974±0.001	0.873±0.017	0.564±0.019	0.144±0.003
Joint Embedding	0.997±0.000	0.981±0.002	0.901±0.010	0.584±0.003	0.146±0.001
Proposed Model	0.997±0.000	0.984±0.001	0.904±0.011	0.578±0.014	0.144±0.002
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.184±0.126	0.340±0.015	0.535±0.060	0.631±0.033	0.249±0.005
Joint Embedding	0.321±0.018	0.449±0.032	0.646±0.048	0.667±0.005	0.253±0.001
Proposed Model	0.345±0.012	0.507±0.036	0.664±0.054	0.656±0.024	0.249±0.003

Table 9.7: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Installer Tag** prediction task at different **FPRs** (*False Positive Rates*). Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

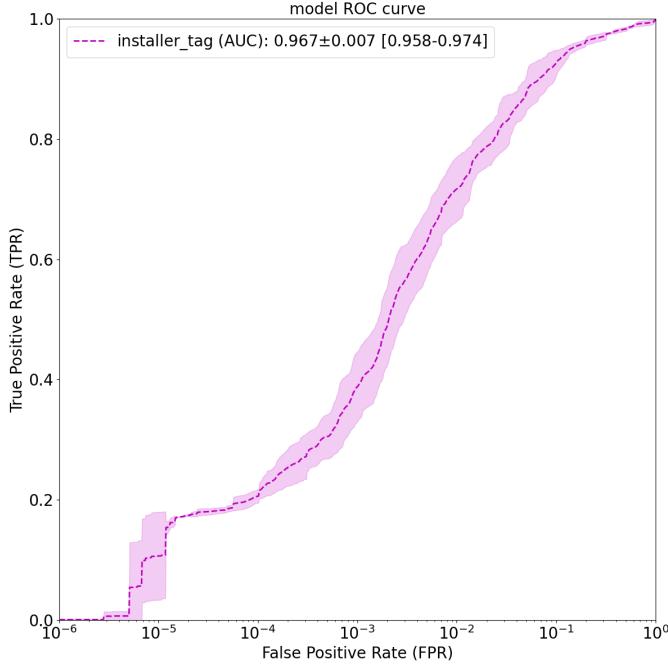


Figure 9.19: ROC curve and AUC statistics of **ALOHA** model for the **Installer Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

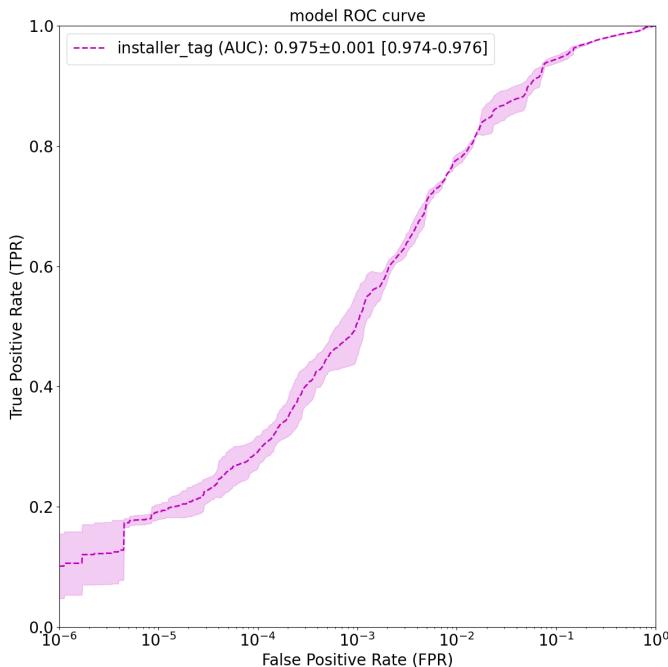


Figure 9.20: ROC curve and AUC statistics of **Joint Embedding** model for the **Installer Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

