
Nice2Predict and JSNice

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Predicting Program Properties from “Big Code”

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Background

Deobfuscating JavaScript

Variable renaming

Minimize file size

Obscure program intent

Syntactic

Type annotation

Types specified in JSDoc comments

Compare Closure compiler, Flow, TypeScript, etc.

Semantic

Obfuscated

Compilation was a success!

Original Size: 247 bytes gzipped (445 bytes uncompressed)

Compiled Size: 122 bytes gzipped (128 bytes uncompressed)

Saved 50.61% off the gzipped size (71.24% without gzip)

The code may also be accessed at [default.js](#).

Compiled Code

Warnings

Errors

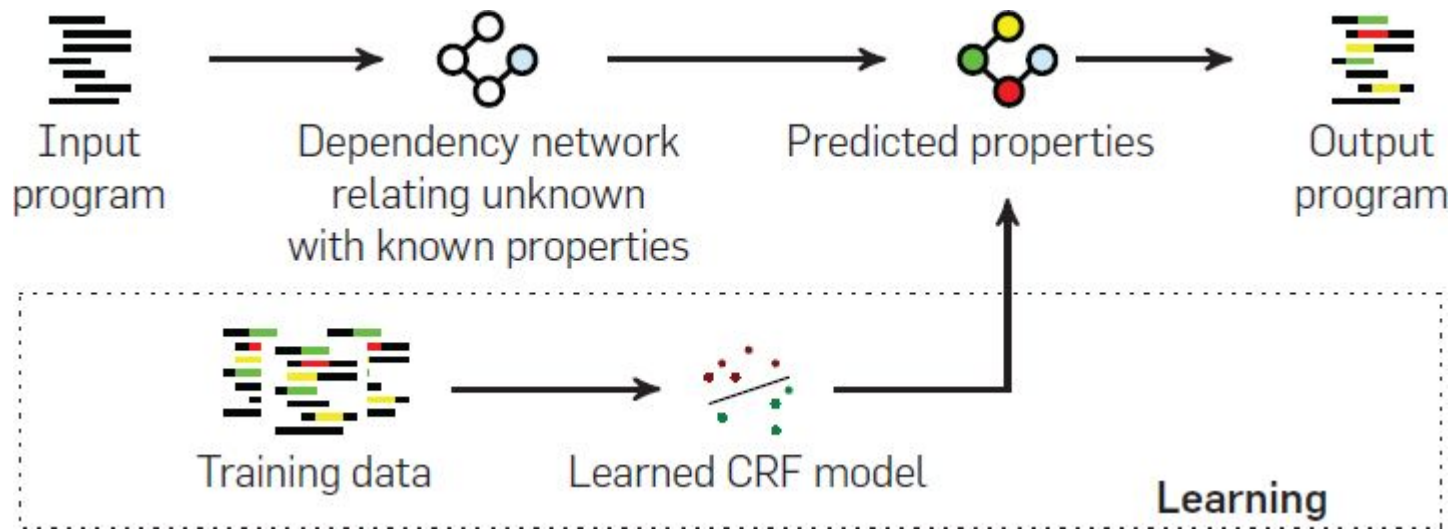
POST data

```
window.a=function(c,d){for(var e=[],f=c.length,b=0;b<f;b+=d)b+d<f?e.push(c.substring(b,b+d)):e.push(c.substring(b,f));return e};
```

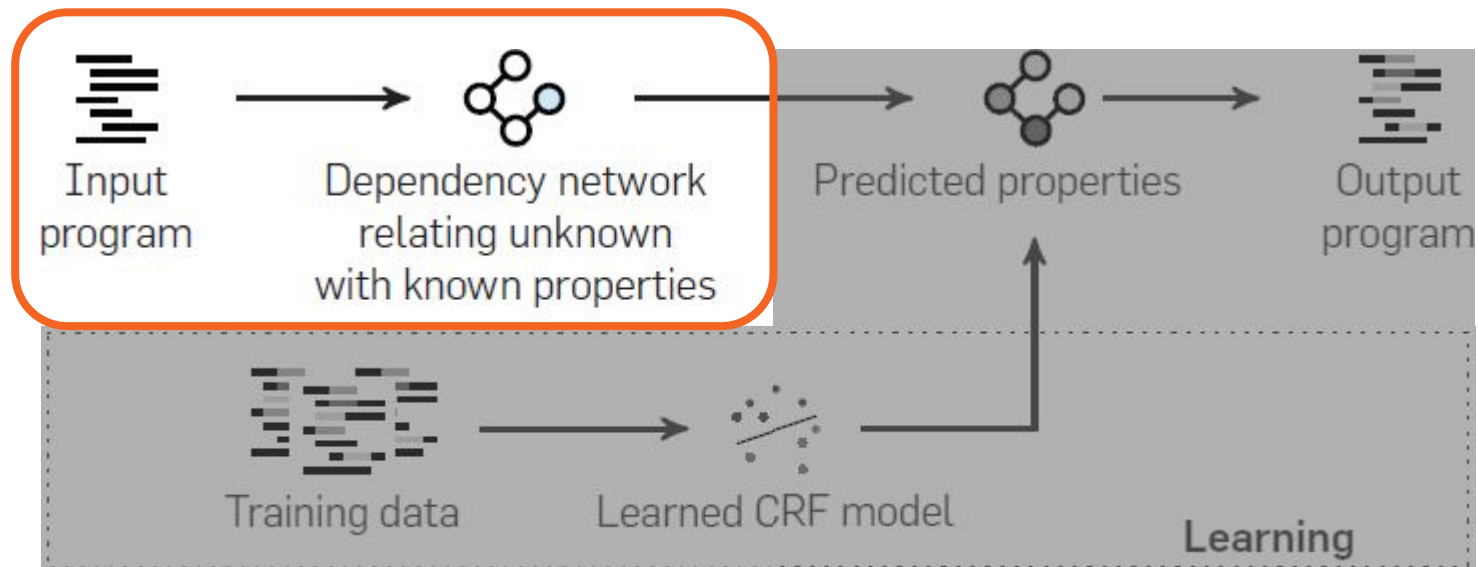
Original: commented, annotated

```
'use strict';  
/**  
 * @param {string} bin  
 * @param {number} size  
 * @return {?} */  
*/  
window.chunkData = function chunkData(bin, size) {  
  /** @type {!Array} */  
  var results = [];  
  var length = bin.length;  
  /** @type {number} */  
  var i = 0;  
  for (; i < length; i = i + size) {  
    if (i + size < length) {  
      results.push(bin.substring(i, i + size));  
    } else {  
      results.push(bin.substring(i, length));  
    }  
  }  
  return results;  
}
```

