

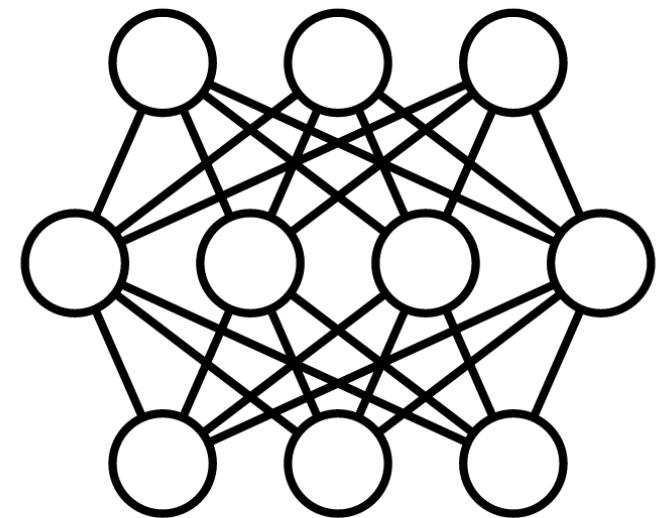
Learning to Optimize Tensor Programs

**Tianqi Chen, Lianmin Zheng, Eddie Yan, Ziheng Jiang, Thierry
Moreau, Luis Ceze, Carlos Guestrin, Arvind Krishnamurthy**

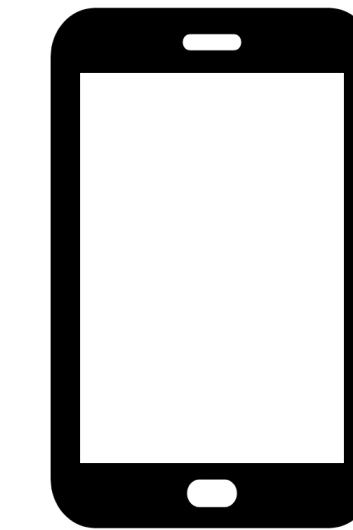
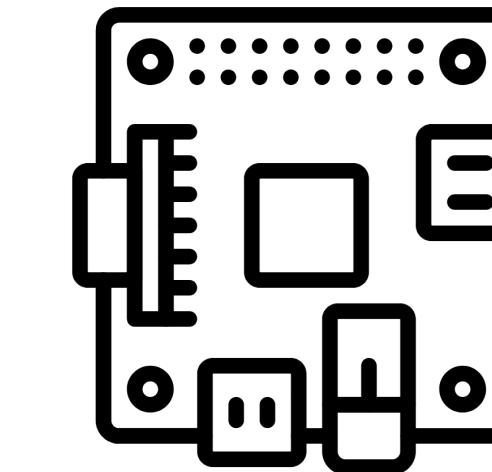
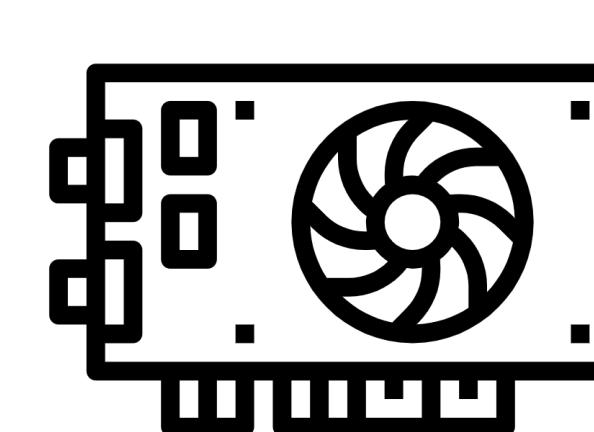
Overview

We want to **efficiently** deploy models to a variety of platforms

Model



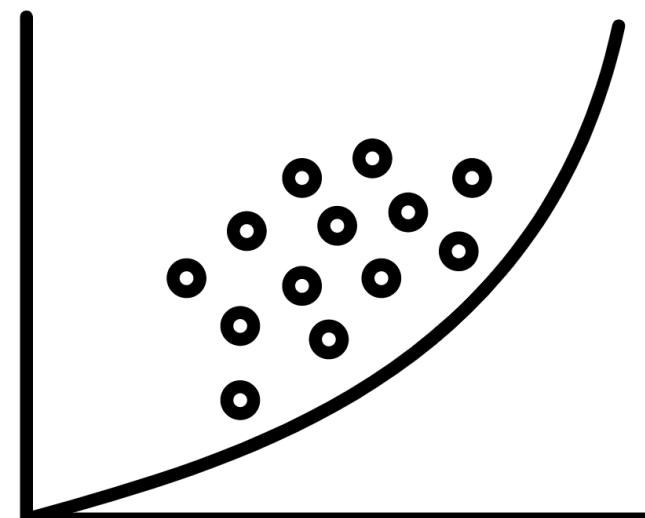
Target



Overview

Recall: TVM allows us to separate the **algorithm** (what is computed) of a program from its **schedule** (how it is computed)

AutoTVM finds the best schedule automatically by using



Cost-model based search
Transfer learning

The problem

$$C_{ij} = \sum_k A_{ki}B_{kj}$$

Possibly millions of programs!

```
for y in range(128):
    for x in range(128):
        C[y][x] = 0
        for k in range(128):
            C[y][x] += A[k][y] * B[k][x]
```

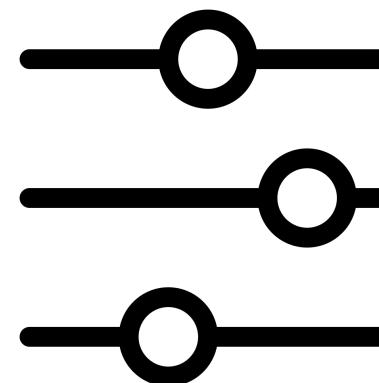
```
for yo in range(1024 / ty):
    for xo in range(1024 / tx):
        C[yo*ty:yo*ty+ty][xo*tx:xo*tx+tx] = 0
        for k in range(1024):
            for yi in range(ty):
                for xi in range(tx):
                    C[yo*ty+yi][xo*tx+xi] +=
                        A[k][yo*ty+yi] * B[k][xo*tx+xi]
```

```
for yo in range(128):
    for xo in range(128):
        intrin.fill_zero(C[yo*8:yo*8+8][xo*8:xo*8+8])
        for ko in range(128):
            intrin.fused_gemm8x8_add(
                C[yo*8:yo*8+8][xo*8:xo*8+8],
                A[yo*8:yo*8+8][xo*8:xo*8+8],
                B[yo*8:yo*8+8][xo*8:xo*8+8])
```

■ ■ ■

How do we efficiently explore the space of possible programs?

Hyperparameter optimization



Find the best set of hyperparameters by optimizing
a **costly to evaluate** fitness function f