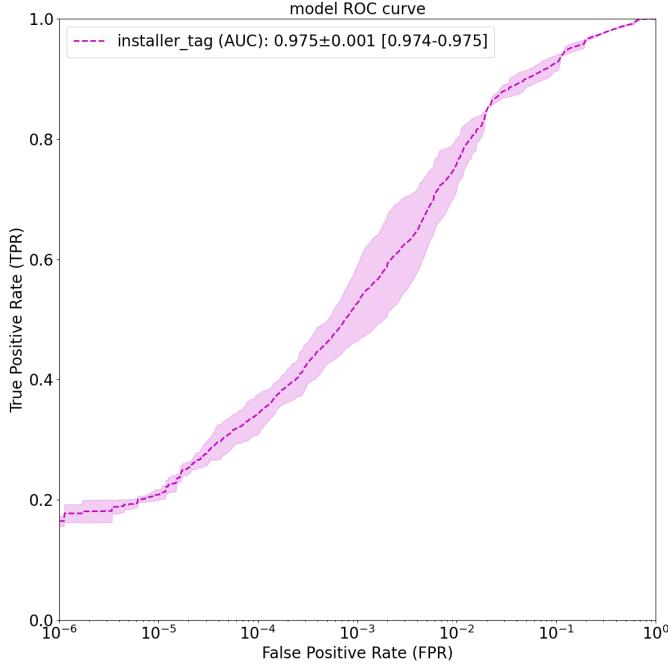


Figure 9.21: ROC curve and AUC statistics of **Proposed Model for the Installer Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.8 Packed tag results

Packed Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.012 ± 0.003	0.204 ± 0.118	0.510 ± 0.069	0.715 ± 0.003	0.942 ± 0.015
Joint Embedding	0.032 ± 0.038	0.338 ± 0.031	0.607 ± 0.029	0.764 ± 0.014	0.944 ± 0.006
Proposed Model	0.095 ± 0.053	0.414 ± 0.073	0.654 ± 0.014	0.735 ± 0.015	0.955 ± 0.008
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	8.4%	26.4%	29.4%	7.0%	22.4%
Error Reduction wrt Joint Embedding	6.5%	11.5%	12.0%	-12.3%	19.6%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	-1666.7%	38.1%	79.7%	-400.0%	46.7%
Std Reduction wrt Joint Embedding	-39.5%	-135.5%	51.7%	-7.1%	-33.3%
Accuracy					
ALOHA (M/B only)	-	-	-	-	-

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Packed Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA	0.865±0.000	0.891±0.016	0.932±0.009	0.952±0.000	0.906±0.002
Joint Embedding	0.868±0.005	0.910±0.004	0.945±0.004	0.959±0.002	0.906±0.001
Proposed Model	0.876±0.007	0.920±0.010	0.952±0.002	0.955±0.002	0.907±0.001
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.012±0.003	0.204±0.118	0.510±0.069	0.715±0.003	0.942±0.015
Joint Embedding	0.032±0.038	0.338±0.031	0.607±0.029	0.764±0.014	0.944±0.006
Proposed Model	0.095±0.054	0.414±0.073	0.654±0.014	0.735±0.015	0.955±0.008
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.994±0.002	0.993±0.006	0.987±0.002	0.919±0.000	0.598±0.004
Joint Embedding	0.994±0.007	0.998±0.000	0.990±0.000	0.924±0.001	0.599±0.002
Proposed Model	0.999±0.000	0.998±0.000	0.990±0.000	0.921±0.002	0.601±0.002
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.023±0.006	0.321±0.175	0.670±0.062	0.804±0.002	0.732±0.008
Joint Embedding	0.060±0.070	0.504±0.035	0.752±0.023	0.836±0.009	0.733±0.003
Proposed Model	0.169±0.088	0.581±0.071	0.788±0.010	0.817±0.010	0.738±0.004

Table 9.8: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Packed Tag** prediction task at different **FPRs** (*False Positive Rates*). Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