Approach



Structured Prediction for Programs



Dependency network

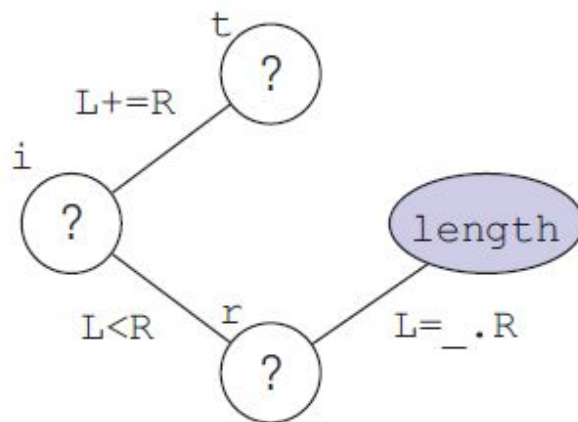
Input: Obfuscated Program

```
function chunkData(e, t) {  
    var n = [];  
    var r = e.length;  
    var i = 0;  
    for (; i < r; i += t) {  
        if (i + t < r) {  
            n.push(e.substring(i, i + t));  
        } else {  
            n.push(e.substring(i, r));  
        }  
    }  
    return n;  
}
```

Unknown properties (variable names):



Known properties (constants, APIs):