Sequential Model-based Global Optimization

1. Choose x^* according to a surrogate model

2. Evaluate $f(x^*)$

3. Add $(x, f(x^*))$ to training data

4. Fit a new surrogate model

5. Repeat

E.g. Gaussian process or
tree-structured Parzen estimator

Assume f is a black-box
and expensive to evaluate

AutoTVM

Expression

$$C_{ij} = \sum_k A_{ki} B_{kj}$$

Schedule space

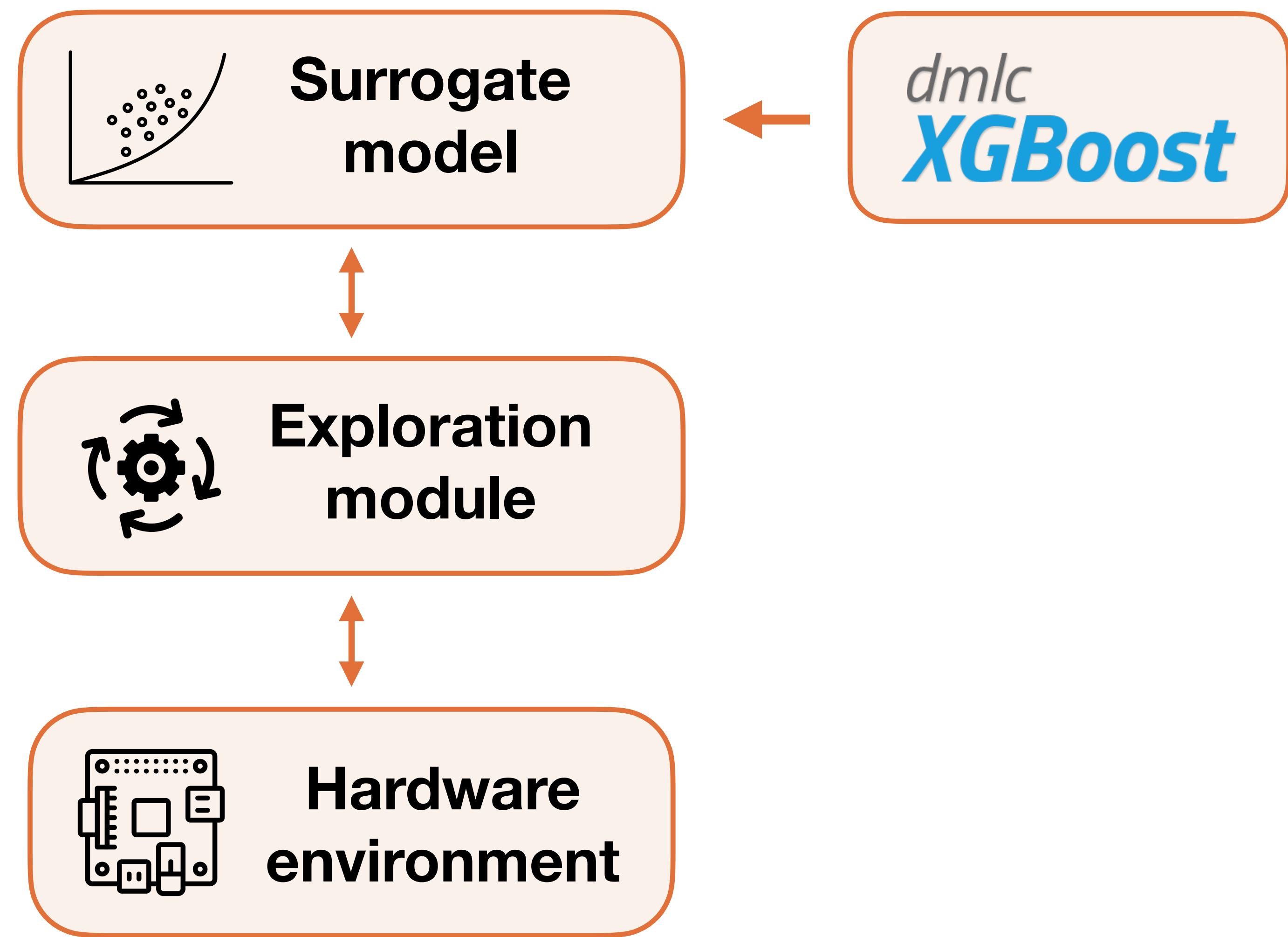
s_1 loop tiling

```
yo, xo, yi, xi = s[C].tile(y, x, ty, tx)
s[C].reorder(yo, xo, k, yi, xi)
```

$x_1 = g(e, s_1)$

```
for yo in range(1024 / ty):
    for xo in range(1024 / tx):
        C[yo*ty:yo*ty+ty][xo*tx:xo*tx+tx] = 0
        for k in range(8):
            for yi in range(8):
                for xi in range(8):
                    yo, xo, ko, yi, xi, ki = s[C].tile(y, x, k, 8, 8)
                    s[C].tensorize(yi, intrin.gemm8x8)
                    A[k][yo*8:yo*8+8][xo*8:xo*8+8] = B[ko*8:ko*8+8][yo*8:yo*8+8]
                    x2 = g(e, s2)
```

```
for yo in range(128):
    for xo in range(128):
        intrin.fill_zero(C[yo*8:yo*8+8][xo*8:xo*8+8])
        for ko in range(128):
            intrin.fused_gemm8x8_add(
                C[yo*8:yo*8+8][xo*8:xo*8+8],
                A[ko*8:ko*8+8][yo*8:yo*8+8],
                B[ko*8:ko*8+8][xo*8:xo*8+8])
```



Feature representation

Extract features from **AST** for every loop variable:

Loop length, vectorizable, parallelizable, ...

For every **buffer**: touch count, reuse ratio, ...

```
for y in range(8):
    for x in range(8):
        C[y][x]=0
        for k in range(8):
            C[y][x]+=A[k][y]*B[k][x]
```

	C	A	B	touched memory	outer loop length
y	64	64	64		y 1
x	8	8	64		x 8
k	1	8	8		k 64

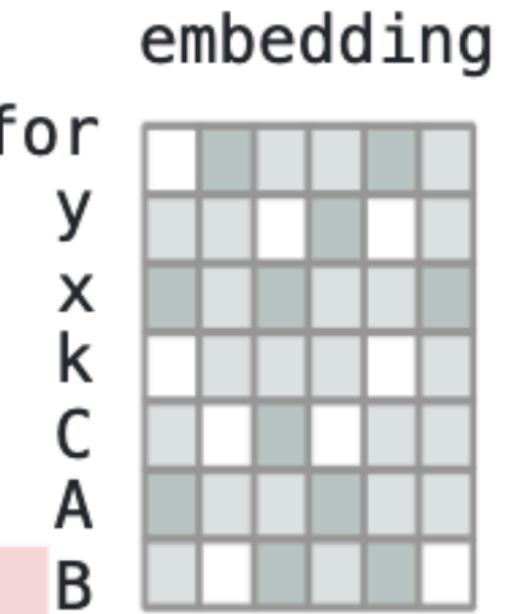
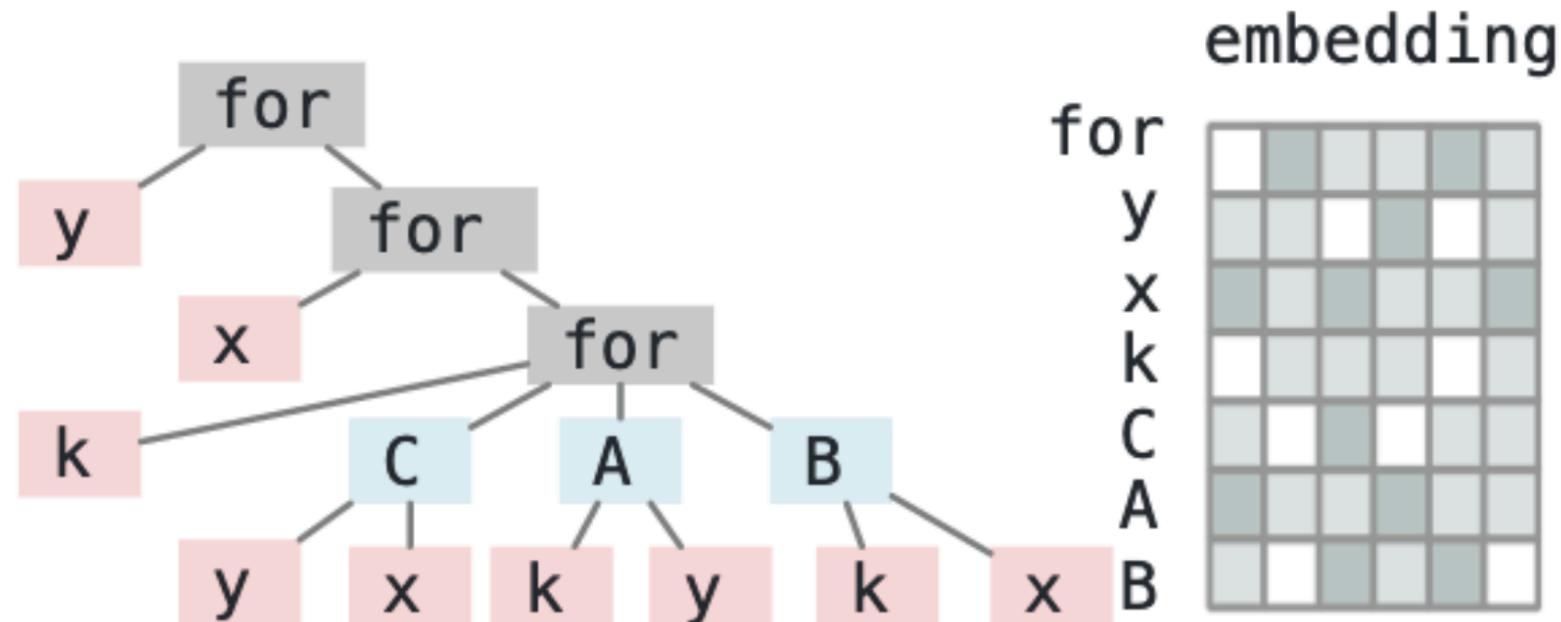
(a) Low level AST

