Boosting Robustness Certification
of Neural Networks

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Abstract

We present a novel approach for the certification of neural networks against adversarial perturbations which combines scalable overapproximation methods with precise (mixed integer) linear programming. This results in significantly better precision than state-of-the-art verifiers on challenging feedforward and convolutional neural networks with piecewise linear activation functions.

1 Introduction

Neural networks are increasingly applied in critical domains such as autonomous driving (Bojarski et al., 2016), medical diagnosis (Amato et al., 2013), and speech recognition (Hinton et al., 2012). However, it has been shown by Goodfellow et al. (2014) that neural networks can be vulnerable against *adversarial* attacks, i.e., imperceptible input perturbations cause neural networks to misclassify. To address this challenge and prove that a network is free of adversarial examples (usually, in a region around a given input), recent work has started investigating the use of certification techniques. Current verifiers can be broadly classified as either complete or incomplete.

Complete verifiers are exact, i.e., if the verifier fails to certify network then the network is non-robust (and vice-versa). Existing complete verifiers are based on Mixed Integer Linear Programming (MILP) (Lomuscio & Maganti, 2017; Fischetti & Jo, 2018; Dutta et al., 2018; Cheng et al., 2017) or SMT solvers (Katz et al., 2017; Ehlers, 2017). Although precise, these can only handle networks with a small number of layers and neurons. To scale, incomplete verifiers usually employ overapproximation methods and hence they are sound but may fail to prove robustness even if it holds. Incomplete verifiers use methods such as duality (Dvijotham et al., 2018), abstract interpretation (Gehr et al., 2018; Singh et al., 2018; 2019), linear approximations (Weng et al., 2018; Wong & Kolter, 2018; Zhang et al., 2018), semidefinite relaxations (Raghunathan et al., 2018), combination of linear and non-linear approximation (Xiang et al., 2017), or search space discretization (Huang et al., 2017). Incomplete verifiers are more scalable than complete ones, but can suffer from precision loss for deeper networks. In principle, incomplete verifiers can be made asymptotically complete by iteratively refining the input space (Wang et al., 2018a;b); however, at the limit, this eliminates any scalability gains and defeats the purpose of using overapproximation in the first place.

This work: boosting complete and incomplete verifiers. A key challenge then is to design a verifier which improves the precision of incomplete methods and the scalability of complete ones. In this work, we make a step towards addressing this challenge based on two key ideas: (i) a combination of state-of-the-art overapproximation techniques used by incomplete methods, including LP relaxations, together with MILP solvers, often employed in complete verifiers; (ii) a novel heuristic, which points to neurons whose approximated bounds should be refined. We implemented these ideas in a system called *RefineAI*, and showed that is is faster than state-of-the-art complete verifiers on small networks while improving precision of existing incomplete verifiers on larger networks.

The recent work of Tjeng et al. (2019) has also explored the combination of linear programming with over-approximation, though it only considers the simple interval based approximation. Our evaluation shows that for the complete verification of a 3x50 network, *RefineAI* is faster, while for the larger 9x200 network their method does not finish in days on images which *RefineAI* verifies in ≈ 14 minutes.

Figure 1: Robustness analysis of a toy example neural network using our refinement-based approach. The *DeepZ* analysis results for the neurons in the blue box are refined using MILP whereas those in green are refined using LP.

Main contributions. Our main contributions are:

- A refinement-based approach for certifying neural network robustness that combines the strengths of fast overapproximation methods with MILP solvers and LP relaxations.
- A novel heuristic for selecting neurons whose bounds should be further refined.
- A complete end-to-end implementation of our approach in a system called *RefineAI*, publicly available at https://github.com/eth-sri/eran.
- An evaluation, showing that *RefineAI* is more precise than existing state-of-the-art incomplete verifiers on larger networks and faster (while being complete) than complete verifiers on smaller networks.

2 OVERVIEW

We now demonstrate how our refinement-based approach improves the precision of a state-of-the-art certifier, thus yielding better results. Our goal is to provide an intuitive understanding of our approach; full formal details are provided in the following section.

Consider the simple fully connected feed forward neural network with ReLU activations shown in Fig. 1. There are two inputs to the network, which are both in the range $[0, 1]$. The network consists of an input layer, two hidden layers, and one output layer. Each layer consist of two neurons each. For our explanation, we separate each neuron into two parts: one represents the output of the affine transformation while the other captures the output of the ReLU activation. The weights for the affine transformation are represented by weights on the edges. The bias for each node is shown above or below it. Our goal is to certify that for any input in $[0, 1] \times [0, 1]$, the output at neuron x_{13} is greater than the output at x_{14}.