Extract properties and build dependency network

—

Extract properties

Extracting Names

Known

Constants

Object Properties

Methods and API Calls

Global Variables

Mostly treated as string constants

Unknown

Local variables

Different scopes -> different properties

Keywords and naming conflicts not allowed in the prediction space Ω_x

Extracting Types

Known

Any **expression** with known type or any constant

Manually provided or built with program analysis

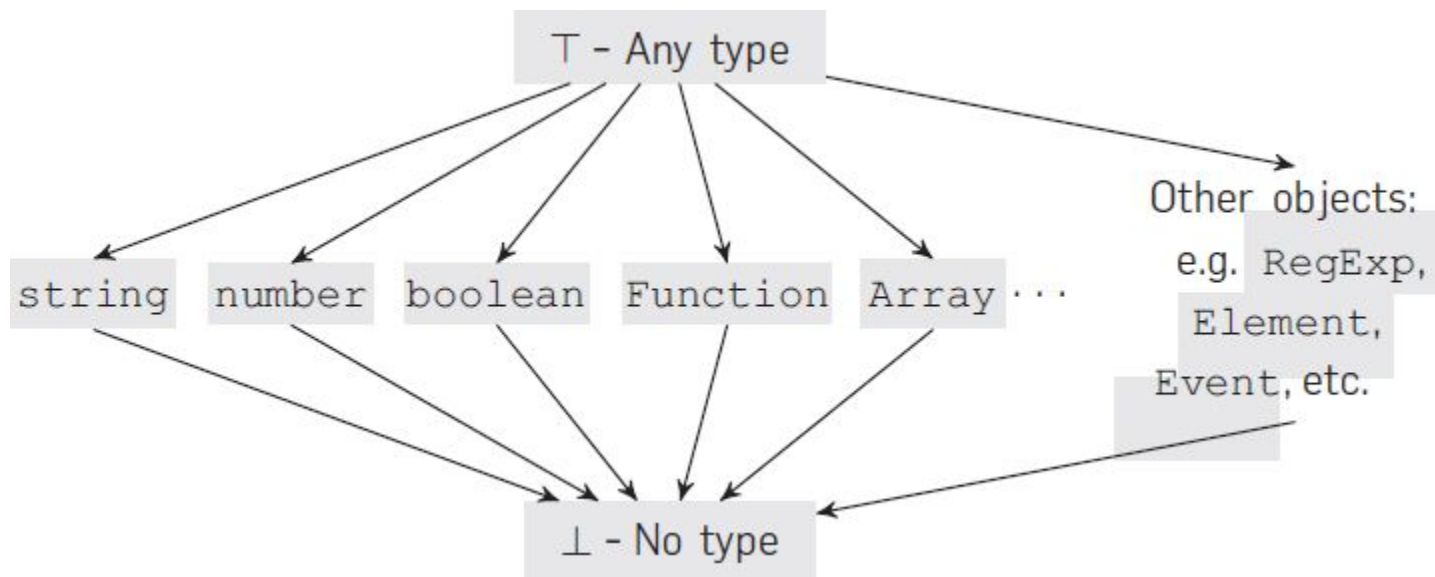
Unknown

Variables with unknown type

$\Omega_x = (JTypes)^n$: no constraints on predictions.

$expr ::= val \mid var \mid expr_1(expr_2) \mid expr_1 \otimes expr_2$ *Expression*
 $val ::= \lambda var : \tau.expr \mid n$ *Value*

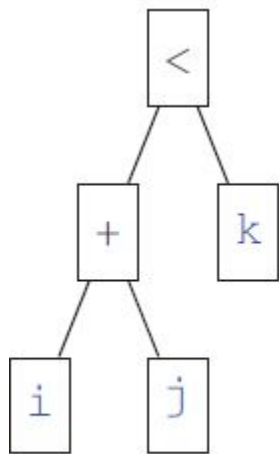
Type Lattice



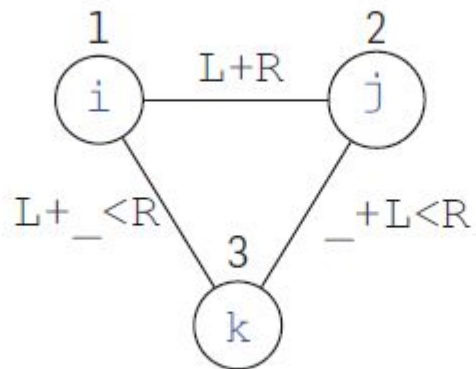
JSTypes is the powerset of all types over this lattice

Build dependency graph

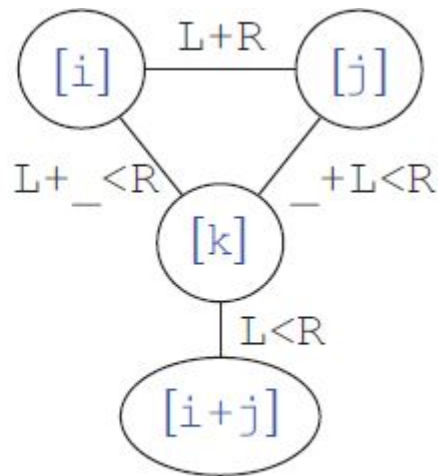
$i + j < k$



AST



Names



Types

Grammar

$$rel_{ast} ::= rel_L(rel_R) \mid rel_L \otimes rel_R$$
$$rel_L ::= \mathbf{L} \mid rel_L(_) \mid _(rel_L) \mid rel_L \otimes _ \mid _ \otimes rel_L$$
$$rel_R ::= \mathbf{R} \mid rel_R(_) \mid _(rel_R) \mid rel_R \otimes _ \mid _ \otimes rel_R$$

Additional relations

ARG_TO_PM

Relates arguments of a function invocation to parameters in the function declaration.

ALIAS

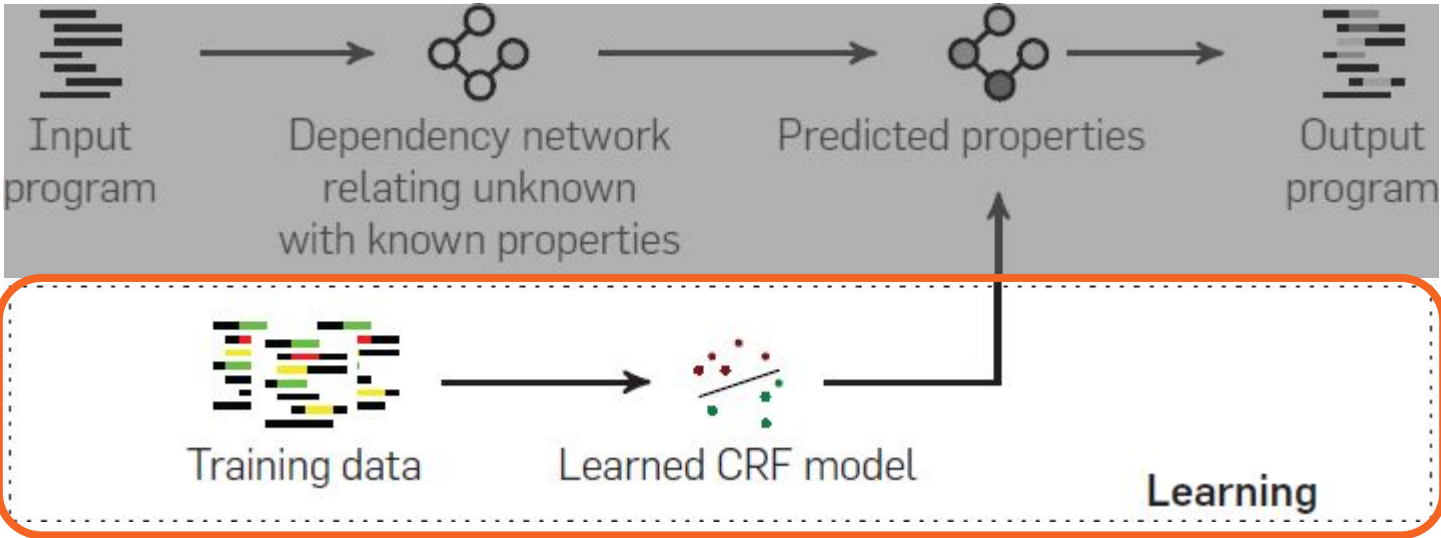
Types only. If a and b are related with r , and c is a variable that aliases b , we add the edge $(a, c, (r, ALIAS))$