(b) Loop context vectors

Feature representation

Alternative representation using TreeGRU

Recursively encode AST into an embedding vector



(c) Vanilla TreeGRU

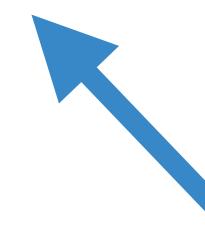
Choosing candidates

Encourage **diversity** by maximizing

$$L(S) = - \sum_{s \in S} \text{surrogate_model}(s) + \alpha \times (\# \text{ configuration components covered by } S)$$



Balance
quality



Balance
diversity

Encourage **exploration** by choosing ϵb candidates randomly

Putting it all together

Algorithm 1: Learning to Optimize Tensor Programs

Input : Transformation space \mathcal{S}_e
Output: Selected schedule configuration s^*
 $\mathcal{D} \leftarrow \emptyset$

while $n_trials < max_n_trials$ **do**

 // Pick the next promising batch

$Q \leftarrow$ run parallel simulated annealing to collect candidates in \mathcal{S}_e using energy function \hat{f}

$S \leftarrow$ run greedy submodular optimization to pick $(1 - \epsilon)b$ -subset from Q by maximizing Equation 3

$S \leftarrow S \cup \{$ Randomly sample ϵb candidates. $\}$

 // Run measurement on hardware environment

for s **in** S **do**

 | $c \leftarrow f(g(e, s)); \mathcal{D} \leftarrow \mathcal{D} \cup \{(e, s, c)\}$

end

 // Update cost model

 update \hat{f} using \mathcal{D}

$n_trials \leftarrow n_trials + b$

end

$s^* \leftarrow$ history best schedule configuration

Use metaheuristic to optimize surrogate



Encourage diversity and exploration

Transfer learning

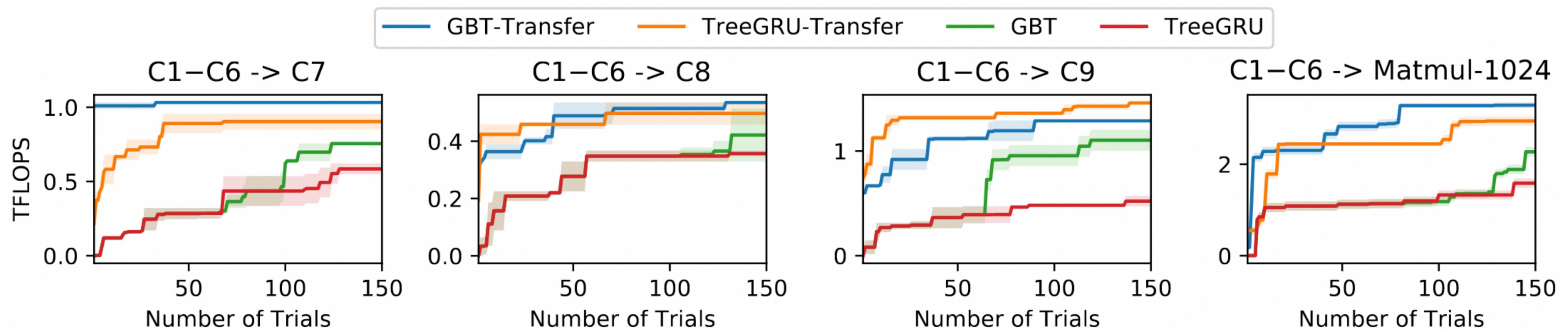
In practice, we need to optimize many workloads

Transfer what has been learned from previous workloads:

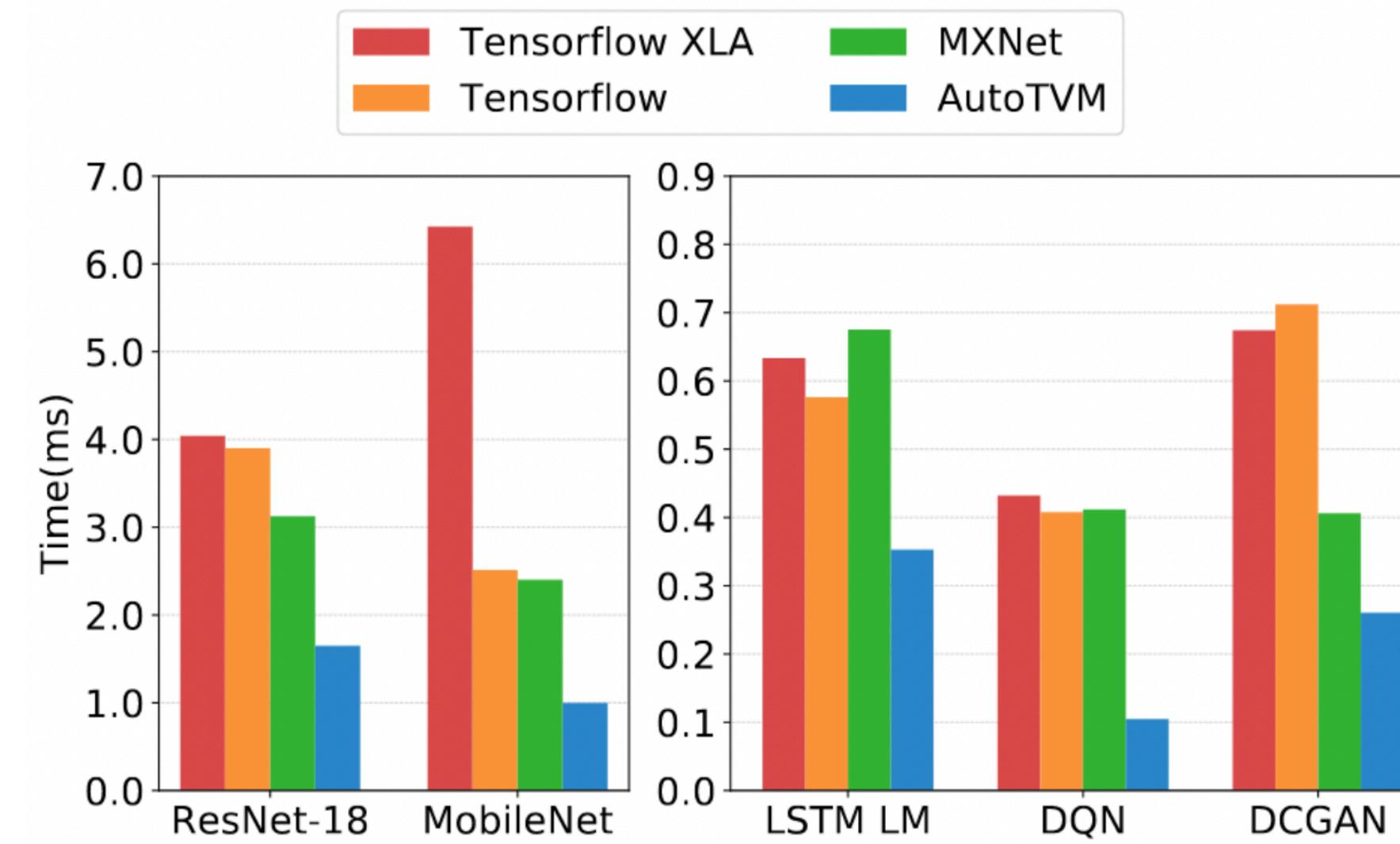
1. Use a **transferable representation**

