

TORCHSCRIPT: OPTIMIZED EXECUTION OF PYTORCH PROGRAMS

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O PyTorch Design Principles

Be Pythonic A first-class member of the python ecosystem, one idiomatic way of doing things.

Put Researchers First Easy APIs for models, data loaders, and optimizers. Hide implementation complexity.

Provide Pragmatic Performance A slowdown of 10% for a simpler API is acceptable; a 2x slowdown is not

Worse is better Save time by keeping the implementation simple, and write new features instead. A simple but incomplete solution is better than a complex one that is hard to maintain

O PyTorch Models are (differentiable) Python programs

```
class LinearLayer(nn.Module):
                                                           class FullBasicModel(nn.Module):
    def init (self, in sz, out sz):
                                                               def __init__(self):
        super(). init__()
                                                                   super(). init ()
        t1 = torch.randn(in sz, out sz)
                                                                   self.conv = nn.Conv2d(1, 128, 3)
        self.w = nn.Parameter(t1)
                                                                   self.fc = LinearLayer(128, 10)
        t2 = torch.randn(out sz)
        self.b = nn.Parameter(t2)
                                                              def forward(self, x):
                                                                   t1 = F.relu(self.conv(x))
    def forward(self, activations):
                                                                   t2 = self.fc(t1)
        t = torch.mm(activations, self.w)
                                                                   return F.softmax(t2)
        return t + self.b
```