MAY_CALL

Names only. If a function variable f may call function g , we add the edge (f, g, MAY_CALL)

MAY_ACCESS

Names only. If in a function variable f , there is an access to an object field named foo , we add the edge (f, foo, MAY_ACCESS)



Learning

Discriminative not Generative

Since predictions are all made *given* a specific observed program x , we are only concerned with the conditional probability $Pr(\mathbf{y} | x)$, not the joint probability $Pr(\mathbf{y}, x)$.

This means we don't need to make any assumptions about the prior probabilities of the observed properties.

—

CRF

CRF

Conditional Random Field

A model for the conditional probability of labels \mathbf{y} given observations x

$$Pr(\mathbf{y} \mid x) = \frac{1}{Z(x)} \exp(\text{score}(\mathbf{y}, x))$$

$Z(x)$ is just a normalization factor:

$$Z(x) = \sum_{\mathbf{y}' \in \Omega_x} \exp(\text{score}(\mathbf{y}', x))$$

score

Sum of feature functions \mathbf{f} associated
with weights \mathbf{w}

$$\text{score}(\mathbf{y}, x) = \sum_{i=1}^k w_i f_i(\mathbf{y}, x) = \mathbf{w}^T \mathbf{f}(\mathbf{y}, x)$$

Feature Functions

Can be anything that controls the likelihood of a given assignment \mathbf{y} .

In practice, this paper just uses the sum of pairwise indicator feature functions over the edges of the dependence network

$$f_i(\mathbf{y}, x) = \sum_{(a,b,rel) \in E^x} \psi_i((\mathbf{y}, \mathbf{z})_a, (\mathbf{y}, \mathbf{z})_b, rel)$$

\mathbf{z} = assignments for known properties

Pairwise Indicator Feature Functions

Independent of the program being queried; defined once for all predictions of a given type (variable names or types).

Preprocessed from all features in the training set (plus an extra feature for equality).

Predicted outputs will be chosen from this same set of possible features.

$\forall \langle l_i^1, l_i^2, rel_i \rangle \in all_features(D):$

$$\psi_i(l^1, l^2, rel) = \begin{cases} 1 & \text{if } l^1 = l_i^1 \text{ and } l^2 = l_i^2 \text{ and } rel = rel_i \\ 0 & \text{otherwise} \end{cases}$$

score

Substituting allows us to simplify the score function. Since $\mathbf{Z}(x)$ does not depend on \mathbf{y}' , this has the same maximum as the full CRF probability.

$$\text{score}(\mathbf{y}, x) = \sum_{(a,b,rel) \in E^x} \sum_{i=1}^k w_i \psi_i((\mathbf{y}, \mathbf{z})_a, (\mathbf{y}, \mathbf{z})_b, rel)$$

—

Learning w

Structured Support Vector Machine (SSVM)*

Generalization of classical SVMs to predict many interdependent labels at once.

Maximize Δ , the margin between $\mathbf{y}^{(j)}$ and every other \mathbf{y}'

$$\forall j, \forall \mathbf{y}' \in \Omega_{x^{(j)}} \quad \text{score}(\mathbf{y}^{(j)}, x^{(j)}) \geq \text{score}(\mathbf{y}', x^{(j)}) + \Delta(\mathbf{y}^{(j)}, \mathbf{y}')$$

[*] TSOCHANTARIDIS, I., JOACHIMS, T., HOFMANN, T., AND ALTUN, Y. Large margin methods for structured and interdependent output variables. *Journal of Machine Learning Research* 6, 2005, 1453–1484

Structured Hinge Loss

$$\mathbf{w}^* = \underset{\mathbf{w}}{\operatorname{argmin}} \sum_{j=1}^t \ell(\mathbf{w}; x^{(j)}, \mathbf{y}^{(j)}) \text{ s.t. } \mathbf{w} \in \mathcal{W}_\lambda \quad (1)$$

where

$$\ell(\mathbf{w}; x^{(j)}, \mathbf{y}^{(j)}) = \max_{\mathbf{y}' \in \Omega_{x^{(j)}}} \mathbf{w}^T [\mathbf{f}(\mathbf{y}', x^{(j)}) - \mathbf{f}(\mathbf{y}^{(j)}, x^{(j)})] + \Delta(\mathbf{y}^{(j)}, \mathbf{y}')$$

Stochastic Gradient Descent*

On every iteration:

- Pick a random program from the training set
- Compute the gradient of the loss function for this program
- Take a step in the negative gradient direction, with step size determined by the learning rate α
- Project back to the feasible region

[*] RATLIFF, N. D., BAGNELL, J. A., AND ZINKEVICH, M. (approximate) subgradient methods for structured prediction. In AISTATS (2007), pp. 380–387.