2. Combine $\hat{f}(x) = \hat{f}^{(\text{global})}(x) + \hat{f}^{(\text{local})}(x)$

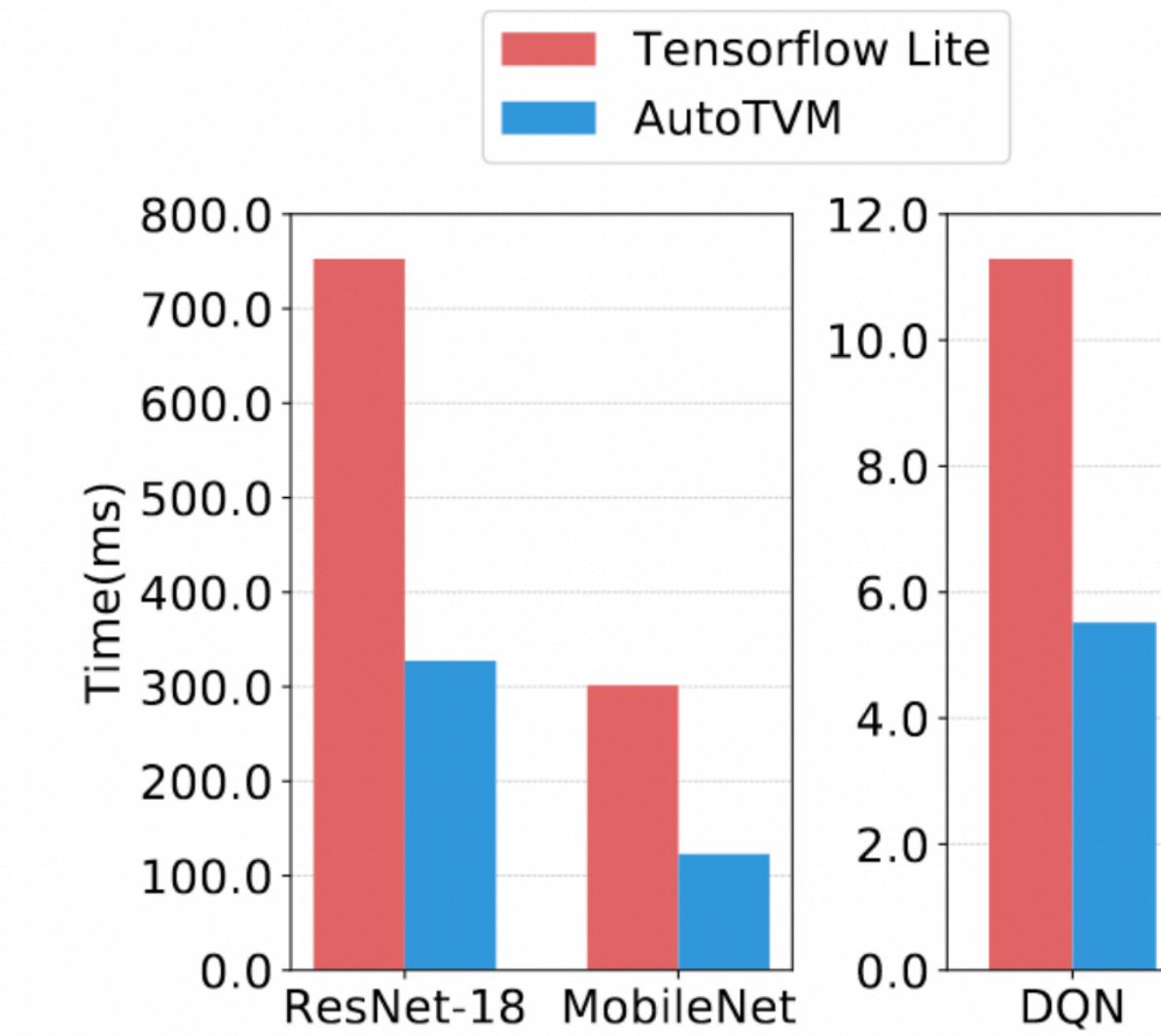
Results



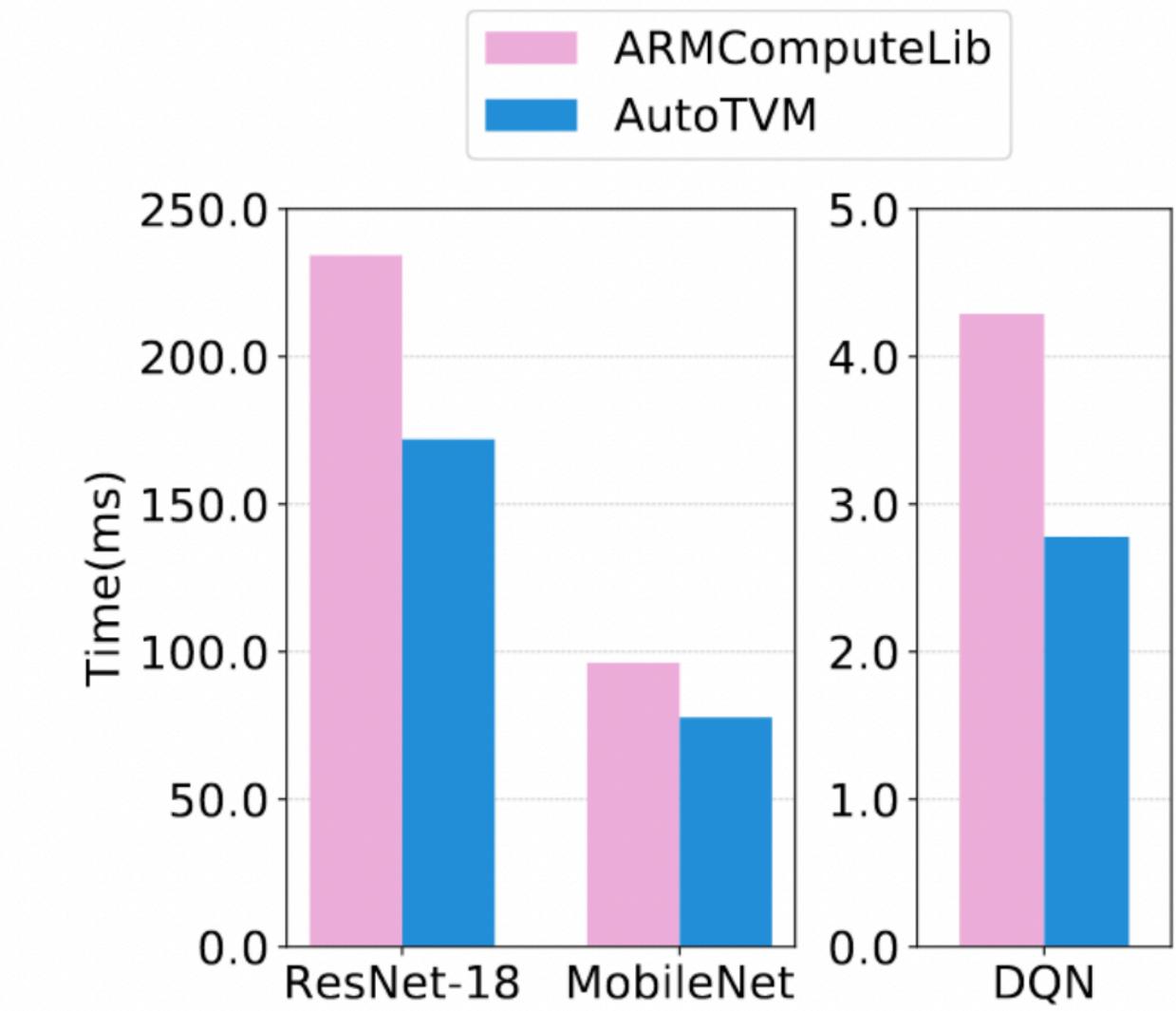
Results



(a) NVIDIA TITAN X End2End



(b) ARM Cortex-A53 End2End



(c) ARM Mali-T860 End2End

Summary

**By using an efficient cost model and transfer learning,
AutoTVM quickly and automatically finds efficient schedules**

...with some caveats

- 1. Need hand-crafted rules to express schedule templates**
- 2. Cost model relies on hand-crafted features**
- 3. Learning does not generalize across different domains**

Discussion questions

- AutoTVM uses transfer learning to speed up the optimization across workloads. Could we use transfer learning to speed up the optimization e.g. across different CPUs? What about from CPUs to GPUs?
- The GBT cost model in AutoTVM relies on a set of hand-crafted features. What are the drawbacks of using such a feature set?
- Hyperparameter optimization algorithms often use uncertainty estimates when choosing candidates. Why is this less important for AutoTVM?