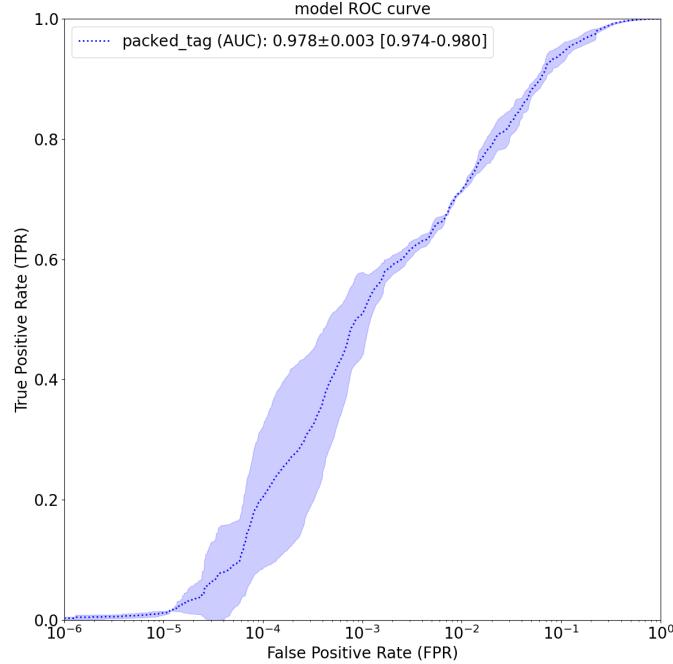


Figure 9.22: ROC curve and AUC statistics of **ALOHA** model for the **Packed Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

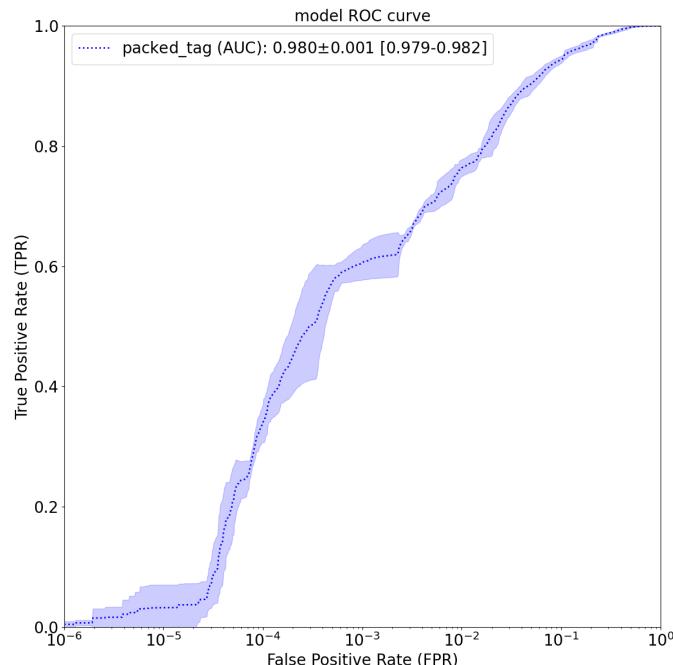


Figure 9.23: ROC curve and AUC statistics of **Joint Embedding** model for the **Packed Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

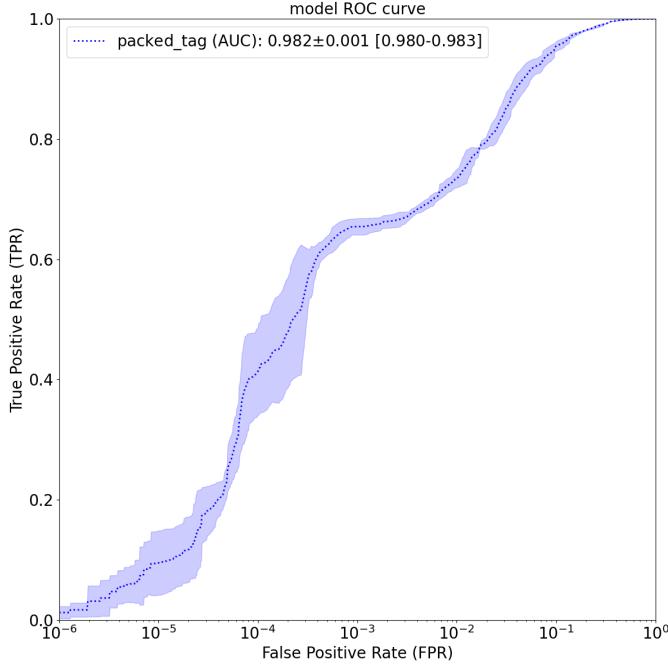


Figure 9.24: ROC curve and AUC statistics of **Proposed Model** for the **Packed Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.9 Ransomware tag results

Ransomware Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.810 ± 0.011	0.831 ± 0.013	0.837 ± 0.009	0.850 ± 0.002	0.921 ± 0.019
Joint Embedding	0.823 ± 0.006	0.843 ± 0.001	0.850 ± 0.002	0.856 ± 0.003	0.941 ± 0.013
Proposed Model	0.827 ± 0.001	0.841 ± 0.004	0.849 ± 0.002	0.856 ± 0.001	0.953 ± 0.006
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	8.9%	5.9%	7.4%	4.0%	40.5%
Error Reduction wrt Joint Embedding	2.3%	-1.3%	-0.7%	0.0%	20.3%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	90.9%	69.2%	77.8%	50.0%	68.4%
Std Reduction wrt Joint Embedding	83.3%	-300.0%	0.0%	66.7%	53.8%
Accuracy					
ALOHA (M/B only)	-	-	-	-	-