We now demonstrate how our certifier functions on this network. We assume that the analysis results after the second affine transformation are refined using a MILP formulation of the network whereas the results after the third affine transformation are refined by an LP formulation of the network. In the next section, we will explain our heuristic for selecting MILP or LP formulations for different neurons in the network. Our analysis leverages the Zonotope domain (Ghorbal et al., 2009) transformers also used in *DeepZ* (Singh et al., 2018), which is a state of the art zonotope

Figure 2: ReLU transformers, computing an affine form. Here, l_x, u_x are the original bounds, whereas l'_x, u'_x are the refined bounds. The slope of the two non-vertical parallel blue lines is $\lambda = u_x/(u_x - l_x)$ and the slope of the two non-vertical parallel green lines is $\lambda' = u'_x/(u'_x - l'_x)$. The blue parallelogram is used to compute an affine form in *DeepZ*, whereas the green parallelogram is used to compute the output of the refined ReLU transformer considered in this work.

certifier for neural network robustness. The Zonotope domain associates an affine form $\hat{x}$ with each neuron x in the network:

$$\hat{x} := [\alpha_0, \beta_0] + \sum_{i=1}^{p} [\alpha_i, \beta_i] \cdot \eta_i$$

Here, $\alpha_0, \beta_0, \alpha_i, \beta_i \in \mathbb{R}$ are real coefficients and $\eta_i \in [a_i, b_i] \subseteq [-1, 1]$ are the noise symbols, which are shared between the affine forms for different neurons. The sharing makes the domain relational and thus it is more precise than the non-relational Box domain. An abstract element in our analysis is an intersection between a Zonotope (given as a list of affine forms) and a bounding box. Thus, for each neuron x, we keep the affine form $\hat{x}$ and an interval $[l_x, u_x]$.

First layer. Our analysis starts by setting

$$\hat{x}_1 = 0.5 + 0.5 \cdot \eta_1, l_1 = 0, u_1 = 1$$

and

$$\hat{x}_2 = 0.5 + 0.5 \cdot \eta_2, l_2 = 0, u_2 = 1,$$

representing the inputs at x_1 and x_2 respectively. Next, an affine transformation is applied on the inputs resulting in the output

$$\hat{x}_3 = \hat{x}_1 + \hat{x}_2 = 1 + 0.5 \cdot \eta_1 + 0.5 \cdot \eta_2, l_3 = 0, u_3 = 2$$

and

$$\hat{x}_4 = \hat{x}_1 - \hat{x}_2 = 0.5 \cdot \eta_1 - 0.5 \cdot \eta_2, l_4 = -1, u_4 = 1.$$

The Zonotope affine transformer is *exact* for this transformation so we do not need any refinement here. Next, the Zonotope ReLU transformer is applied. We note that as $l_3 \geq 0$, the neuron x_3 provably takes only non-negative values. Thus, the ReLU Zonotope transformer outputs $\hat{x}_5 = \hat{x}_3$ and we set $l_5 = l_3, u_5 = l_3$ which is the exact result. For x_4, $l_4 < 0$ and $u_4 > 0$ and thus neuron x_4 can take both positive and negative values. The corresponding output does not have a closed affine form and hence the approximation in blue shown in Fig. 2 is used to compute the result. This approximation minimizes the area of the result in the input-output plane and introduces a new noise symbol $\eta_3 \in [-1, 1]$. The result is

$$\hat{x}_6 = 0.25 + 0.25 \cdot \eta_1 - 0.25 \cdot \eta_2 + 0.25 \cdot \eta_3, l_6 = -0.5, u_6 = 1.$$

Note that the Zonotope approximation for x_6 from Fig. 2 permits negative values whereas x_6 can only take non-negative values in the concrete. This approximation typically accumulates as the analysis progresses deeper into the network, resulting in overall imprecision and failure to prove properties that may actually hold.

MILP-based refinement at second layer. Next, the analysis handles the second affine transformation and computes

$$\hat{x}_7 = \hat{x}_5 - \hat{x}_6 + 1 = 1.75 + 0.25 \cdot \eta_1 + 0.75 \cdot \eta_2 - 0.25 \cdot \eta_3, l_7 = 0.5, u_7 = 3$$

and

$$\hat{x}_8 = \hat{x}_5 - \hat{x}_6 - 1 = -0.25 + 0.25 \cdot \eta_1 + 0.75 \cdot \eta_2 - 0.25 \cdot \eta_3, l_8 = -1.5, u_8 = 1.$$

Here, x_7 is provably positive, whereas x_8 can take both positive and negative values. Due to the approximation for x_6, the bounds for x_7 and x_8 are imprecise. Notice that the Zonotope ReLU transformer for x_8 will introduce more imprecision and although the ReLU transformer for provably positive inputs such as x_7 does not lose precision with respect to the input, it still propagates the imprecision in the computation of the abstract values for x_7.