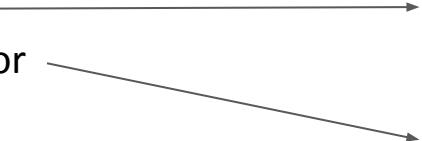
Ansor: Generating High-Performance Tensor Programs for Deep Learning

Lianmin Zheng, Chengfan Jia, Minmin Sun, Zhao Wu, Cody Hao Yu, Ameer Haj-Ali, Yida Wang,
Jun Yang, Danyang Zhuo, Koushik Sen, Joseph E. Gonzalez, Ion Stoica

Background

- Low-latency execution of Deep Neural Networks (DNN)
- Existing deep learning frameworks map the operators in DNNs to vendor-provided kernel libraries to achieve high performance
 - Growing diversity of hardware platforms
 - Hard to manually tune these libraries
- High-performance tensor programs are needed

Background

- Template-guided search
 - TVM
 - FlexTensor
 - Template design is a complicated process
 - **Manual** templates only cover **limited program structures**
 - Fails to include optimizations involving multiple operators
- Sequential construction based search
 - Halide
 - The cost model trained on complete programs **cannot accurately predict** the final performance of **incomplete** programs
 - The fixed order of sequential decisions **limits the search space**
 - Sequential construction based search is not scalable

Challenges - Ansor

Main deficiencies of existing solutions:

- Predefined manually-written templates
- Aggressive pruning and evaluating incomplete programs
- Limited rules are used to construct the search space

Search space not large enough!

Challenges:

- ★ Automatically construct a search space that is large enough
- ★ Search efficiently without comparing incomplete programs
- ★ Prioritize subgraphs critical to the end-to-end performance when optimizing an entire DNN with many subgraphs

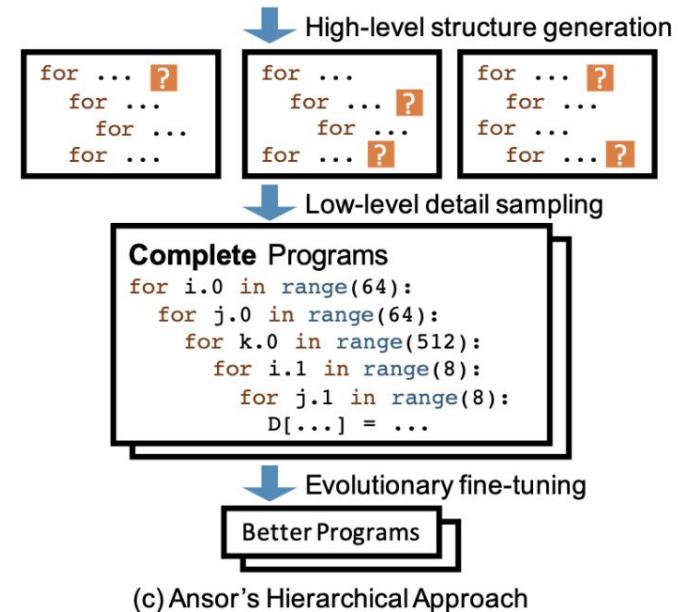
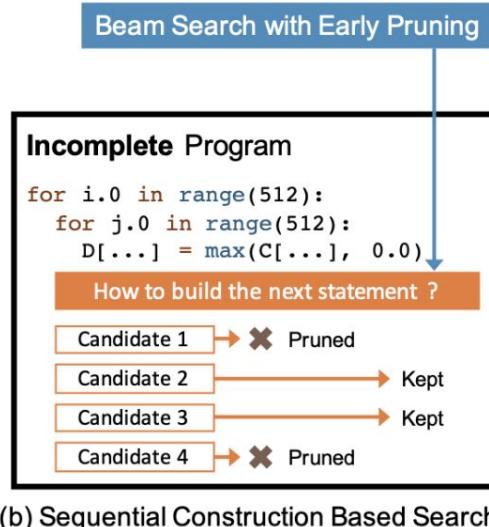
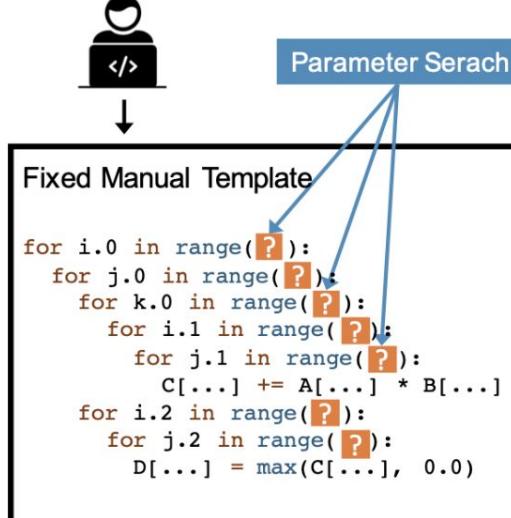


Figure 2: Search strategy comparison

Ansor

- Input: set of DNNs
 - Partitioned into small subgraphs
- Three components:
 - Program Sampler
 - Performance Tuner
 - Task scheduler

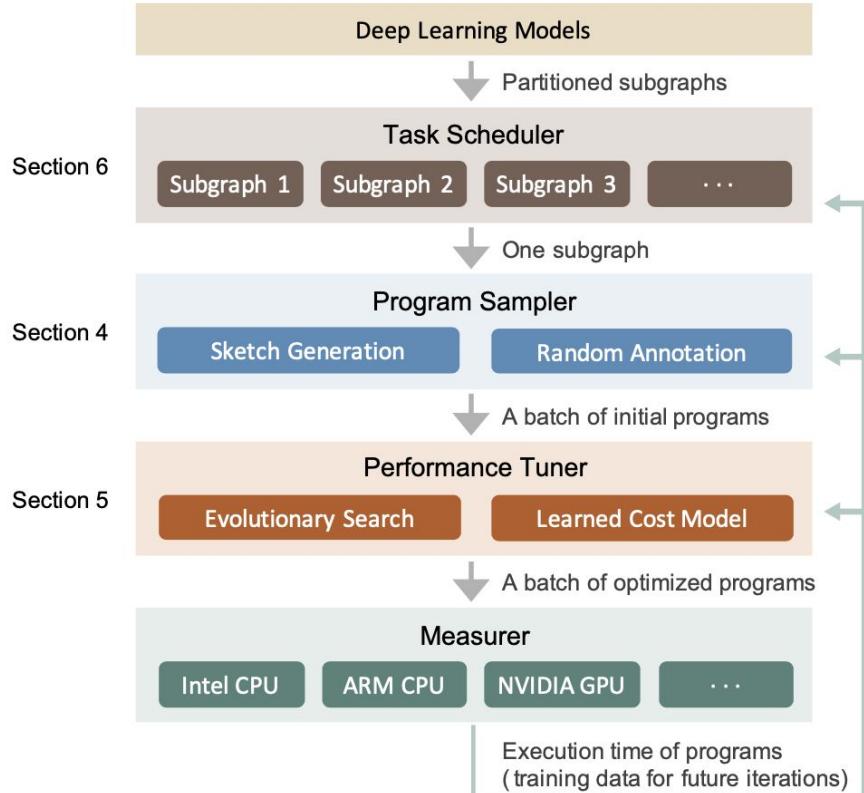


Figure 4: System Overview

Program Sampling

Program Sampler

Sketch Generation

Random Annotation

- Limited search space
 - Automatically expand the search space by recursively applying a set of flexible derivation rules
- Evaluating incomplete programs
 - Randomly sample complete programs in the search space