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Ransomware Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA	0.990±0.001	0.991±0.001	0.991±0.000	0.983±0.000	0.901±0.001
Joint Embedding	0.991±0.000	0.992±0.000	0.991±0.000	0.983±0.000	0.902±0.001
Proposed Model	0.991±0.000	0.992±0.000	0.991±0.000	0.983±0.000	0.903±0.000
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.810±0.011	0.831±0.013	0.837±0.009	0.850±0.002	0.921±0.019
Joint Embedding	0.823±0.006	0.843±0.001	0.850±0.002	0.856±0.003	0.941±0.013
Proposed Model	0.827±0.001	0.841±0.004	0.849±0.002	0.856±0.001	0.953±0.006
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	1.000±0.000	0.998±0.000	0.979±0.000	0.824±0.000	0.337±0.005
Joint Embedding	1.000±0.000	0.998±0.000	0.979±0.000	0.825±0.000	0.342±0.003
Proposed Model	1.000±0.000	0.998±0.000	0.979±0.000	0.825±0.000	0.345±0.001
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.895±0.007	0.906±0.008	0.902±0.006	0.837±0.001	0.493±0.007
Joint Embedding	0.903±0.004	0.914±0.001	0.910±0.001	0.840±0.002	0.502±0.005
Proposed Model	0.905±0.001	0.913±0.002	0.910±0.001	0.841±0.001	0.506±0.002

Table 9.9: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Ransomware Tag** prediction task at different **FPRs** (*False Positive Rates*). Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

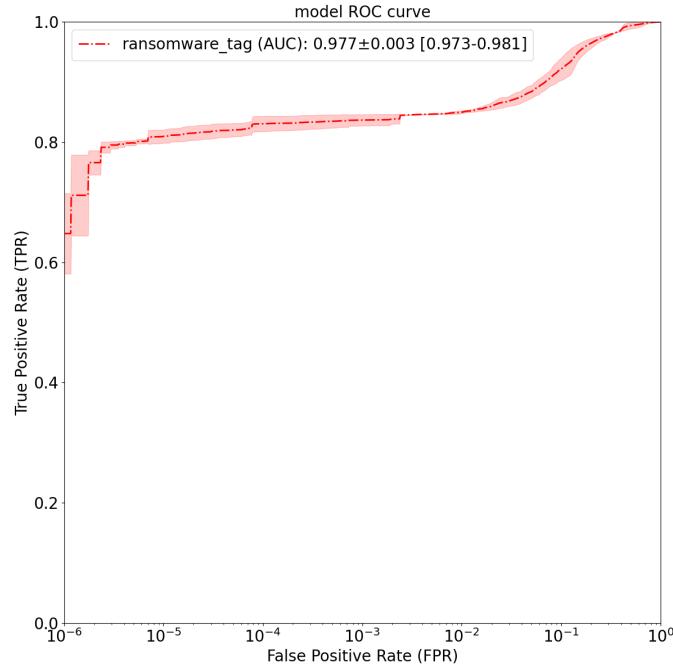


Figure 9.25: ROC curve and AUC statistics of **ALOHA** model for the **Ransomware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