Computing the gradient

$$\mathbf{y}_{best} \leftarrow \operatorname{argmax}_{\mathbf{y}' \in \Omega_{x^{(j)}}} (\operatorname{score}(\mathbf{y}', x^{(j)}) + \Delta(\mathbf{y}^{(j)}, \mathbf{y}')), \quad (2)$$

resulting in the gradient \mathbf{g}

$$\mathbf{g} \leftarrow \mathbf{f}(\mathbf{y}_{best}, x^{(j)}) - \mathbf{f}(\mathbf{y}^{(j)}, x^{(j)})$$

Regularization

Enforces non-negativity

Minimizes overfitting

Operates on vector components independently

$\lambda = 2.0$ for names, 5.0 for types

$$\text{Proj}_{\mathcal{W}_\lambda}(\mathbf{w}) = \mathbf{w}' \quad \text{such that} \quad w'_i = \max(0, \min(1/\lambda, w_i))$$

Training

Initialize: $w_i = 1/2\lambda$, $\alpha = 0.1$

If the number of wrong labels does not decrease, halve α

Up to 24 iterations over the data

Parallelized across multiple threads*

[*] ZINKEVICH, M., WEIMER, M., LI, L., AND SMOLA, A. J. Parallelized stochastic gradient descent. In NIPS (2010), pp. 2595–2603.

Training runtime

32-core machine with four 2.13GHz Xeon processors, running Ubuntu 12.04 with 64-Bit OpenJDK Java 1.7.0_51.

- Training for name prediction: ~10 hours.
 - 57 minutes to compile the input code and generate networks for the input programs
 - 23 minutes per SSVM (sub-) gradient descent optimization pass
 - Training for type prediction
 - 57 minutes for compilation and network construction
 - 2 minutes and 16 seconds per SSVM (sub-)gradient descent optimization pass
-