Hierarchical representation of the search space

- Top level: sketch
 - Generate sketches
- Lower level: annotation
 - Randomly annotate the sketches

Program Sampling

Program Sampler

Sketch Generation

Random Annotation

No	Rule Name	Condition	Application
1	Skip	$\neg IsStrictInlinable(S, i)$	$S' = S; i' = i - 1$
2	Always Inline	$IsStrictInlinable(S, i)$	$S' = Inline(S, i); i' = i - 1$
3	Multi-level Tiling	$HasDataReuse(S, i)$	$S' = MultiLevelTiling(S, i); i' = i - 1$
4	Multi-level Tiling with Fusion	$HasDataReuse(S, i) \wedge HasFusableConsumer(S, i)$	$S' = FuseConsumer(MultiLevelTiling(S, i), i); i' = i - 1$
5	Add Cache Stage	$HasDataReuse(S, i) \wedge \neg HasFusableConsumer(S, i)$	$S' = AddCacheWrite(S, i); i = i'$
6	Reduction Factorization	$HasMoreReductionParallel(S, i)$	$S' = AddRfactor(S, i); i' = i - 1$
...	User Defined Rule

Table 1: Derivation rules used to generate sketches

Example Input 1:

* The mathematical expression:

$$C[i, j] = \sum_k A[i, k] \times B[k, j]$$

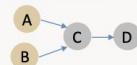
$$D[i, j] = \max(C[i, j], 0.0)$$

where $0 \leq i, j, k < 512$

* The corresponding naive program:

```
for i in range(512):
    for j in range(512):
        for k in range(512):
            C[i, j] += A[i, k] * B[k, j]
for i in range(512):
    for j in range(512):
        D[i, j] = max(C[i, j], 0.0)
```

* The corresponding DAG:



Example Input 2:

* The mathematical expression:

$$B[i, l] = \max(A[i, l], 0.0)$$

$$C[i, k] = \begin{cases} B[i, k], & k < 400 \\ 0, & k \geq 400 \end{cases}$$

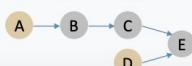
$$E[i, j] = \sum_k C[i, k] \times D[k, j]$$

where $0 \leq i < 8$, $0 \leq j < 4$,
 $0 \leq k < 512$, $0 \leq l < 400$

* The corresponding naive program:

```
for i in range(8):
    for l in range(400):
        B[i, l] = max(A[i, l], 0.0)
for i in range(8):
    for k in range(512):
        C[i, k] = B[i, k] if k < 400 else 0
for i in range(8):
    for j in range(4):
        for k in range(512):
            E[i, j] += C[i, k] * D[k, j]
```

* The corresponding DAG:



Generated sketch 1

```
for i_0 in range(TILE_I0):
    for j_0 in range(TILE_J0):
        for i_1 in range(TILE_I1):
            for j_1 in range(TILE_J1):
                for k_0 in range(TILE_K0):
                    for i_2 in range(TILE_I2):
                        for j_2 in range(TILE_J2):
                            for k_1 in range(TILE_K1):
                                for i_3 in range(TILE_I3):
                                    for j_3 in range(TILE_J3):
                                        C[...] += A[...] * B[...]
for i_4 in range(TILE_I2 * TILE_I3):
    for j_4 in range(TILE_J2 * TILE_J3):
        D[...] = max(C[...], 0.0)
```

Generated sketch 2

```
for i in range(8):
    for k in range(512):
        C[i, j] = max(A[i, k], 0.0) if k < 400 else 0
for i_0 in range(TILE_I0):
    for j_0 in range(TILE_J0):
        for i_1 in range(TILE_I1):
            for j_1 in range(TILE_J1):
                for k_0 in range(TILE_K0):
                    for i_2 in range(TILE_I2):
                        for j_2 in range(TILE_J2):
                            for k_1 in range(TILE_K1):
                                for i_3 in range(TILE_I3):
                                    for j_3 in range(TILE_J3):
                                        E.cache[...] += C[...] * D[...]
for i_4 in range(TILE_I2 * TILE_I3):
    for j_4 in range(TILE_J2 * TILE_J3):
        E[...] = E.cache[...]
```

Generated sketch 3

```
for i in range(8):
    for k in range(512):
        C[i, k] = max(A[i, k], 0.0) if k < 400 else 0
for i in range(8):
    for j in range(4):
        for k_o in range(TILE_K0):
            for k_i in range(TILE_K1):
                E.rf[...] += C[...] * D[...]
for i in range(8):
    for j in range(4):
        for k_i in range(TILE_K1):
            E[...] += E.rf[...]
```

Sampled program 1

```
parallel i_0@j_0@i_1@j_1 in range(256):
    for k_0 in range(32):
        for i_2 in range(16):
            unroll k_1 in range(16):
                unroll i_3 in range(4):
                    vectorize j_3 in range(16):
                        C[...] += A[...] * B[...]
for i_4 in range(64):
    vectorize j_4 in range(16):
        D[...] = max(C[...], 0.0)
```

Sampled program 2

```
parallel i_2 in range(16):
    for j_2 in range(128):
        for k_1 in range(512):
            for i_3 in range(32):
                vectorize j_3 in range(4):
                    C[...] += A[...] * B[...]
parallel i_4 in range(512):
    for j_4 in range(512):
        D[...] = max(C[...], 0.0)
```

Sampled program 3

```
parallel i_0 in range(8):
    for k in range(512):
        C[i, j] = max(A[i, k], 0.0)
        if k < 400 else 0
for k_0 in range(512):
    vectorize j_3 in range(4):
        E.cache[...] += C[...] * D[...]
vectorize j_4 in range(4):
    E[...] = E.cache[...]
```

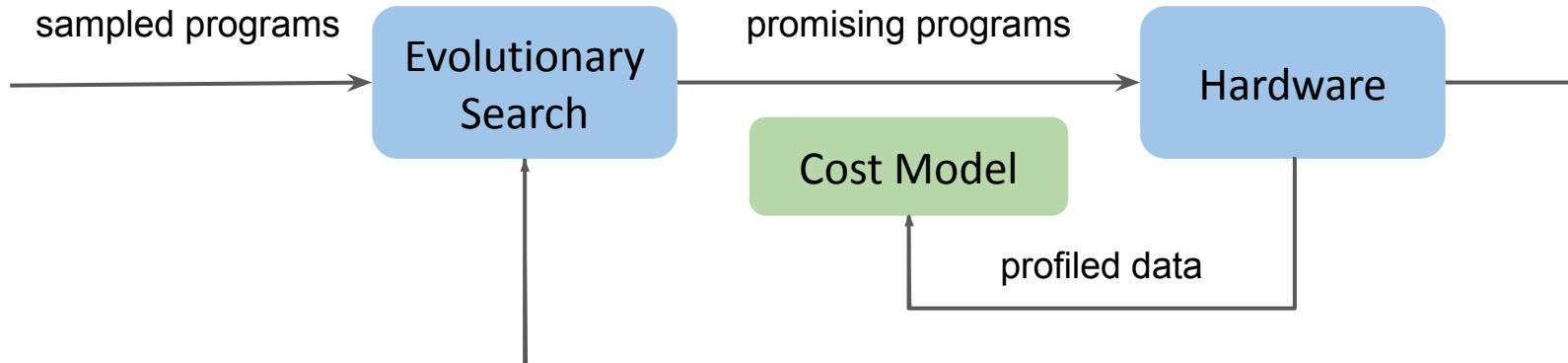
Sampled program 4

```
parallel i in range(8):
    for k in range(512):
        C[i, k] = ...
for j in range(4):
    unroll k_o in range(32):
        vectorized k_i in range(16):
            E.rf[...] += C[...] * D[...]
parallel i in range(8):
    for j in range(4):
        unroll k_i in range(16):
            E[...] += E.rf[...]
```