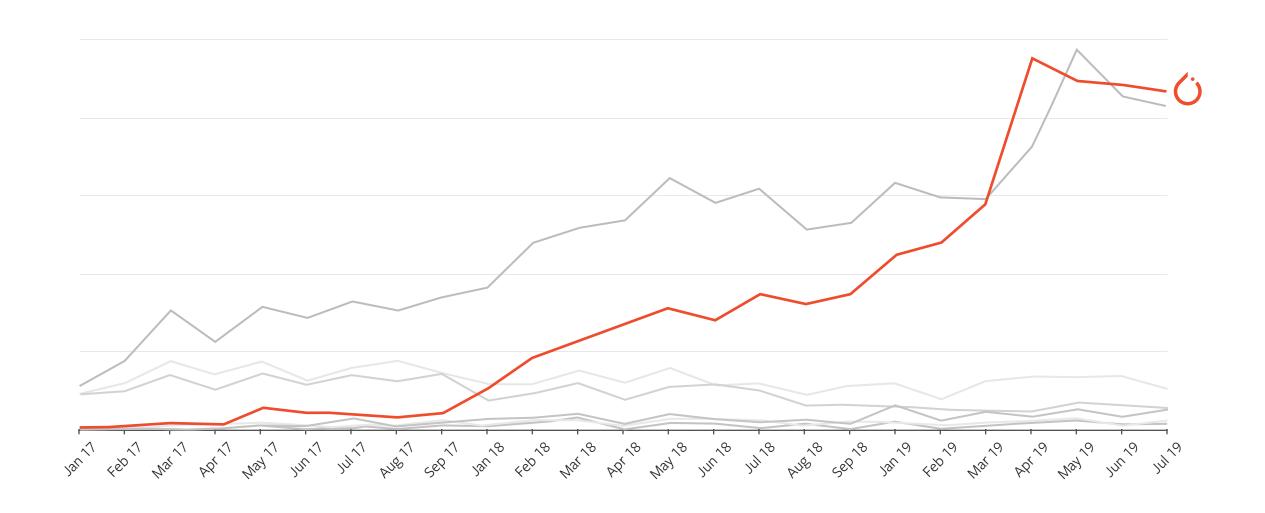
Why? Pythonic

- + Debuggable print and pdb
- + Hackable use any Python library

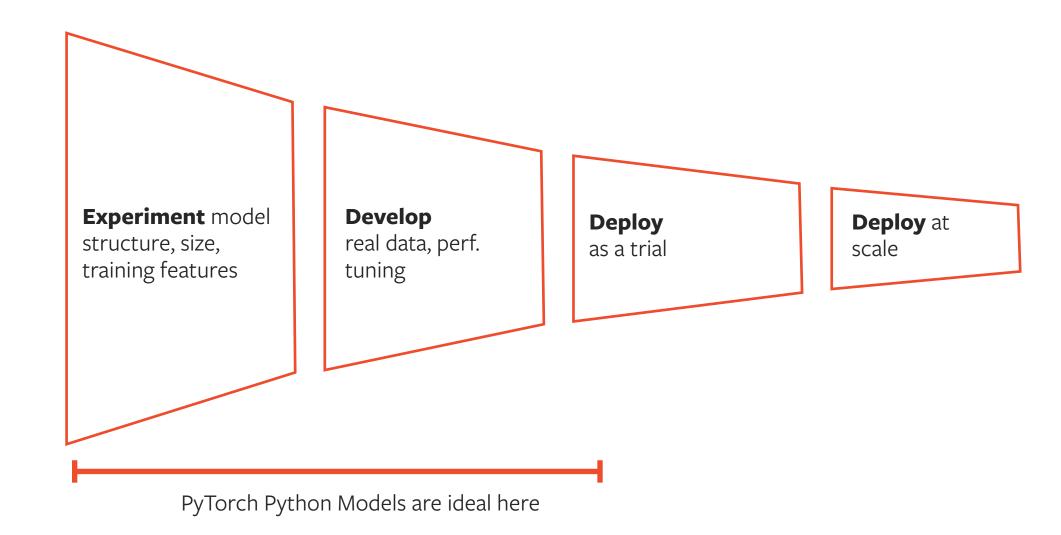
Uses well-understood object-oriented programming abstractions



GROWTH IN ARXIV MENTIONS IN RESEARCH PAPERS









REQUIREMENTS FOR DEPLOYING MODELS



PORTABILITY

Models should run anywhere



PERFORMANCE

Whole-program optimization



PROBLEM STATEMENT — WE NEED A SYSTEM THAT CAN:

1

CAPTURE THE STRUCTURE OF PYTORCH PROGRAMS.

USE THAT STRUCTURE TO OPTIMIZE.



PROBLEM STATEMENT — WE NEED A SYSTEM THAT CAN:

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CAPTURE THE STRUCTURE OF PYTORCH PROGRAMS.

TORCHSCRIPT

2

USE THAT STRUCTURE TO OPTIMIZE.

JIT COMPILER



PyTorch

Models are Python *programs*

- + Simple
- + Debuggable print and pdb
- + Hackable use any Python library
- Needs Python to run
- Difficult to optimize and parallelize

TorchScript

Models are Python programs, TorchScript

an optimizable subset of Python

- + Same "models are programs" approach
- + Production deployment
- + No Python dependency
- + Optimizable



Authoring TorchScript

Write model directly in a subset of Python

- AST-driven transformation
- Control-flow is preserved
- print statements can be used for debugging
- Remove the annotations to debug using standard Python tools.

```
class RNN(nn.Module):
  def __init__(self, W_h, U_h, W_y, b_h, b_y):
    super(RNN, self).__init__()
    self.W h = nn.Parameter(W h)
    self.U h = nn.Parameter(U h)
    self.W y = nn.Parameter(W y)
    self.b h = nn.Parameter(b h)
    self.b y = nn.Parameter(b y)
  def forward(self, x, h):
    V = []
    for t in range(x.size(0)):
      h = torch.tanh(x[t] @ self.W h + h @ self.U h + self.b h)
     y += [torch.tanh(h @ self.W y + self.b y)]
      if t % 10 == 0:
        print("stats: ", h.mean(), h.var())
    return torch.stack(y), h
script_rnn = torch.jit.script(RNN(W_h, U_h, W_y, b_h, b_y))
# save the compiled code and parameters so they can run elsewhere
script rnn.save("my rnn.pt")
```



Loading a model without Python

Torch Script models can be saved to a model archive, and loaded in a python-free executable using a C++ API.