Thus, to reduce precision loss, in our method we refine the bounds for both x_7 and x_8 by formulating the network up to the second affine transformation as a MILP instance based on a formulation from Tjeng et al. (2019) and compute bounds for x_7 and x_8, respectively. The MILP solver improves the lower bounds for x_7 and x_8 to 1 and -1, respectively, which then updates the corresponding lower bounds in our abstraction, i.e., $l_7 = 1$ and $l_8 = -1$. Next, the ReLU transformer is applied. Since x_7 is provably positive, we get $\hat{x}_9 = \hat{x}_7, l_9 = l_7, u_9 = u_7$.

We note that x_8 can take both positive and negative values and is therefore approximated. However, the ReLU transformer now uses the refined bounds instead of the original bounds and thus the approximation shown in green from Fig. 2 is used. This approximation has smaller area in the input-output plane compared to the blue one and thus reduces the approximation error. The result is

$$\hat{x}_{10} = 0.125 + 0.125 \cdot \eta_1 + 0.375 \cdot \eta_2 - 0.125 \cdot \eta_3 + 0.25 \cdot \eta_4, l_8 = -0.5, u_8 = 1.$$

LP-based refinement at final layer. Continuing with the analysis, we now process the final affine transformation, which yields

$$\hat{x}_{11} = 4.875 + 0.625 \cdot \eta_1 + 1.875 \cdot \eta_2 - 0.625 \cdot \eta_3 + 0.25 \cdot \eta_4, l_{11} = 1.75, u_{11} = 8.25$$

and

$$\hat{x}_{12} = 2.625 + 0.125 \cdot \eta_1 + 0.375 \cdot \eta_2 + -0.125 \cdot \eta_3 - 0.25 \cdot \eta_4, l_{12} = 1.75, u_{12} = 3.5.$$

Due to the approximations from previous layers, the computed values can be imprecise. We note that as the analysis proceeds deeper into the network, refining bounds with MILP becomes expensive. Thus, we refine the bounds by encoding the network up to the third affine transformation using the faster LP relaxation of the network based on Ehlers (2017) and compute the bounds for x_{11} and x_{12}, respectively. This leads to the following results: $l_{11} = 3.25$, $l_{12} = 2$, and $u_{12} = 3$. As both x_{13} and x_{14} are provably positive, the subsequent ReLU transformations set $\hat{x}_{13} = \hat{x}_{11}, l_{13} = l_{11}, u_{13} = u_{11}$ and $\hat{x}_{14} = \hat{x}_{12}, l_{14} = l_{12}, u_{14} = u_{12}$.

Proving robustness. Since the lower bound l_{13} for x_{13} is greater than the upper bound u_{14} for x_{14}, our analysis can prove that the given neural network provides the same label for all inputs in $[0, 1] \times [0, 1]$ and is thus robust. In contrast, *DeepZ* without our refinement would compute

$$\hat{x}_{13} = 4.95 + 0.6 \cdot \eta_1 + 1.8 \cdot \eta_2 - 0.6 \cdot \eta_3 + 0.3 \cdot \eta_4, l_{13} = 1.65, u_{13} = 8.25$$

and

$$\hat{x}_{14} = 2.55 + 0.15 \cdot \eta_1 + 0.45 \cdot \eta_2 - 0.15 \cdot \eta_3 - 0.3 \cdot \eta_4, l_{14} = 1.5, u_{14} = 3.6.$$

As a result, *DeepZ* fails to prove that x_{13} is greater than x_{14}, and thus fails to prove robustness.

Generalization to other abstractions. We note that our refinement-based approach is not restricted to the Zonotope domain. It can be easily extended for refining the results computed by other abstractions such as Polyhedra (Singh et al., 2017) or the abstraction used in *DeepPoly* (Singh et al., 2019). For example, the ReLU transformer in Singh et al. (2019) also depends on the bounds of input neurons and thus it will benefit from the precise bounds computed using our refinement. Since *DeepPoly* often produces more precise results than *DeepZ*, we believe that a straightforward extension of this work to *DeepPoly* will further improve certification results.

3   Our Approach

In this section, we describe our approach in more formal terms. As in the previous section, we will consider affine transformations and ReLU activations as separate layers. As illustrated earlier, the key idea will be to combine abstract interpretation with exact and inexact MILP formulation of the network which is then solved using a solver, in order to compute more precise results for neuron bounds. We begin by describing the ingredients of abstract interpretation.

Our approach requires an abstract domain $\mathbb{A}_n$ (i.e., some set whose elements can be encoded symbolically) such as Interval, Zonotope, the abstraction in *DeepPoly*, or Polyhedra. An abstract domain has a bottom element $\perp \in \mathbb{A}_n$ as well as the following components:

- A (potentially non-computable) concretization function $\gamma_n \colon \mathbb{A}_n \to \mathcal{P}(\mathbb{R}^n)$ that associates with each abstract element $a \in \mathbb{A}_n$ a set of concrete points from $\mathbb{R}^n$. We have, $\gamma_n(\perp) = \emptyset$.