Performance Fine-Tuning



- Evolutionary search
- Learned cost model



Performance Fine-Tuning - Evolutionary Search

- Select some programs from the current generation according to certain probabilities
- Apply one of the evolution operations to generate a new program
 - Tile size mutation
 - Parallel mutation
 - Pragma mutation
 - Computation location mutation
 - Node-based crossover

Apply to general tensor programs and handle a search space with complicated dependency

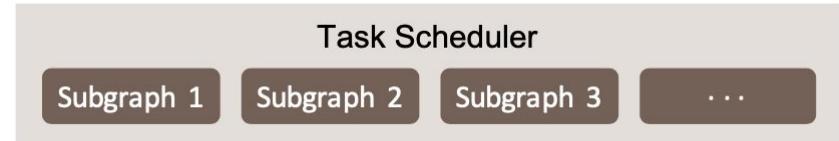
Perform out-of-order modifications to programs, addressing the sequential limitations

Performance Fine-Tuning - Learned Cost Model

$$loss(f, P, y) = w_p (\sum_{s \in S(P)} f(s) - y)^2 = y (\sum_{s \in S(P)} f(s) - y)^2$$

- Has great **portability** since a single model design can be reused for different hardware backends
- Gives relatively **accurate** estimations of the fitness of programs
- Querying the model is actually orders of magnitudes **faster** than the actual measurement

Task Scheduler



- Select the subgraphs that are more important to the overall performance
- A task: the process performed to generate high-performance programs for a subgraph
- One unit of time resources: one iteration of selecting a task, generating a batch of promising programs for the subgraph, and measuring the program on hardware

Task Scheduler

$t = (t_1, t_2, \dots, t_n)$, initialized to $t = (1, 1, \dots, 1)$

Objective: minimize the end-to-end cost, i.e. $\text{minimize } f(g_1(t), g_2(t), \dots, g_3(t))$

$$f_1 = \sum_{j=1}^m \sum_{i \in S(j)} w_i \times g_i(t)$$

$$f_2 = \sum_{j=1}^m \max(\sum_{i \in S(j)} w_i \times g_i(t), L_j)$$

$$f_3 = -\left(\prod_{j=1}^m \frac{B_j}{\sum_{i \in S(j)} w_i \times g_i(t)}\right)^{\frac{1}{m}}$$

$$f_4 = \sum_{j=1}^m \sum_{i \in S(j)} w_i \times \max(g_i(t), ES(g_i, t))$$

Table 2: Examples of objective functions for multiple neural networks

Task Scheduler

Approximate the gradient of f to choose the task $i = \operatorname{argmax}_i |\frac{\partial f}{\partial t_i}|$

$$\begin{aligned}\frac{\partial f}{\partial t_i} \approx \frac{\partial f}{\partial g_i} & \left(\alpha \frac{g_i(t_i) - g_i(t_i - \Delta t)}{\Delta t} + \right. \\ & \left. (1 - \alpha) \left(\min \left(-\frac{g_i(t_i)}{t_i}, \beta \frac{C_i}{\max_{k \in N(i)} V_k} - g_i(t_i) \right) \right) \right)\end{aligned}$$

Evaluation - Single Operator Benchmark

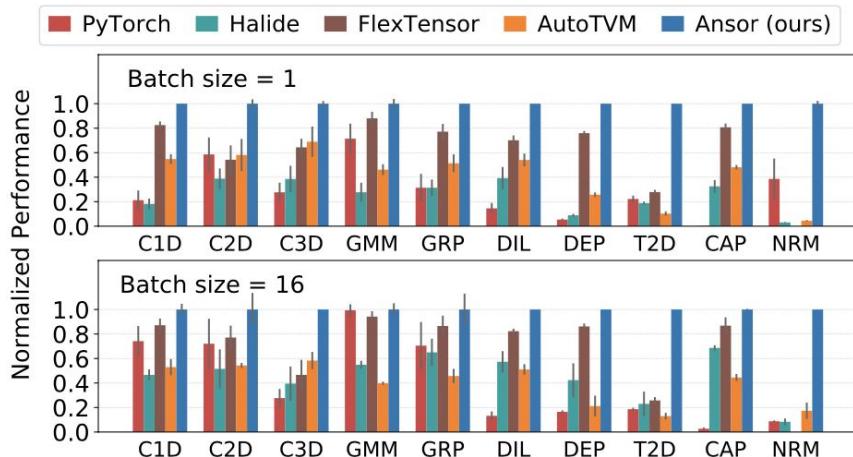


Figure 6: Single operator performance benchmark on a 20-core Intel-Platinum-8269CY

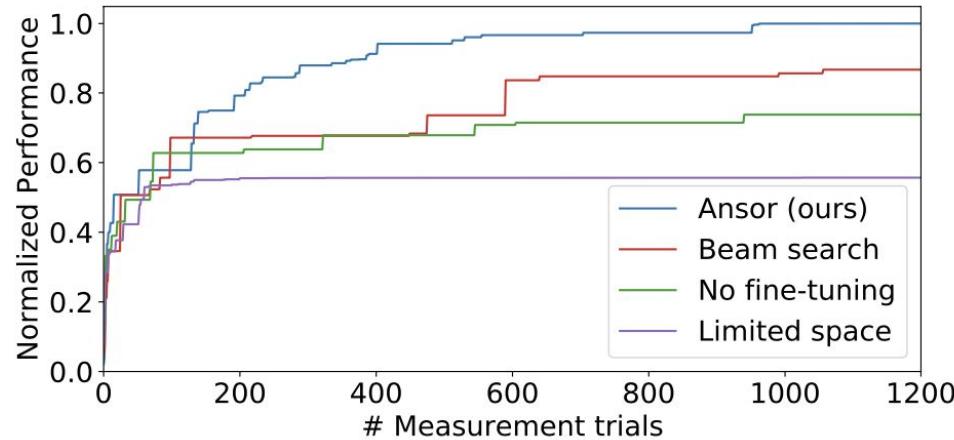


Figure 7: Ablation study of four variants of Ansor on a convolution operator

Evaluation - Subgraph Benchmark

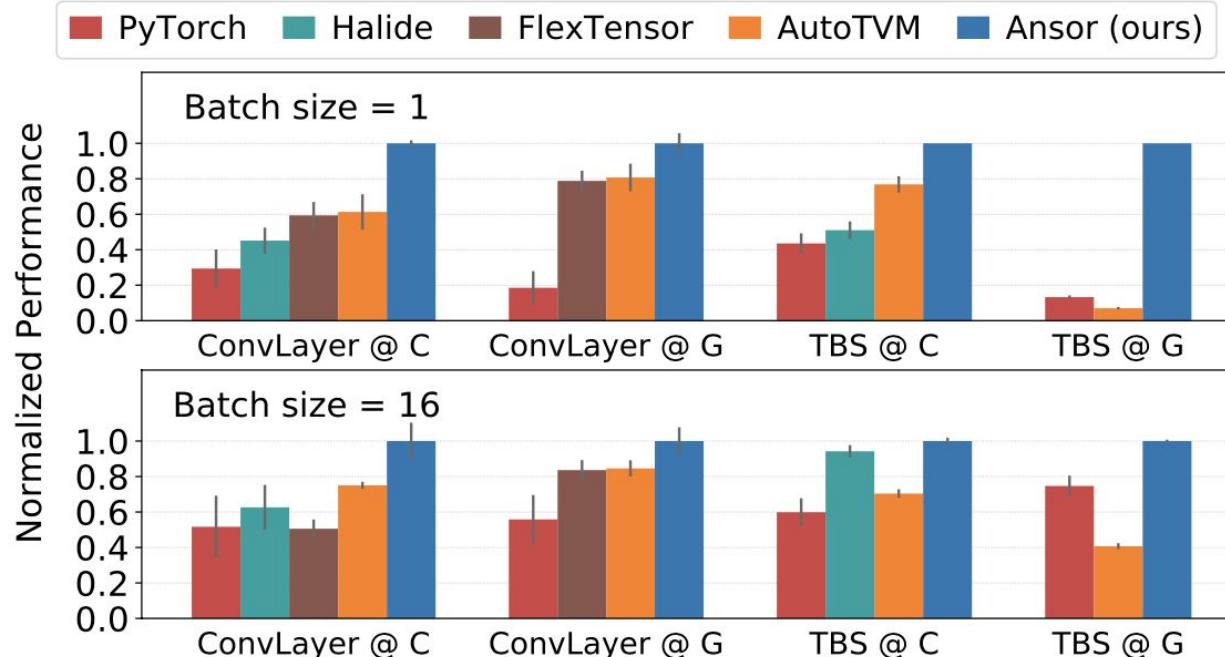


Figure 8: Subgraph performance benchmark on a 20-core Intel-Platinum-8269CY and an NVIDIA V100