Our C++ Tensor API is the same as our Python API, so you can do preprocessing and post processing before calling the model.



What subset of PyTorch is valid Torch Script?

- ✓ Static typing and type inference of all values
- √ Tensors and numeric primitives
- √ If statements
- √ Loops (and break, continue, return)
- ✓ User-defined classes with fixed attributes
- ✓ Tuples, Lists
- ✓ print and strings
- √ Gradients propagation through script functions

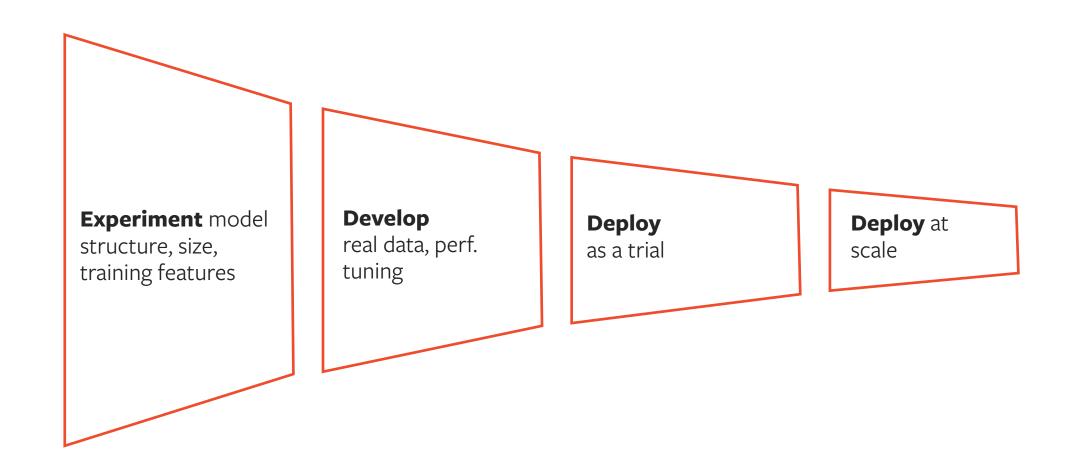
- ✓ In-place updates to tensors or lists
- ✓ All standard library nn. Modules like nn. Conv

- **X** Inheritance
- ✗ More complicated control-flow (e.g. generators)

For more details https://pytorch.org/docs/master/jit.html#torch-script-language-reference

C

Pay for what you use: Models only need to be in TorchScript for deployment.



O Python initialization, TorchScript inference

```
# 1. Define your model
class MyMod(torch.nn.Module):
    def __init__(self):
        ...
    def forward(self):
        ...

# 2. Create an instance of your model, and run init
my_nn_module = MyMod()
# 3. Convert your model to TorchScript
my_script_module = torch.jit.script(my_nn_module)
# 4. Run inference
output = my_script_module(input)
```

Model *initialization* is Python. *Inference* is TorchScript.