- An abstraction function $\alpha_n \colon \mathcal{P}(\mathbb{R}^n) \to \mathbb{A}_n$, where $\mathbb{X} \subseteq \gamma_n(\alpha_n(\mathbb{X}))$ for all $\mathbb{X} \in \mathcal{P}(\mathbb{R}^n)$. We assume that $\alpha_n(\prod_i [l_i, u_i])$ is computable for all $\boldsymbol{l}, \boldsymbol{u} \in \mathbb{R}^n$, where

$$\prod_i [l_i, u_i] = \{\boldsymbol{x} \in \mathbb{R}^n \mid l_i \le x_i \le u_i\}.$$

- A bounding box function $\iota_n \colon \mathbb{A}_n \to \mathbb{R}^n \times \mathbb{R}^n$, where $\gamma_n(a) \subseteq \prod_i [l_i, y_i]$ for $(\boldsymbol{l}, \boldsymbol{u}) = \iota_n(a)$ for all $a \in \mathbb{A}_n$.

- A meet operation $a \sqcap L$ for each $a \in \mathbb{A}_n$ and linear constraints L over n real variables, where $\{x \in \gamma_n(a) \mid L(x)\} \subseteq \gamma_n(a \sqcap L)$.

- An affine abstract transformer $T^{\#}_{\boldsymbol{x} \mapsto \boldsymbol{Ax}+\boldsymbol{b}} \colon \mathbb{A}_m \to \mathbb{A}_n$ for each transformation of the form $(\boldsymbol{x} \mapsto \boldsymbol{Ax} + \boldsymbol{b}) \colon \mathbb{R}^m \to \mathbb{R}^n$, where

$$\{\boldsymbol{Ax} + \boldsymbol{b} \mid \boldsymbol{x} \in \gamma_n(a)\} \subseteq \gamma_n(T^{\#}_{\boldsymbol{x} \mapsto \boldsymbol{Ax}+\boldsymbol{b}}(a))$$

for all $a \in \mathbb{A}_m$.

- A ReLU abstract transformer $T^{\#}_{\mathrm{ReLU}|_{\prod_i [l_i, u_i]}} \colon \mathbb{A}_n \to \mathbb{A}_n$, where

$$\{\mathrm{ReLU}(\boldsymbol{x}) \mid \boldsymbol{x} \in \gamma_n(a) \cap \prod_i [l_i, u_i]\} \subseteq T^{\#}_{\mathrm{ReLU}|_{\prod_i [l_i, u_i]}}(a)$$

for all abstract elements $a \in \mathbb{A}_n$ and for all lower and upper bounds $\boldsymbol{l}, \boldsymbol{u} \in \mathbb{R}^n$ on input activations of the ReLU operation.

Abstract interpretation.   As first shown by Gehr et al. (2018), any such abstract domain induces a method for robustness certification of neural networks with ReLU activations.

For example, assume that we want to certify that a given neural network $f \colon \mathbb{R}^m \to \mathbb{R}^n$ considers class i more likely than class j for all inputs $\bar{\boldsymbol{x}}$ with $||\bar{\boldsymbol{x}} - \boldsymbol{x}||_\infty \le \epsilon$ for a given $\boldsymbol{x}$ and ϵ. We can first use the function α_m to compute a symbolic overapproximation of the set of possible input activations $\bar{\boldsymbol{x}}$, namely

$$a_{\mathrm{in}} = \alpha_m(\{\bar{\boldsymbol{x}} \in \mathbb{R}^m \mid ||\bar{\boldsymbol{x}} - \boldsymbol{x}||_\infty \le \epsilon\}).$$

Given that the neural network can be written as a composition of affine functions and ReLU layers, we can then propagate the abstract element a_{in} through the corresponding abstract transformers to obtain a symbolic overapproximation a_{out} of the concrete outputs of the neural network.

For example, if the neural network $f(\boldsymbol{x}) = \boldsymbol{A}' \cdot \mathrm{ReLU}(\boldsymbol{Ax} + \boldsymbol{b}) + \boldsymbol{b}'$ has a single hidden layer with h hidden neurons, we first compute $a' = T^{\#}_{\boldsymbol{x} \mapsto \boldsymbol{Ax}+\boldsymbol{b}}(a_{\mathrm{in}})$, which is a symbolic overapproximation of the input to the ReLU activation function. We then compute $(\boldsymbol{l}, \boldsymbol{u}) = \iota_h(a')$ to obtain opposite corners of a bounding box of all possible ReLU input activations, such that we can apply the ReLU abstract transformer:

$$a'' = T^{\#}_{\mathrm{ReLU}|_{\prod_i [l_i, u_i]}}(a').$$

Finally, we apply the affine abstract transformer again to obtain $a_{\mathrm{out}} = T^{\#}_{\boldsymbol{x} \mapsto \boldsymbol{A}'\boldsymbol{x}+\boldsymbol{b}'}(a'')$. Using our assumptions, we can conclude that the set $\gamma_n(a_{\mathrm{out}})$ contains all output activations that f can possibly produce when given any of the inputs $\bar{\boldsymbol{x}}$. Therefore, if $a_{\mathrm{out}} \sqcap (x_i \le x_j) = \perp$, we have proved the property: for all $\bar{\boldsymbol{x}}$, the neural network considers class i more likely than class j.

Incompleteness. While this approach is sound (i.e., whenever we prove the property, it actually holds), it is incomplete (i.e., we might not prove the property, even if it holds), because the abstract transformers produce a superset of the set of concrete outputs that the corresponding concrete transformers produce. This is in particular quite imprecise for deep neural networks, because the overapproximations introduced in each layer accumulate.

Refinement. To combat spurious overapproximation, we use mixed integer linear programming (MILP) to compute refined lower and upper bounds l, u after applying each affine abstract transformer (except for the first layer). We then refine the abstract element using the meet operator and the linear constraints $l_i \leq x_i \leq u_i$ for all input activations i, i.e., we replace the current abstract element a by $a' = a \sqcap (\bigwedge_i l_i \leq x_i \leq u_i)$, and continue analysis with the refined abstract element.

Furthermore, and importantly, for ReLU, we obtain a refined abstract transformer using the new (and tighter) lower and upper bounds.

To enable refinement, we augment each abstract element with variables and constraints. There is one variable for each neuron, and the constraints fully capture the behaviour of the neural network up to the last layer whose abstract transformer has been executed. We write $x_i^{(k)}$ to denote the variable corresponding to the activation of the i-th neuron in the k-th layer, where the input layer has $k = 0$. Similarly, we will write $l_i^{(k)}$ and $u_i^{(k)}$ to denote the best derived lower and upper bounds for this neuron.

From the input layer, we get constraints of the form $l_i^0 \leq x_i^{(0)} \leq u_i^0$, from affine layers, we get constraints of the form $x_i^{(k)} = \sum_j a_{ij}^{(k-1)} x_j^{(k-1)} + b_i^{(k-1)}$ and from ReLU layers we get constraints of the form $x_i^{(k)} = \max(0, x_i^{(k-1)})$.

MILP. Let $\varphi^{(k)}$ denote the conjunction of all constraints up to and including those from layer k. To obtain the best possible lower and upper bounds for layer k with p neurons, we need to solve the following $2 \cdot p$ optimization problems:

$$l_i^{(k)} = \min_{\substack{x_1^{(0)}, \ldots, x_p^{(k)} \\ \text{s.t. } \varphi^{(k)}(x_1^{(0)}, \ldots, x_p^{(k)})}} x_i^{(k)}, \text{ for } i = 1, \ldots, p,$$

$$u_i^{(k)} = \max_{\substack{x_1^{(0)}, \ldots, x_p^{(k)} \\ \text{s.t. } \varphi^{(k)}(x_1^{(0)}, \ldots, x_p^{(k)})}} x_i^{(k)}, \text{ for } i = 1, \ldots, p.$$

As was shown by Tjeng et al. (2019), those optimization problems can be encoded exactly as MILPs, such that the bounds can be computed with off-the-shelf MILP solvers. Note that the encoding of subsequent layers depends on the bounds computed in previous layers, where tighter bounds produce simpler MILP instances.

LP relaxation. While not introducing any approximation, unfortunately, current MILP solvers do not scale to larger deep neural networks. It becomes increasingly more expensive to refine bounds with the MILP-based formulation as the analysis proceeds deeper into the network. However, for soundness it is not crucial that the produced bounds are the best possible: for example, plain abstract interpretation uses sound bounds produced by the bounding box function ι instead. Therefore, for deeper layers in the network, we explore the trade-off between precision and scalability by also considering an intermediate method, which is faster than exact MILP, but also more precise than abstract interpretation. We relax the constraints in $\varphi^{(k)}$ in the same way as Ehlers (2017) to obtain a set of weaker linear constraints $\varphi_{\text{LP}}^{(k)}$. We then use the solver to solve the relaxed optimization problems that are constrained by $\varphi_{\text{LP}}^{(k)}$ instead of $\varphi^{(k)}$, producing possibly looser bounds $l^{(k)}$ and $u^{(k)}$. Note that the encoding of subsequent layers depends on the bounds computed in previous layers, where tighter bounds reduce the amount of newly introduced approximation.