Model Sizes

Names

7,627,484 features

145.5MB

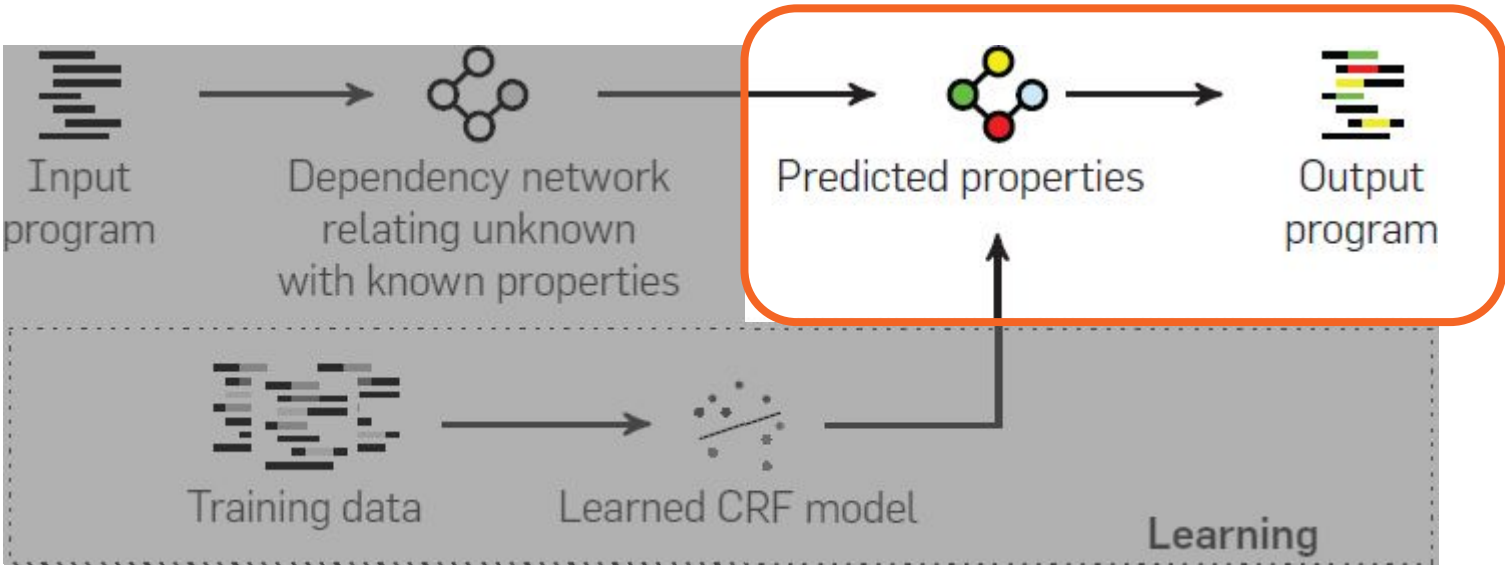
Dictionary of all names and types

16.8MB

Types

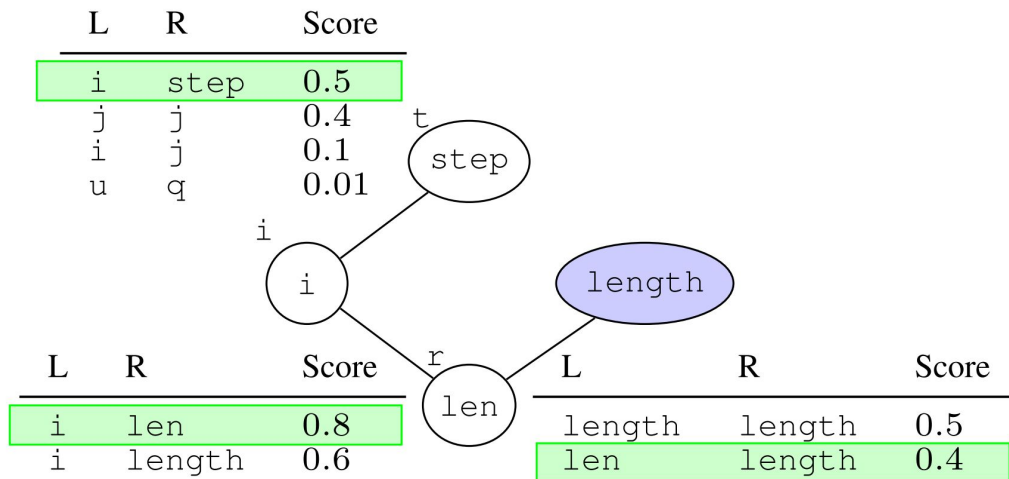
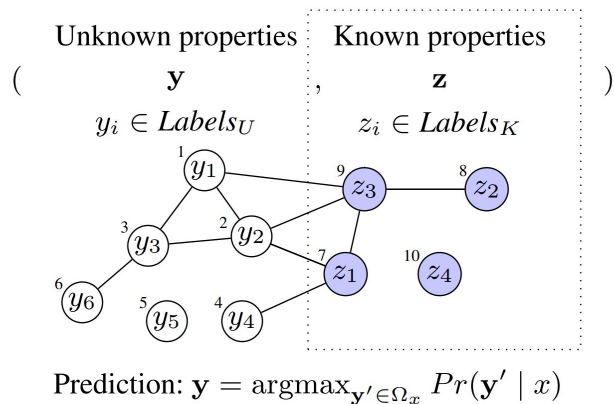
70,052 features

1.3MB



MAP inference

Prediction



$$\mathbf{y} = \operatorname{argmax}_{\mathbf{y}' \in \Omega_x} Pr(\mathbf{y}' | x) = \operatorname{argmax}_{\mathbf{y}' \in \Omega_x} score(\mathbf{y}', x) = \operatorname{argmax}_{\mathbf{y}' \in \Omega_x} \mathbf{w}^T \mathbf{f}(\mathbf{y}', x)$$

Algorithm 1: Greedy Inference Algorithm

Input: network $G^x = \langle V^x, E^x \rangle$ of program x ,
initial assignment of n unknown properties $\mathbf{y}_0 \in \Omega_x$,
known properties \mathbf{z}
pairwise feature functions ψ_i and their learned weights w_i
Output: $\mathbf{y} \approx \operatorname{argmax}_{\mathbf{y}' \in \Omega_x} (\operatorname{score}(\mathbf{y}', x))$

```
1 begin
2    $\mathbf{y} \leftarrow \mathbf{y}_0$ 
3   for  $pass \in [1..num\_passes]$  do
4     // for each node with unknown property in the graph  $G^x$ 
5     for  $v \in [1..n]$  do
6        $E_v \leftarrow \{(v, -, -) \in E^x\} \cup \{(-, v, -) \in E^x\}$ 
7        $score_v \leftarrow \operatorname{scoreEdges}(E_v, (\mathbf{y}, \mathbf{z}))$ 
8       for  $l' \in \operatorname{candidates}(v, (\mathbf{y}, \mathbf{z}), E_v)$  do
9          $l \leftarrow y_v$  // get current label of  $v$ 
10         $y_v \leftarrow l'$  // change label of  $v$  in  $\mathbf{y}$ 
11         $score'_v \leftarrow \operatorname{scoreEdges}(E_v, (\mathbf{y}, \mathbf{z}))$ 
12        if  $\mathbf{y} \in \Omega_x \wedge score'_v > score_v$  then
13          |  $score_v \leftarrow score'_v$ 
14        else
15          |  $y_v \leftarrow l$  // no score improvement: revert label.
16  return  $\mathbf{y}$ 
```

scoreEdges()

Score the subset of the network adjacent to the current node

$$\text{scoreEdges}(E, A) = \sum_{(a,b,rel) \in E} \sum_{i=1}^k w_i \psi_i(A_a, A_b, rel)$$

candidates()

Takes the s labels with the highest corresponding weights for each edge

Beam size s controls precision vs. running time

$s = 64$, experimentally determined (see Results section)

Decrease the beam size by a factor of 16 if a node has more than 32 adjacent nodes

Optimize by edges (pairs) instead of single nodes

“At almost no computation cost, we also perform optimizations on pairs of nodes in addition to individual nodes. In this case, for each edge in G^x , we use the s best scoring features on the same type of edge in the training set and attempt to set the labels of the two elements connected by the edge to the values in each triple.”