Evaluation - End-to-End Network Benchmark

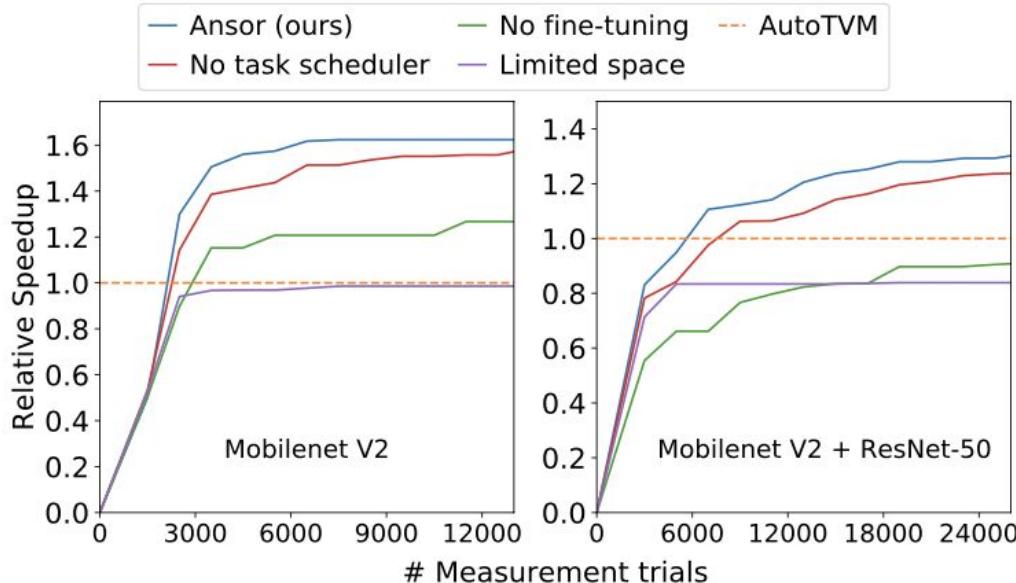
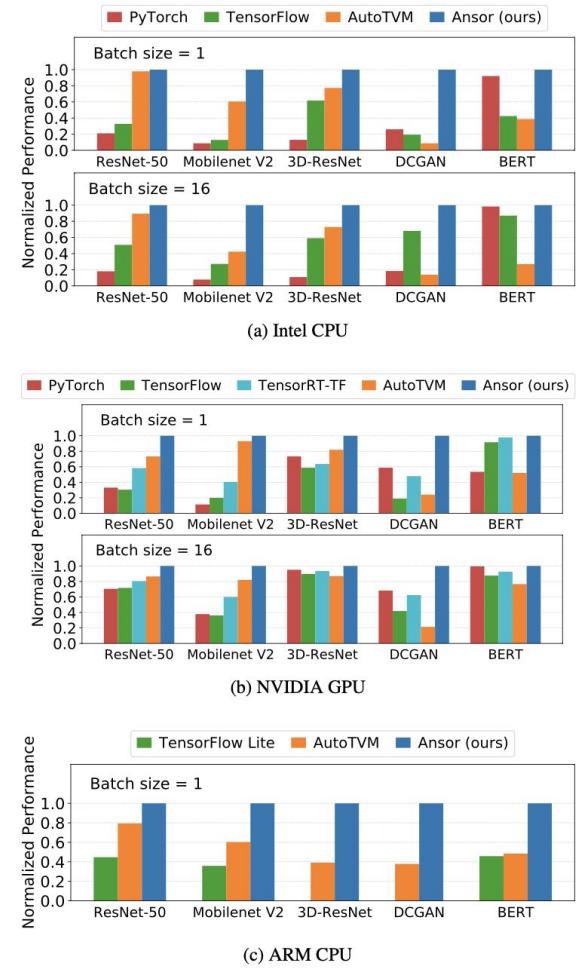


Figure 10: Network performance auto-tuning curve

Figure 9: Network inference performance benchmark on three hardware platforms



Evaluation - Search Time

	AutoTVM	Ansol	Time-saving
ResNet-50	21,220	6,403	3.3 ×
Mobilenet-V2	31,272	1,892	16.5 ×
3D-ResNet	5,158	1,927	2.7 ×
DCGAN	3,003	298	10.1 ×
BERT	6,220	496	12.5 ×

(a) The number of measurements.

	AutoTVM	Ansol	Time-saving
ResNet-50	39,250	4,540	8.6 ×
Mobilenet-V2	58,468	660	88.6 ×
3D-ResNet	7,594	2,296	3.3 ×
DCGAN	4,914	420	11.7 ×
BERT	12,007	266	45.1 ×

(b) Wall-clock time (seconds)

Table 3: The number of measurements and wall-clock time used for Ansol to match the performance of AutoTVM on the Intel CPU (batch size=1)

Evaluation - Cost Model Evaluation

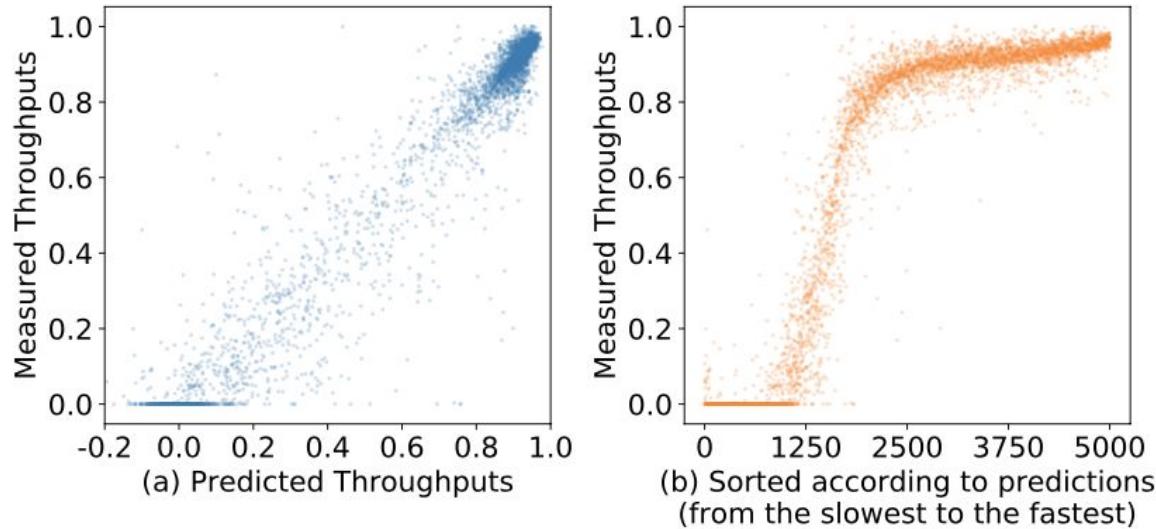


Figure 11: Measured throughputs vs. predicted throughputs.

Summary

Ansor: an automated search framework that generates high-performance tensor programs for deep neural networks

Compared with previous existing search strategies:

- Explores more optimization combinations by sampling programs from a hierarchical representation of the search space.
- Fine-tunes the sampled programs with evolutionary search and uses a learned cost model to identify the best programs.
- Utilizes a task scheduler to simultaneously optimize multiple subgraphs in deep neural networks.

Discussion

1. Besides XGBoost and the TreeGRU used in AutoTVM, is there any other model that might further improve the performance?
2. As stated in the paper, Ansor currently only supports dense operators. To support sparse operators that are commonly used in sparse neural networks and graph neural networks, how can we redesign the search space?