Anytime MILP relaxation. MILP solvers usually provide the option to provide an explicit timeout after which the solver must terminate. In return, the solver may not be able to solve the instance

exactly, but it will instead provide lower and upper bounds on the objective function in a best-effort fashion. This provides another way to compute sound but inexact bounds $\boldsymbol{l}^{(k)}$ and $\boldsymbol{u}^{(k)}$.

In practice, we choose a fraction $\alpha \in (0, 1]$ of neurons in a given layer k and compute bounds for them using MILP with a timeout T in a first step. In the second step, for a fraction $\delta \in [0, 1 - \alpha]$ of neurons in the layer, we set the timeout to $\beta \cdot \overline{T}$, where $\overline{T}$ is the average time taken by the MILP solver to solve one of the instances from the first step and $\beta \in [0, 1]$ is a parameter.

Neuron selection heuristic. To select the α-fraction of neurons for the first step of the anytime MILP relaxation for the k-th layer, we rank the neurons. If the next layer is a ReLU layer, we first ignore all neurons whose activations can be proven to be non-positive using abstract interpretation (i.e., using the bounds produced by ι), because in this case it is already known that ReLU will map the activation to 0. The remaining neurons are ordered in up to two different ways, once by width (i.e. neuron i has key $u_i^{(k)} - l_i^{(k)}$), and possibly once by the sum of absolute output weights. i.e., if the next layer is a fully connected layer $\boldsymbol{x} \mapsto \boldsymbol{Ax} + \boldsymbol{b}$, the key of neuron i is $\sum_j |A_{i,j}|$. If the next layer is a ReLU layer, we skip the ReLU layer and use the weights from the fully connected layer that follows it (if any). The two ranks of a neuron in both orders are added, and the α-fraction with smallest rank sum is selected and their bounds are refined with a timeout of T whereas the next δ-fraction of neurons are refined with a timeout of $\beta \cdot \overline{T}$.

End-to-end approach. To certify robustness of deep neural networks, we combine MILP, LP relaxation, and abstract interpretation. We first pick numbers of layers $k_{\text{MILP}}, k_{\text{LP}}, k_{\text{AI}}$ that sum to the total number of layers of the neural network. For the analysis of the k_{MILP} layers, we refine bounds using anytime MILP relaxation with the neuron selection heuristic. As an optimization, we do not perform refinement after the abstract transformer for the first layer in case it is an affine transformation as the abstract domain computes the tightest possible bounding box for an affine transformation of a box (this is always the case in our experiments). For the k_{LP} layers, we refine bounds using LP relaxation combined with the neuron selection heuristic, i.e., the network up to the layer to be refined is encoded using linear constraints. For the remaining k_{AI} layers, we use abstract interpretation without additional refinement (however, this also benefits from refinement that was performed in previous layers), and compute the bounds using ι.

Final property certification. Let k be the index of the last layer and p be the number of output classes. We can encode the final certification problem using the output abstract element a_{out} obtained after applying the abstract transformer for the last layer in the network. If we want to prove that class i is assigned a higher probability than class j, it suffices to show that $a_{\text{out}} \sqcap (x_i^{(k)} \leq x_j^{(k)}) = \bot$. If this fails, one can resort to complete verification using MILP: the property is satisfied if and only if the set of constraints $\varphi^{(k)}(x_1^{(0)}, \ldots, x_p^{(k)}) \wedge (x_i^{(k)} \leq x_j^{(k)})$ is unsatisfiable.

4 EVALUATION

We now evaluate the effectiveness of our approach for the problem of robustness verification of ReLU-based feedforward and convolutional neural networks. Our results show that our approach enables faster complete verification than the state-of-the-art complete certifier from Tjeng et al. (2019) and produces more precise results than state-of-the-art incomplete certifiers: *DeepZ* (Singh et al., 2018) and *DeepPoly* (Singh et al., 2019), when complete certification becomes infeasible.

We implemented our approach in the form of a Python analyzer which we call *RefineAI*. *RefineAI* uses the Zonotope transformers from *DeepZ* for performing abstract interpretation over the networks and Gurobi (Gurobi Optimization, LLC, 2018) for solving MILP and LP instances. *RefineAI* builds on the Zonotope and Interval abstractions implemented in the ELINA library (eli, 2018; Singh et al., 2017). All of our code, neural networks, and images used in our experiments are publicly available at https://github.com/eth-sri/eran.

Evaluation datasets. We used the popular MNIST (Lecun et al., 1998) and CIFAR10 (Krizhevsky, 2009) datasets in our experiments. Both datasets include a training set of $60\,000$ and a testing set of $10\,000$ images. All images in the MNIST dataset are grayscale images of size 28×28 pixels whereas those in CIFAR10 are RGB images of size 32×32.

Table 1: Neural network architectures used in our experiments.