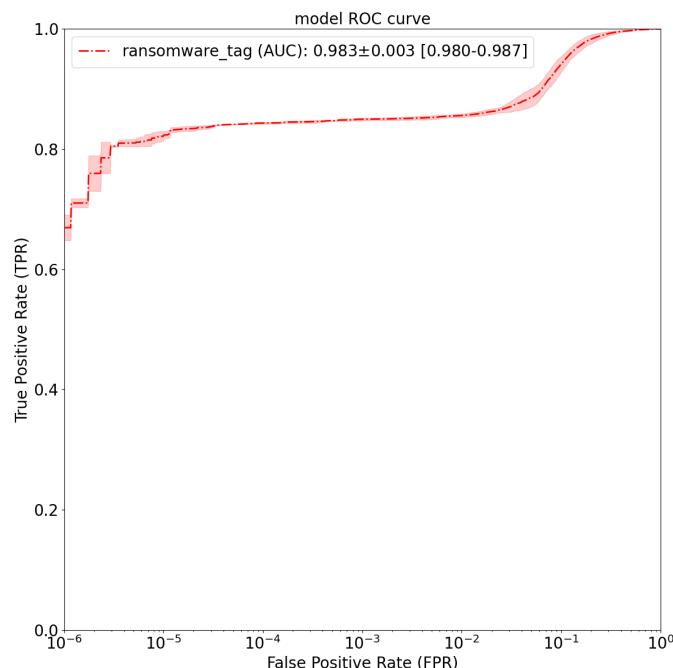


Figure 9.26: ROC curve and AUC statistics of **Joint Embedding** model for the **Ransomware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

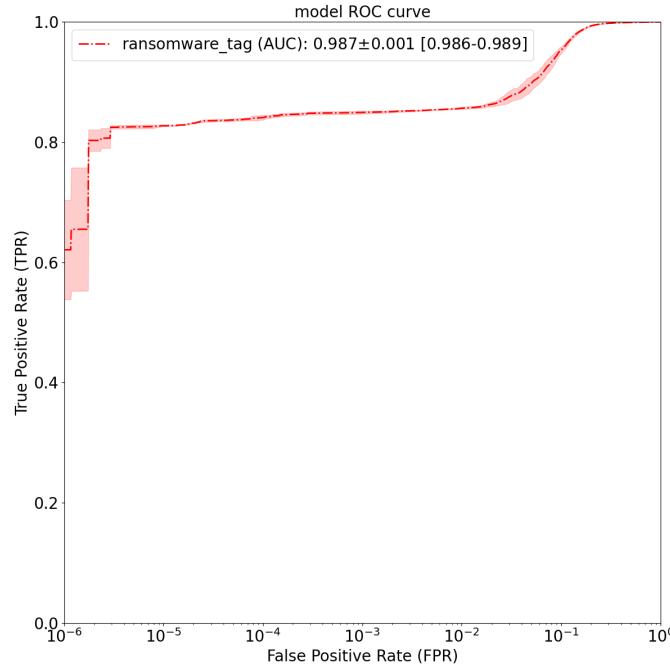


Figure 9.27: ROC curve and AUC statistics of **Proposed Model** for the **Ransomware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.10 Spyware tag results

Spyware Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.088±0.052	0.196 ± 0.128	0.239 ± 0.165	0.582 ± 0.031	0.902 ± 0.024
Joint Embedding	0.013 ± 0.004	0.214 ± 0.101	0.398 ± 0.151	0.688 ± 0.033	0.922 ± 0.009
Proposed Model	0.044 ± 0.026	0.246 ± 0.159	0.363 ± 0.060	0.660 ± 0.003	0.887 ± 0.016
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	-4.8%	6.2%	16.3%	18.7%	-15.3%
Error Reduction wrt Joint Embedding	3.1%	4.1%	-5.8%	-9.0%	-44.9%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	50.0%	-24.2%	63.6%	90.3%	33.3%
Std Reduction wrt Joint Embedding	-550.0%	-57.4%	60.3%	90.9%	-77.8%
Accuracy					

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Spyware Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.903±0.006	0.914±0.014	0.918±0.017	0.946±0.003	0.900±0.003
Joint Embedding	0.894±0.000	0.916±0.011	0.935±0.016	0.958±0.004	0.902±0.001
Proposed Model	0.898±0.003	0.919±0.017	0.931±0.006	0.955±0.000	0.899±0.002
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.088±0.052	0.196±0.128	0.239±0.166	0.582±0.031	0.902±0.024
Joint Embedding	0.012±0.004	0.214±0.101	0.398±0.151	0.688±0.033	0.922±0.009
Proposed Model	0.044±0.026	0.246±0.159	0.363±0.060	0.660±0.003	0.887±0.016
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.998±0.002	0.993±0.005	0.950±0.037	0.874±0.006	0.519±0.007
Joint Embedding	0.997±0.003	0.994±0.004	0.975±0.012	0.892±0.004	0.524±0.002
Proposed Model	0.996±0.004	0.990±0.011	0.977±0.004	0.888±0.000	0.515±0.004
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.158±0.091	0.307±0.195	0.352±0.234	0.698±0.025	0.659±0.012
Joint Embedding	0.025±0.008	0.340±0.146	0.548±0.167	0.777±0.022	0.669±0.004
Proposed Model	0.082±0.048	0.366±0.223	0.526±0.065	0.757±0.002	0.651±0.008