Not clear how this fits with the algorithm pseudocode

Results

Datasets

Training

10,517 JavaScript projects from
GitHub

No overlap with eval set

324,501 files

Filtered minified and obfuscated files

Evaluation

50 JavaScript projects with the
highest number of commits from
BitBucket

2,710 files

381,243 LOC, Largest file 3,055

383.5 (109.5) arcs and 29.2 (12.6)
random variables for names (types)
on average in the eval set

Parameter selection

10-fold cross-validation

1% sample of the training data

$\lambda = 2.0$ for names, 5.0 for types

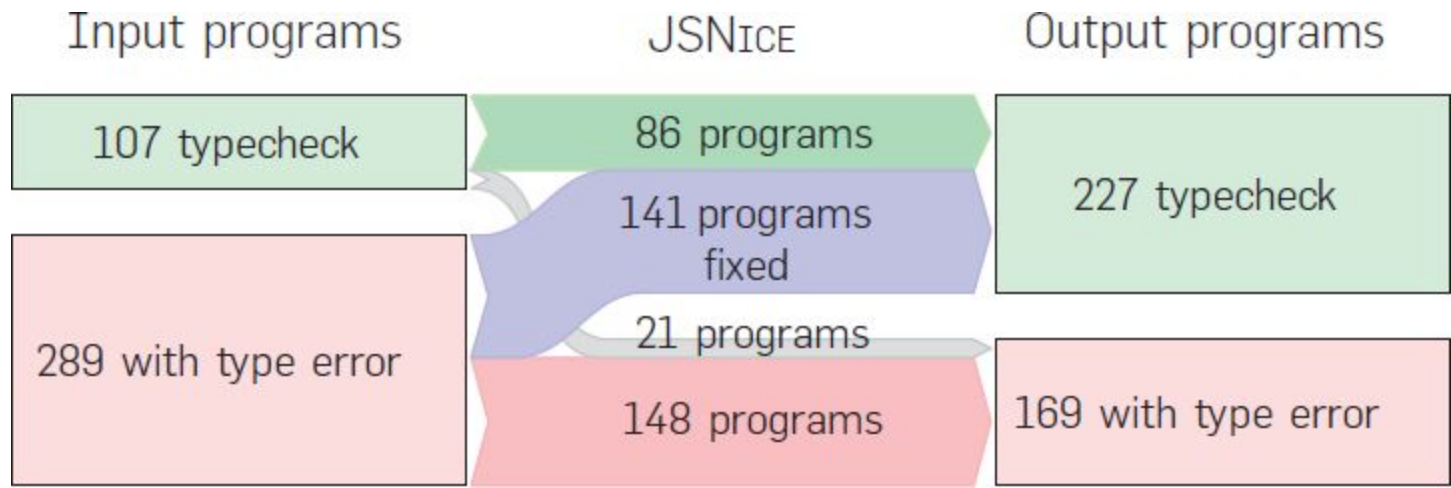
Margin Δ should be applied

Results

| System | Names Accuracy | Types Precision | Types Recall |
|---------------------------|-------------------|--------------------|-----------------|
| all training data | 63.4% | 81.6% | 66.9% |
| 10% of training data | 54.5% | 81.4% | 64.8% |
| 1% of training data | 41.2% | 77.9% | 62.8% |
| all data, no structure | 54.1% | 84.0% | 56.0% |
| baseline - no predictions | 25.3% | 37.8% | 100% |

Beam size and prediction time

| Beam parameter | Name prediction | | Type prediction | |
|-----------------------|-----------------|---------|-----------------|-------|
| | Accuracy | Time | Precision | Time |
| b | | | | |
| 4 | 57.9% | 43ms | 80.6% | 36ms |
| 8 | 59.2% | 60ms | 80.9% | 39ms |
| 16 | 62.8% | 62ms | 81.6% | 33ms |
| 32 | 63.2% | 80ms | 81.3% | 37ms |
| 64 (JSNICE) | 63.4% | 114ms | 81.6% | 40ms |
| 128 | 63.5% | 175ms | 82.0% | 42ms |
| 256 | 63.5% | 275ms | 81.6% | 50ms |
| Naïve greedy, no beam | 62.8% | 115.2 s | 81.7% | 410ms |



Typechecking

Demo!?

Discussion

A note on name inference

“We note that our name inference process is independent of what the minified names are. In particular, the process will return the same names regardless of which minifier was used to obfuscate the original program (provided these minifiers always rename the *same set* of variables).”