Dataset	Model	Type	#Neurons	#layers	Defense
MNIST	3×50	fully connected	160	3	None
	5×100	fully connected	510	5	DiffAI
	6×100	fully connected	610	6	None
	9×100	fully connected	910	9	None
	6×200	fully connected	1 210	6	None
	9×200	fully connected	1 810	9	None
	ConvSmall	convolutional	3 604	3	None
	ConvBig	convolutional	34 688	6	DiffAI
	ConvSuper	convolutional	88 500	6	DiffAI
CIFAR10	6×100	fully connected	610	6	None
	ConvSmall	convolutional	4 852	3	DiffAI

Neural networks. Table 1 shows 11 different MNIST and CIFAR10 feedforward (FNNs) and convolutional networks (CNNs) with ReLU activations used in our experiments. Out of these 4 were trained to be robust against adversarial attacks using DiffAI (Mirman et al., 2018) whereas the remaining 7 did not have any adversarial training. The largest network in our experiments contains > 88K neurons whereas the deepest network contains 9 layers.

Robustness properties. We consider the adversarial region defined by the L_∞-norm (Carlini & Wagner, 2017) perturbation. The region is parameterized by $\epsilon \in \mathbb{R}$ and contains all images in which each pixel $\overline{x}_i$ has distance $\leq \epsilon$ from the corresponding pixel x_i in the original input image x. Our goal is to certify that the neural network produces the same label as on the original image on all images in the adversarial region. Thus, the adversarial attack considered here is untargeted.

Experimental setup. All experiments for the 3×50 FNN and all CNNs were carried out on a 2.6 GHz Intel Xeon CPU E5-2690 with 512 GB of main memory; the remaining FNNs were evaluated on a 3.3 GHz 10 Core Intel i9-7900X Skylake CPU with a main memory of 64 GB.

Benchmarks. We selected the first 100 images from the test set of both MNIST and CIFAR10. Any image on which the neural network misclassified without perturbation was not considered for robustness certification in our experiments. We consider complete certification with *RefineAI* on the 3×50 network and chose an ϵ for which the incomplete certifier *DeepZ* certified $< 40\%$ of all candidate images. We consider incomplete certification for the remaining networks and chose an ϵ for which complete certification with *RefineAI* becomes infeasible.

4.1 COMPLETE CERTIFICATION

We first consider the exact robustness certification of the MNIST 3×50 networks against L_∞-norm attack with $\epsilon = 0.03$. *DeepZ* verifies 37% of images in the test set for this value of ϵ. *RefineAI* first runs *DeepZ* analysis on the whole network collecting the bounds for all neurons in the network. If *DeepZ* fails to certify, then the collected bounds are used to encode the robustness certification as a MILP instance (discussed in section 3). We compare *RefineAI* against the approach from Tjeng et al. (2019) which is a state-of-the-art complete certifier. This approach is also MILP-based like ours, but it uses Interval analysis and LP to determine neuron bounds. We implemented the Interval analysis and LP-based analysis to determine the initial bounds. We call the MILP solver only If LP analysis (or Interval analysis) fails to certify. All complete certifiers certify the neural network to be robust against L_∞-norm perturbations for 85% of the images. The average runtimes of *RefineAI*, MILP with presolve bounds from Interval analysis, and MILP with bounds from LP analysis are 28, 123, and 35 seconds respectively. Based on our result, we believe that the Zonotope analysis offers a good middle ground between the speed of the Interval analysis and the precision of LP for bound computation as it produces precise bounds faster than LP.

Table 2: Precision and runtime of *RefineAI* vs. *DeepZ* and *DeepPoly*.

Dataset	Model	ϵ	DeepZ		DeepPoly		RefineAI	
			precision(%)	time(s)	precision(%)	time(s)	precision(%)	time(s)
MNIST	5×100	0.07	38	0.6	53	0.3	53	381
	6×100	0.02	31	0.6	47	0.2	67	194
	9×100	0.02	28	1.0	44	0.3	59	246
	6×200	0.015	13	1.8	32	0.5	39	567
	9×200	0.015	12	3.7	30	0.9	38	826
	ConvSmall	0.12	7	1.4	13	6.0	21	748
	ConvBig	0.2	79	7	78	61	80	193
	ConvSuper	0.1	97	133	97	400	97	665
CIFAR10	6×100	0.0012	31	4.0	46	0.6	46	765
	ConvSmall	0.03	17	5.8	21	20	21	550

4.2 INCOMPLETE CERTIFICATION

We next consider incomplete certification for the remaining networks and compare against *DeepZ* and *DeepPoly*. We measure the average runtime per image in seconds and the precision measured by the % of images for which the certifier certified the network to be robust. We note that *DeepZ* has same precision as Fast-Lin (Weng et al., 2018) and *DeepPoly* has the same precision as CROWN (Zhang et al., 2018). The ϵ values used for defining the adversarial regions are shown in Table 2. The ϵ values for networks trained to be robust are larger than for networks that are not. We note that running the Interval analysis to obtain initial bounds loses too much precision on these large networks with ϵ values considered in our experiments. As a result, the approach from Tjeng et al. (2019) has to rely on applying LP per neuron to obtain precise bounds for the MILP solver which does not scale. For example, on the 9×200 network, determining bounds with LP already takes > 20 minutes (without calling the MILP solver which is more expensive than LP) whereas *RefineAI* has an average running time of ≈ 14 minutes.