Table 9.10: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Spyware Tag** prediction task at different **FPRs (False Positive Rates)**. Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

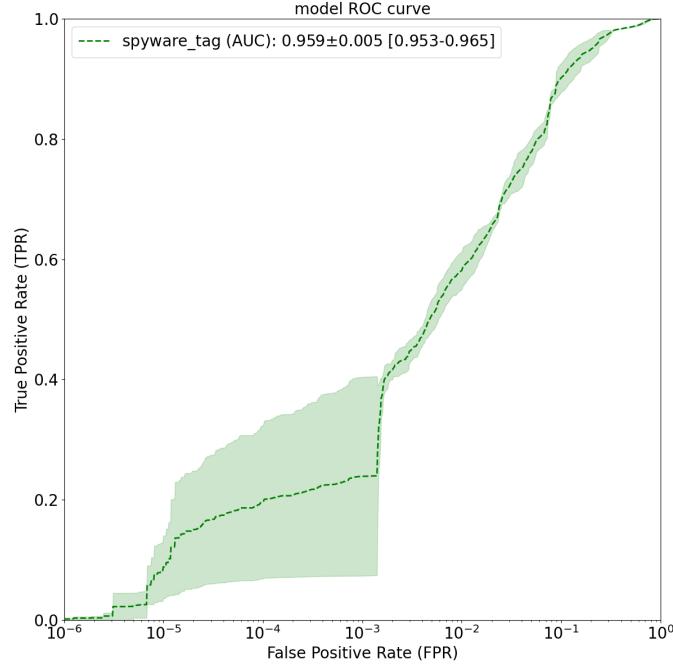


Figure 9.28: ROC curve and AUC statistics of **ALOHA** model for the **Spyware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

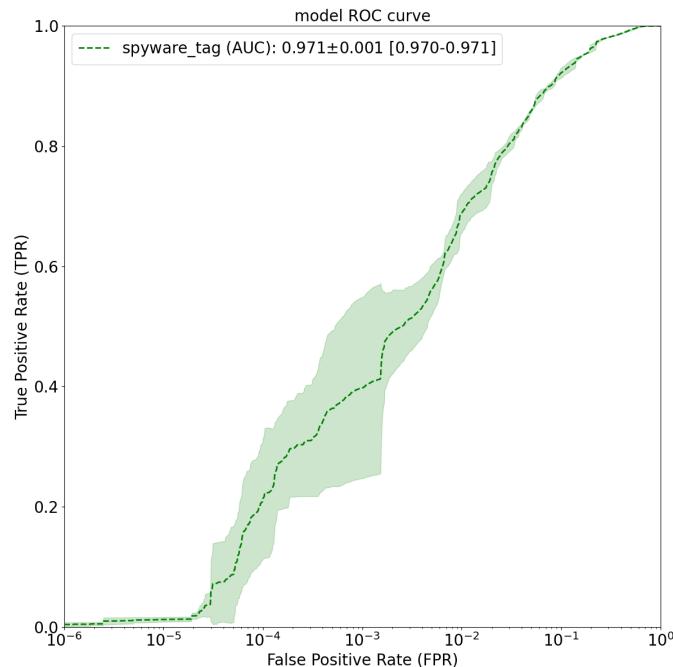


Figure 9.29: ROC curve and AUC statistics of **Joint Embedding** model for the **Spyware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

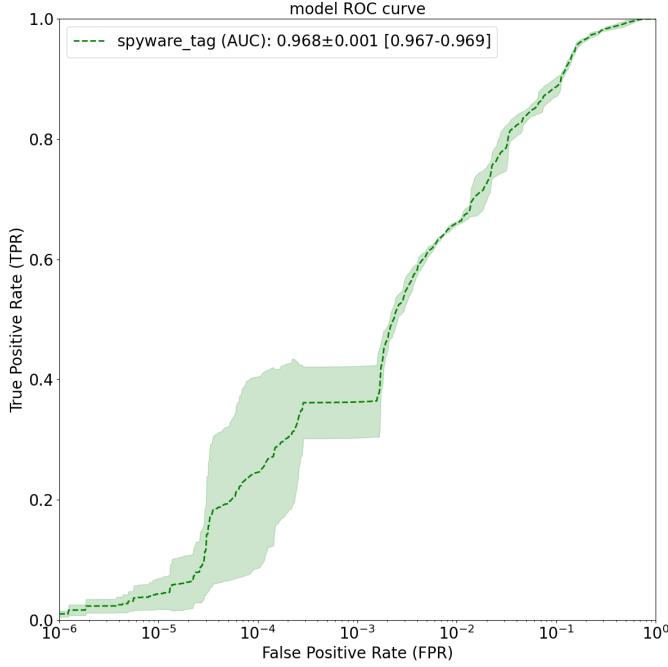


Figure 9.30: ROC curve and AUC statistics of **Proposed Model** for the **Spyware Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