class ResNet(torch.nn.Module):

```
# Initialization code, written in Python
def __init__(self, block, layers, num_classes=1000):
    super(ResNet, self). init ()
    self.inplanes = 64
    self.conv1 = nn.Conv2d(3, 64, kernel_size=7, stride=2, padding=3,
                           bias=False)
    self.bn1 = nn.BatchNorm2d(64)
    self.relu = nn.ReLU(inplace=True)
    self.maxpool = nn.MaxPool2d(kernel size=3, stride=2, padding=1)
    self.layer1 = self. make layer(block, 64, layers[0])
    self.layer2 = self. make layer(block, 128, layers[1], stride=2)
    self.layer3 = self. make layer(block, 256, layers[2], stride=2)
    self.layer4 = self. make layer(block, 512, layers[3], stride=2)
    self.avgpool = nn.AdaptiveAvgPool2d((1, 1))
    self.fc = nn.Linear(512 * block.expansion, num classes)
    . . .
def make layer(self, block, planes, blocks, stride=1):
    downsample = None
    if stride != 1 or self.inplanes != planes * block.expansion:
        downsample = nn.Sequential(
            conv1x1(self.inplanes, planes * block.expansion, stride),
            nn.BatchNorm2d(planes * block.expansion),
    lavers = []
    layers.append(block(self.inplanes, planes, stride, downsample))
    self.inplanes = planes * block.expansion
    for in range(1, blocks):
        layers.append(block(self.inplanes, planes))
    return nn.Sequential(*layers)
```

```
# model code, written in TorchScript
def forward(self, x):
    x = self.conv1(x)
    x = self.bn1(x)
    x = self.relu(x)
    x = self.maxpool(x)

    x = self.layer1(x)
    x = self.layer2(x)
    x = self.layer3(x)
    x = self.layer4(x)

    x = self.avgpool(x)
    x = x.view(x.size(0), -1)
    x = self.fc(x)
```

Python initialization. TorchScript inference.



```
class RNN(nn.Module):
 def __init__(self, W_h, U_h, W_y, b_h, b_y):
    super(RNN, self). init ()
    self.W h = nn.Parameter(W h)
    self.U_h = nn.Parameter(U_h)
    self.W y = nn.Parameter(W y)
    self.b h = nn.Parameter(b h)
    self.b y = nn.Parameter(b y)
 def forward(self, x, h):
   y = []
    for t in range(x.size(0)):
     h = torch.tanh(x[t] @ self.W_h + h @ self.U_h + self.b_h)
     y += [torch.tanh(h @ self.W y + self.b y)]
     if t % 10 == 0:
        print("stats: ", h.mean(), h.var())
    return torch.stack(y), h
```

Control flow in forward always corresponds to dynamic execution in the model



Converting nn.Modules to TorchScript

```
script_rnn = torch.jit.script(RNN(W_h, U_h, W_y, b_h, b_y))
```

torch.jit.script takes a *fully initialized* nn.Module and converts it to TorchScript. The result is an instance of ScriptModule.

- 1. Parameters (self.weight, self.bias) are preserved
- 2. Submodules (self.layer1) are recursively converted
- 3. Attributes (self.training) are converted, if possible
- 4. Methods are converted into TorchScript, starting with the top-level module's forward method, and recursively converting any method it reaches. @torch.jit.export can set additional entry points for conversion

Model structure is *preserved* during conversion including: Function calls, objects, control-flow, leading to accurate stack traces.



CASE STUDY

Recurrent Neural Network Grammars

December 4, 2018

Improving Semantic Parsing for Task Oriented Dialog

Conversational Al Workshop at NeurIPS 2018

By: Arash Einolghozati, Panupong Pasupat, Sonal Gupta, Rushin Shah, Mrinal Mohit, Mike Lewis, Luke Zettlemoyer

- Complex dynamic behavior based on the inputs
- Typically written in pure C++



```
def forward(
    self,
    tokens: torch.Tensor,
    seq_lens: torch.Tensor,
    dict_feat: Tuple[torch.Tensor, torch.Tensor, torch.Tensor],
    actions: List[List[int]],
    contextual_token_embeddings: torch.Tensor,
    beam_size: int = 1,
    top_k: int = 1,
) -> List[Tuple[torch.Tensor, torch.Tensor]]:
    actions_idx = actions[0]
    assert len(actions_idx) > 0, "actions must be provided for training"
    token_embeddings = self.embedding(
        tokens, dict_feat, contextual_token_embeddings
    beam = [self.gen_init_state(tokens, token_embeddings)]
    all finished = False
    while not all_finished:
        # Stores plans for expansion as (score, state, action)
        plans : List[Plan] = []
        all_finished = True
        # Expand current beam states
        for state in beam:
            # Keep terminal states
            if state.finished():
                plans.append(Plan(state.neg_prob, const.TERMINAL_ELEMENT, state))
            else:
                all finished = False
                plans.extend(self.gen_plans(state))
```

heam clear()