Parameter values. We experimented with different values of the analysis parameters $k_{\text{MILP}}, k_{\text{LP}}, k_{\text{AI}}, \alpha, \delta, \beta, T$ and chose those values that offered the best tradeoff between performance and precision for the robustness certification for each neural network. We refine the output of all affine transformations in the network that are followed by a ReLU except the first one. In a given layer, we consider all neurons μ that can take non-negative values after applying the affine transformation as candidates for refinement.

For the MNIST 5×100, 6×100, 9×100, 6×200, and 9×200 FNNs, we refine the candidate neurons in the output of the affine transformations in layer 2-4 with MILP and those in the remaining layers using LP. For MILP based refinement, we use $\alpha = \frac{\mu}{5^{k-2} \cdot p}$ in layer k with p neurons. For LP based refinement, we use $\alpha = \frac{\mu}{2^{k-5} \cdot p}$. We use timeout $T = 1$ second, $\beta = 0.5$, and $\delta = \frac{\mu}{p} - \alpha$ for both MILP and LP based refinements. For the CIFAR10 6×100 FNN, we use the same values except that we use $\alpha = \frac{\mu}{2^{k-2} \cdot p}$ for MILP refinement and set $T = 6$ seconds for both MILP and LP based refinement as it is more time consuming to refine for CIFAR10 networks.

For the CNNs, we do not refine convolutional layers as the number of candidates in these layers are very large and instead refine all candidate neurons in fully connected layers with larger timeout to compensate for more difficult problem instances for the solver. For the MNIST ConvSmall, ConvBig and CIFAR10 ConvSmall networks, we refine all the candidate neurons with MILP using a timeout $T = 10$ seconds. For the MNIST ConvSuper network, we refine similarly but use LP instead of MILP with a timeout of $T = 15$ seconds.

Results for incomplete certification. Table 2 shows the precision and the average runtime of all three certifiers. It can be seen that *RefineAI* is significantly more precise than both *DeepZ* and *DeepPoly*. It certifies significantly more images than *DeepZ* on all networks except the MNIST ConvSuper network. This is because that for the ϵ considered, *DeepZ* is already very precise. *Re-*

fineAI certifies network to be more robust on more images than *DeepPoly* on 6 out of 10 networks. It can be seen that the number of neurons in the network is not the determining factor for the average runtime of *RefineAI*. We observe that *RefineAI* runs faster on the networks trained to be robust and the top three networks with the largest runtime for *RefineAI* are all networks not trained to be robust. This is because robust networks are relatively easier to certify and produce only a small number of candidate neurons for refinement, which are easier to refine by the solver. For example, even though the same parameter values are used for refining the results on the MNIST ConvSmall and ConvBig networks, the average runtime of *RefineAI* on the robustly trained ConvBig network with around 35K neurons, 6 layers and a perturbation region defined using $\epsilon = 0.2$ is almost 4 times less than on the non-robust ConvSmall network with only $3,604$ neurons, 3 layers and a smaller $\epsilon = 0.12$.

4.3 EFFECT OF NEURON SELECTION HEURISTIC

We use the neuron selection heuristic from section 3 to determine neurons which need to be refined more than others for FNNs as refining all neurons in a layer with MILP can significantly slow down the analysis. To check whether our heuristic is able to identify important neurons, we ran the analysis on the MNIST 9×200 FNN by keeping all analysis parameters the same, except instead of selecting the neurons with the smallest rank sum first we selected the neurons with the largest rank sum first (thus refining neurons more if our heuristic deeps them unimportant). With this change, the average runtime does not change significantly. However, the modified analysis loses precision and fails to certify two images that the analysis refining with our neuron selection heuristic certifies.

5 CONCLUSION

In this paper, we presented a refinement-based approach for increasing the precision of incomplete verifiers for feedforward and convolutional neural networks with ReLU activations. This enabled us to certify robustness properties that existing state-of-the-art complete verifiers cannot handle (due to scalability issues) whereas state-of-the-art incomplete verifiers could not prove as they use overapproximations that are too coarse. We believe combining the strengths of overapproximation methods with that of mixed integer linear programming as done in this work is a promising direction for further advancing the state-of-the-art in neural network certification.

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