9.1.11 Worm tag results

Worm Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
TPR					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.004 ± 0.003	0.062 ± 0.020	0.289 ± 0.108	0.605 ± 0.012	0.877 ± 0.061
Joint Embedding	0.117 ± 0.035	0.191 ± 0.098	0.581 ± 0.012	0.656 ± 0.004	0.770 ± 0.015
Proposed Model	0.104 ± 0.054	0.204 ± 0.119	0.508 ± 0.094	0.657 ± 0.007	0.785 ± 0.011
Error Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Error Reduction wrt ALOHA	10.0%	15.1%	30.8%	13.2%	-74.8%
Error Reduction wrt Joint Embedding	-1.5%	1.6%	-17.4%	0.3%	6.5%
Std Reduction wrt ALOHA (M/B only)	-	-	-	-	-
Std Reduction wrt ALOHA	-1700.0%	-495.0%	13.0%	41.7%	82.0%
Std Reduction wrt Joint Embedding	-54.3%	-21.4%	-683.3%	-75.0%	26.7%
Accuracy					
ALOHA (M/B only)	-	-	-	-	-

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Worm Tag	FPR				
	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}
ALOHA	0.842±0.000	0.851±0.003	0.887±0.017	0.929±0.002	0.896±0.010
Joint Embedding	0.860±0.005	0.872±0.015	0.933±0.002	0.937±0.001	0.879±0.002
Proposed Model	0.858±0.009	0.874±0.019	0.921±0.015	0.937±0.001	0.882±0.002
Recall					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.004±0.003	0.062±0.020	0.289±0.108	0.605±0.012	0.877±0.061
Joint Embedding	0.117±0.035	0.191±0.098	0.581±0.012	0.656±0.004	0.770±0.015
Proposed Model	0.104±0.054	0.204±0.119	0.508±0.094	0.657±0.007	0.785±0.011
Precision					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.980±0.016	0.990±0.004	0.980±0.006	0.919±0.002	0.622±0.017
Joint Embedding	1.000±0.000	0.998±0.001	0.991±0.000	0.925±0.000	0.592±0.005
Proposed Model	0.999±0.000	0.997±0.002	0.989±0.002	0.925±0.001	0.596±0.004
F1 Score					
ALOHA (M/B only)	-	-	-	-	-
ALOHA	0.009±0.005	0.116±0.036	0.436±0.124	0.730±0.009	0.727±0.033
Joint Embedding	0.209±0.054	0.310±0.133	0.732±0.010	0.767±0.003	0.669±0.009
Proposed Model	0.185±0.086	0.324±0.155	0.666±0.083	0.768±0.005	0.678±0.007

Table 9.11: Mean and standard deviation results (TPR, Accuracy, Recall, Precision and F1-Score) of the different models for the **Worm Tag** prediction task at different **FPRs** (*False Positive Rates*). Results were aggregated over 3 training runs with different weight initializations and minibatch orderings. Best results are shown in **bold**. Under **TPR** results are also presented the percentage reduction in mean detection error and in ROC curve standard deviation introduced by the *Proposed Model* with respect to both *ALOHA* model and *Joint Embedding*.