COMPLEX CONTROL FLOW

```
while not all finished:
    for state in beam:
        if state.finished():
        else:
    for plan in plans[:beam_size]:
        beam.append(self.execute_plan(plan, actions_idx, beam_size))
```



USE COMMON DATA STRUCTURES

```
beam = [self.gen_init_state(tokens, token embeddings)]
    plans : List[Plan] = []
            plans.append(Plan(state.neg_prob, const.TERMINAL_ELEMENT, state))
            plans.extend(self.gen_plans(state))
    beam.clear()
    # Take actions to regenerate the beam
    plans.sort()
    for plan in plans[:beam_size]:
        beam.append(self.execute_plan(plan, actions_idx, beam_size))
beam.sort()
```



DEFINE YOUR OWN CLASSES

```
@torch.jit.scriptbeam:
class Plan:
    def <u>if</u>init<u>te</u>(self, hscore: float, action: int, state: ParserState):
        self.score = score an (state.neg prob, const.TERMINAL ELEMENT, state))
        self.action = action
        self.state = state
     def __lt__(self, other):
    beam#ctype: (Plan) -> bool
        return self.score < other.score
```



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```
import torch
class MyModule(torch.nn.Module):
   def init (self, N, M, state: List[Tensor]):
       super(MyModule, self). init ()
       self.weight = torch.nn.Parameter(torch.rand(N, M))
       self.state = state
   def forward(self, input):
       self.state.append(input)
       if input.sum() > 0:
            output = self.weight.mv(input)
       else:
            output = self.weight + input
       return output
# Compile the model code to a static representation
my module = MyModule(3, 4, [torch.rand(3, 4)])
my script module = torch.jit.script(my module)
# Save the compiled code and model data
# so it can be loaded elsewhere
my script module.save("my script module.pt")
```

TorchScript IR

```
graph(%self : ClassType<MyModule>,
     %input.1 : Tensor):
 %16 : int = prim::Constant[value=1]()
 %6 : None = prim::Constant()
 %8 : int = prim::Constant[value=0]()
 %2 : Tensor[] = prim::GetAttr[name="state"](%self)
 %4 : Tensor[] = aten::append(%2, %input.1)
 %7 : Tensor = aten::sum(%input.1, %6)
 %9 : Tensor = aten::gt(%7, %8)
 %10 : bool = aten::Bool(%9)
 %output : Tensor = prim::If(%10)
    block0():
     %11 : Tensor = prim::GetAttr[name="weight"](%self)
     %output.1 : Tensor = aten::mv(%11, %input.1)
     -> (%output.1)
    block1():
     %14 : Tensor = prim::GetAttr[name="weight"](%self)
     %output.2 : Tensor = aten::add(%14, %input.1, %16)
     -> (%output.2)
 return (%output)
```

O Improving Performance with TorchScript

Standard Compiler Passes

- Dead code elimination
- Constant propagation
- Common sub-expression elimination
- Loop unrolling

Tensor Optimizations

- Algebraic peephole optimizations
- Batching of matrix multiplications
- Point-wise fusions of element-wise operations

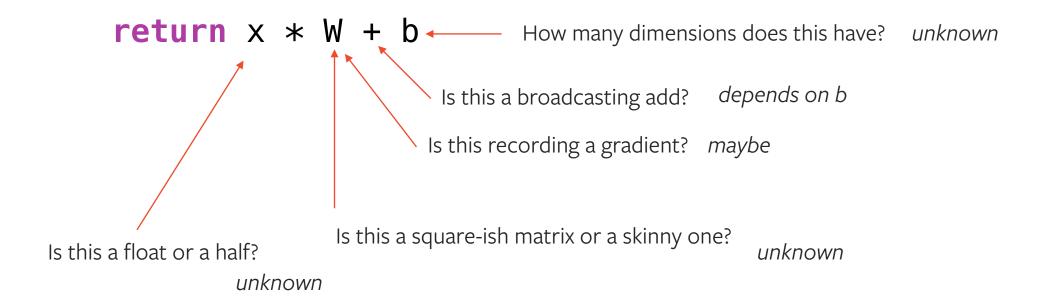
Runtime Optimization

- No global interpreter lock (GIL)
- fork/wait parallelism at the language level