However...

```
1 function chunkData(e, t) {
2   var n = [];
3   var a = e.length;
4   var i = 0;
5   for (; i < a; i += t) {
6     if (i + t < a) {
7       n.push(e.substring(i, i + t));
8     } else {
9       n.push(e.substring(i, a));
10    }
11  }
12  return n;
13 }
14 }
```

```
1 function chunkData(▼xs, ▼i) {
2   var ▼data = [];
3   var ▼end = ▼xs.length;
4   var ▼offset = 0;
5   for (; ▼offset < ▼end; ▼offset += ▼i) {
6     if (▼offset + ▼i < ▼end) {
7       ▼data.push(▼xs.substring(▼offset, ▼offset +
8 ▼i));
9     } else {
10      ▼data.push(▼xs.substring(▼offset, ▼end));
11    }
12  }
13  return ▼data;
14 }
```

```
1 function chunkData(a1, t) {
2   var n = [];
3   var a = a1.length;
4   var i = 0;
5   for (; i < a; i += t) {
6     if (i + t < a) {
7       n.push(a1.substring(i, i + t));
8     } else {
9       n.push(a1.substring(i, a));
10    }
11  }
12  return n;
13 }
14 }
```

```
1 function chunkData(▼str, ▼step) {
2   var ▼buffer = [];
3   var ▼length = ▼str.length;
4   var ▼i = 0;
5   for (; ▼i < ▼length; ▼i += ▼step) {
6     if (▼i + ▼step < ▼length) {
7       ▼buffer.push(▼str.substring(▼i, ▼i + ▼step));
8     } else {
9       ▼buffer.push(▼str.substring(▼i, ▼length));
10    }
11  }
12  return ▼buffer;
13 }
```

Other concerns

Function names treated as givens, might be over-weighted

Different behavior between Proprietary <http://jsnice.org/> and
Open Source <http://www.nice2predict.org/>

Future Work

<http://apk-deguard.com/>

“Similarly to JSNice, DeGuard is based on powerful probabilistic graphical models learned from thousands of open source programs. Using these models, DeGuard recovers important information in Android APKs, including method and class names as well as third-party libraries. DeGuard can reveal string decoders and classes that handle sensitive data in Android malware.”

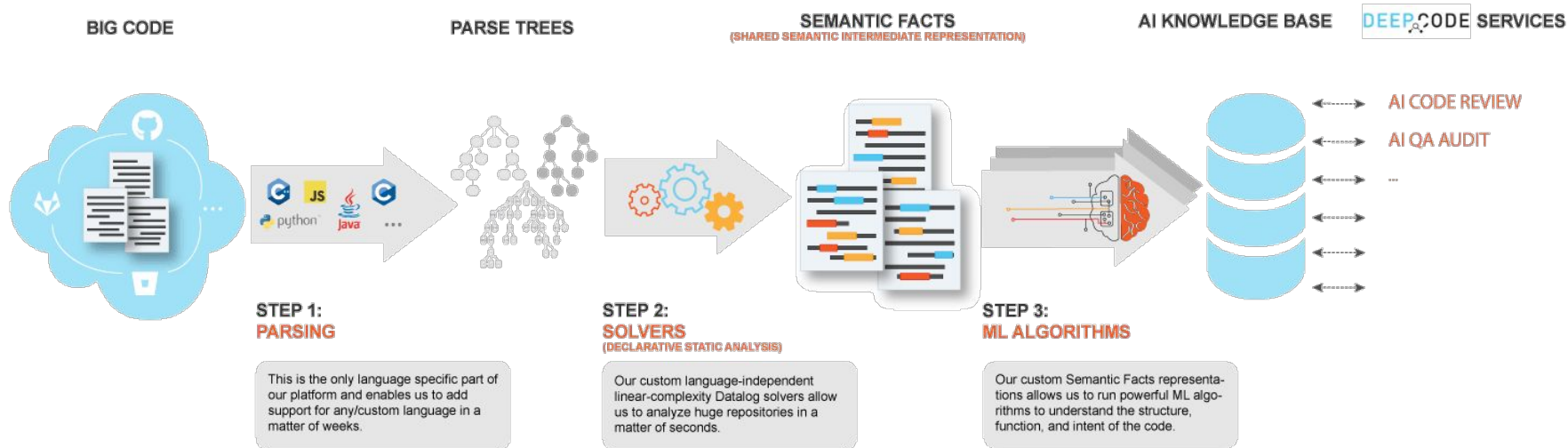
<https://www.sri.inf.ethz.ch/deguard>

<https://debin.ai/>

“DEBIN is a novel system for predicting debug information in stripped binaries. It is able to distinguish register-allocated and memory-allocated variables with decision-tree-based classification. Moreover, it is capable of predicting meaningful names and types for variables and functions through structured prediction with probabilistic graphical models. These models are learned from thousands of non-stripped binary in open source packages. The system can be further used for malware inspection.”

<https://github.com/eth-sri/debin>

<https://www.deepcode.ai/>



Beyond layout obfuscation

Can you start with a control flow graph instead of an AST?

Analyze code changes*

[*] Rumen Paletov, Petar Tsankov, Veselin Raychev, and Martin Vechev. 2018. Inferring crypto API rules from code changes. In Proceedings of the 39th ACM SIGPLAN Conference on Programming Language Design and Implementation (PLDI 2018). ACM, New York, NY, USA, 450-464. DOI: <https://doi.org/10.1145/3192366.3192403>

Questions?

The End



Thank you very much!