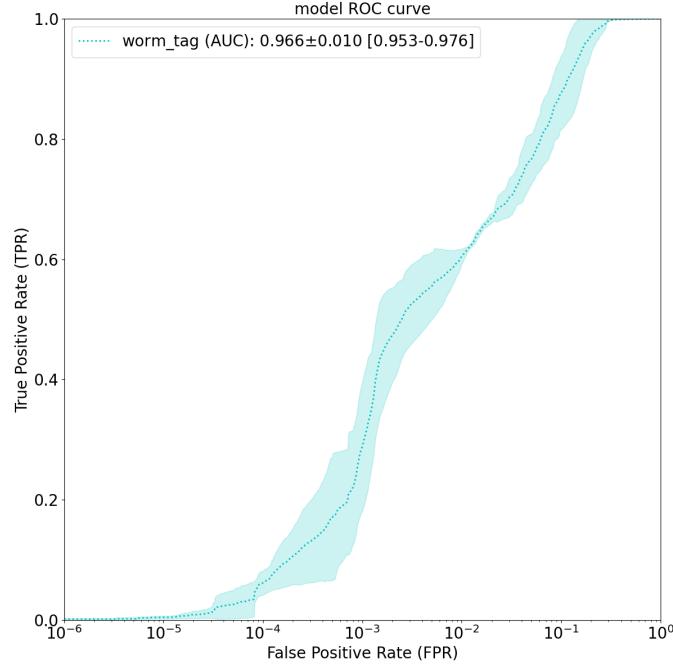


Figure 9.31: ROC curve and AUC statistics of **ALOHA** model for the **Worm Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

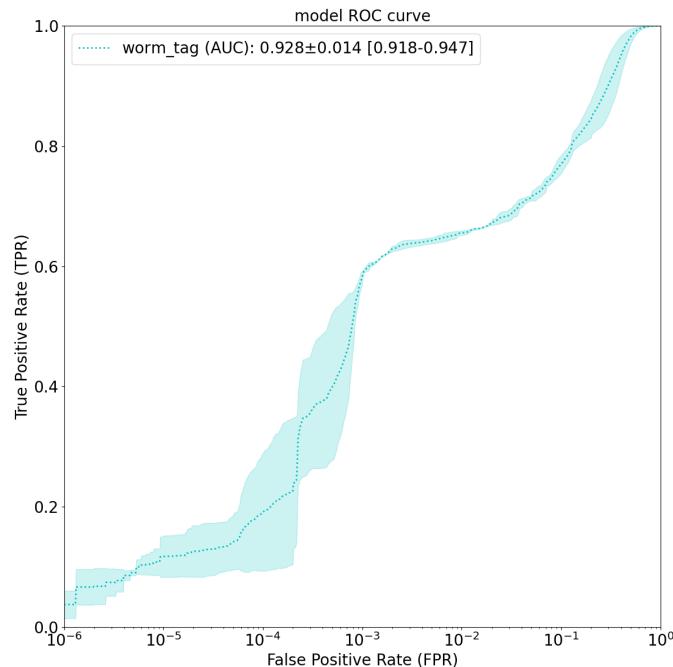


Figure 9.32: ROC curve and AUC statistics of **Joint Embedding** model for the **Worm Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

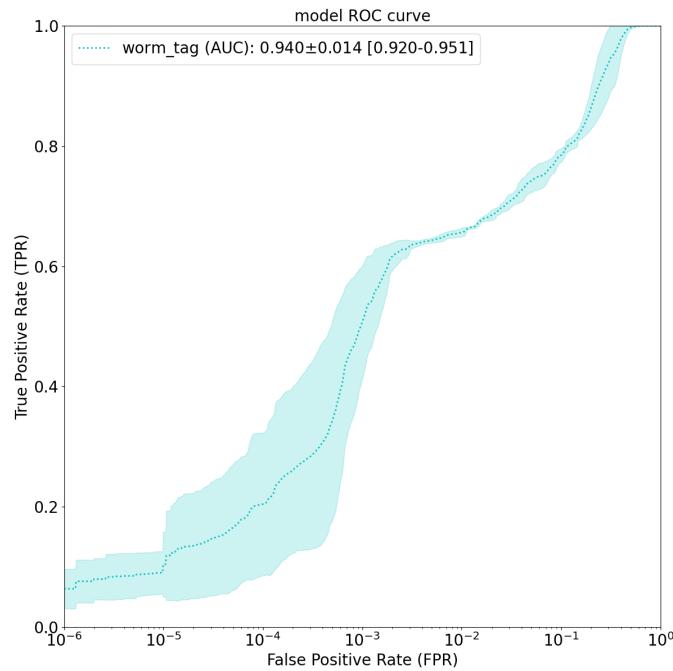


Figure 9.33: ROC curve and AUC statistics of **Proposed Model** for the **Worm Tag**. The line represents the *mean* TPR at a given FPR, while the shaded region represents the *standard deviation*. Statistics were computed over **3** training runs, each with random parameter initialization.

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