```
@torch.jit.script
def LSTMCellS(x, hx, cx, w_ih, w_hh, b_ih, b_hh):
    x mm = x.mm(w ih.t())
    h mm = hx.mm(w hh.t())
    gates = x_mm + h_mm + b_ih + b_hh
    ingate, forgetgate, cellgate, outgate = gates.chunk(4, 1)
    ingate = torch.sigmoid(ingate)
    forgetgate = torch.sigmoid(forgetgate)
    cellgate = torch.tanh(cellgate)
    outgate = torch.sigmoid(outgate)
    cy = (forgetgate * cx) + (ingate * cellgate)
    hy = outgate * torch.tanh(cy)
    return hy, cy
graph(%x : Float(*, *)
      %hx : Float(*, *)
      %cx : Float(*, *)
      %w_ih : Float(*, *)
      %w_hh : Float(*, *)
      %b_ih : Float(*)
      %b_hh : Float(*)) {
    %9 : Float(*, *) = aten::t(%w_ih)
    %10 : Float(*, *) = aten::mm(\frac{1}{8}x, %9)
    %11 : Float(*, *) = aten::t(%w_hh)
%12 : Float(*, *) = aten::mm(%hx, %11)
    %77 : Tensor[] = prim::ListConstruct(%b_hh, %b_ih, %10, %12)
    %78 : Tensor[] = aten::broadcast_tensors(%77)
    %79 : Tensor, %80 : Tensor, %81 : Tensor, %82 : Tensor = prim::ListUnpack(%78)
    %hy: Float(*, *), %cy: Float(*, *) = prim::FusionGroup_0(%cx, %82, %81, %80, %79)
    %30 : (Float(*, *), Float(*, *)) = prim::TupleConstruct(%hy, %cy)
    return (%30);
```

Optimization through the dynamic behavior in Torch

def linear(x: Tensor, W: Tensor, b: Tensor) -> Tensor:



O Challenge Broadcasting

def should_i_fuse(x: Tensor, y: Tensor, z: Tensor) -> Tensor:

return x + y + z

Scenario 1

X : Float[1000]

Y: Float[1000]

Z: Float[1000]

Fused:

2000 ops

3000 reads from memory

Unfused:

2000 ops

4000 reads from memory

Scenario 2

X : Float[3]

Y: Float[3]

Z: Float[3x1000]

Fused:

6000 ops

3006 reads from memory

_ Unfused:

3000 ops

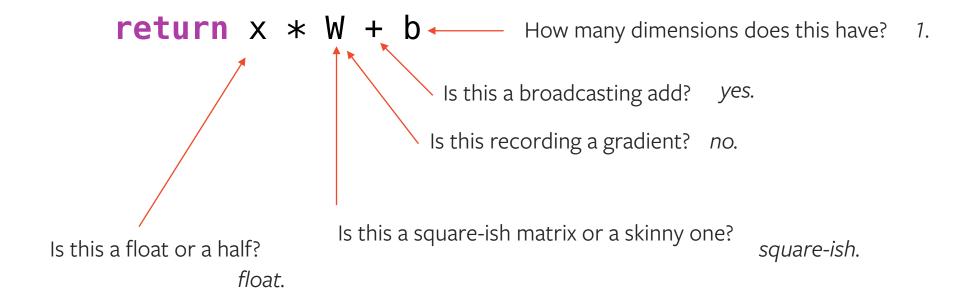
3006 reads from memory

OProfile Guided Optimization of TorchScript

While unknown statically, properties are very stable in practice.

We use **profile-guided** execution of TorchScript programs with **guarded optimistic** optimizations.

def linear(x: Tensor, W: Tensor, b: Tensor) -> Tensor:



Code in bold is fusible but we must know it runs on the GPU, it is floating point, and how tensors its are broadcast

```
graph(%self : torch .LSTMCell,
    %input.1 : Tensor,
    %state.1 : (Tensor, Tensor)):
%23 : int = prim::Constant[value=4]() # ../eg.py:24:60
%24 : int = prim::Constant[value=1]() # ../eq.py:24:63
%hx.1 : Tensor, %cx.1 : Tensor = prim::TupleUnpack(%state.1)
%7 : Tensor = prim::GetAttr[name="weight ih"](%self)
%8 : Tensor = aten::t(\%7) # ../eq.py:22:33
%9 : Tensor = aten::mm(%input.1, %8) # ../eq.py:22:17
%10 : Tensor = prim::GetAttr[name="bias ih"](%self)
%12 : Tensor = aten::add(%9, %10, %24) # ../eg.py:22:17
%14 : Tensor = prim::GetAttr[name="weight hh"](%self)
%15 : Tensor = aten::t(%14) # ../eq.py:23:30
%16 : Tensor = aten::mm(%hx.1, %15) # ../eg.py:23:17
%18 : Tensor = aten::add(%12, %16, %24) # ../eg.py:22:17
%19 : Tensor = prim::GetAttr[name="bias hh"](%self)
%gates.1 : Tensor = aten::add(%18, %19, %24) # ../eg.py:22:1
. . .
```

```
graph(%self : torch .LSTMCell,
    %input.1 : Tensor,
   %state.1 : (Tensor, Tensor)):
%4 : int = prim::Constant[value=1]() # ../eg.py:24:63
%hx.1 : Tensor, %cx.1 : Tensor = prim::TupleUnpack(%state.1)
%7 : Tensor = prim::GetAttr[name="weight ih"](%self)
%8 : Tensor = prim::profile(%7)
%9 : Tensor = aten::t(%8) # ../eg.py:22:33
%10 : Tensor = prim::profile(%input.1)
%11 : Tensor = prim::profile(%9)
12: Tensor = aten::mm(10, 11) # ../eq.py:22:17
%13 : Tensor = prim::GetAttr[name="bias ih"](%self)
%14 : Tensor = prim::profile(%12)
%15 : Tensor = prim::profile(%13)
%16 : Tensor = aten::add(%14, %15, %4) # ../eg.py:22:17
%17 : Tensor = prim::GetAttr[name="weight hh"](%self)
%18 : Tensor = prim::profile(%17)
%19 : Tensor = aten::t(%18) # ../eq.py:23:30
%20 : Tensor = prim::profile(%hx.1)
%21 : Tensor = prim::profile(%19)
22: Tensor = aten::mm(20, 21) # ../eq.py:23:17
%23 : Tensor = prim::profile(%16)
%24 : Tensor = prim::profile(%22)
%25 : Tensor = aten::add(%23, %24, %4) # ../eg.py:22:17
%26 : Tensor = prim::GetAttr[name="bias hh"](%self)
%27 : Tensor = prim::profile(%25)
%28 : Tensor = prim::profile(%26)
%gates.1 : Tensor = aten::add(%27, %28, %4) # ../eg.py:22:17
```

(1) Insert profiling code at every use of a Tensor

```
graph(%self : __torch__.LSTMCell,
    %input.1 : Tensor,
    %state.1 : (Tensor, Tensor)):
%3 : int = prim::Constant[value=4]() # ../eq.py:24:60
%4 : int = prim::Constant[value=1]() # ../eg.py:24:63
%hx.1 : Tensor, %cx.1 : Tensor = prim::TupleUnpack(%state.1)
%7 : Tensor = prim::GetAttr[name="weight ih"](%self)
%8 : Float(40, 10) = prim::profile(%7)
%9 : Tensor = aten::t(%8) # ../eq.py:22:33
%10 : Float(8, 10) = prim::profile(%input.1)
%11 : Float(10, 40) = prim::profile(%9)
12: Tensor = aten::mm(10, 11) # ../eq.py:22:17
%13 : Tensor = prim::GetAttr[name="bias ih"](%self)
%14 : Float(8, 40) = prim::profile(%12)
%15 : Float(40) = prim::profile(%13)
%16 : Tensor = aten::add(%14, %15, %4) # ../eg.py:22:17
%17 : Tensor = prim::GetAttr[name="weight hh"](%self)
%18 : Float(40, 10) = prim::profile(%17)
%19: Tensor = aten::t(%18) # ../eg.py:23:30
%20 : Float(8, 10) = prim::profile(%hx.1)
%21 : Float(10, 40) = prim::profile(%19)
%22 : Tensor = aten::mm(%20, %21) # ../eg.py:23:17
%23 : Float(8, 40) = prim::profile(%16)
%24 : Float(8, 40) = prim::profile(%22)
%25 : Tensor = aten::add(%23, %24, %4) # ../eg.py:22:17
%26 : Tensor = prim::GetAttr[name="bias hh"](%self)
%27 : Float(8, 40) = prim::profile(%25)
%28 : Float(40) = prim::profile(%26)
%gates.1 : Tensor = aten::add(%27, %28, %4) # ../eg.py:22:17
```

(2) Run the graph a few times to record sizes

```
graph(%self : __torch__.LSTMCell,
   %input.1 : Tensor,
   %state.1 : (Tensor, Tensor)):
   %3 : int = prim::Constant[value=4]() # ../eg.py:24:60
   %4 : int = prim::Constant[value=1]() # ../eg.py:24:63
   %hx.1 : Tensor, %cx.1 : Tensor = prim::TupleUnpack(%state.1)
   %7 : Tensor = prim::GetAttr[name="weight ih"](%self)
                                                                    guard failure!
   %8 : Float(40, 10) = prim::guard(%7)
   %9 : Tensor = aten::t(%8) # ../eq.py:22:33
   %10 : Float(8, 10) = prim::quard(%input.1)
   %11 : Float(10, 40) = prim::guard(%9)
   %12 : Tensor = aten::mm(%10, %11) # ../eq.py:22:17
                                                                 Unoptimized Fallback
    \cdot
                                                                 %15: Tensor = aten::t(%14) # ../eq.py
                                                                 %16 : Tensor = aten::mm(%hx.1, %15) #
                                                                 %18 : Tensor = aten::add(%12, %16, %24
                                                                 . . .
```

(3) Replace profile nodes with guards. If a guard fails during execution, we fallback to the unoptimized code.

```
(
```

```
graph(%self : torch .LSTMCell,
      %input.1 : Tensor,
      %state.1: (Tensor, Tensor)):
 %98 : Float(40) = prim::guard(%26, %25, %93)
 %122 : Tensor[] = prim::ListConstruct(%25, %98)
 %123 : Tensor[] = aten::broadcast tensors(%122)
 %124 : Tensor, %125 : Tensor = prim::ListUnpack(%123)
 %hy.1 : Float(8, 10), %cy.1 : Float(8, 10) = prim::FusionGroup 1(%93, %125, %124)
 %60 : (Tensor, Tensor) = prim::TupleConstruct(%hy.1, %cy.1)
 %62 : (Tensor, (Tensor, Tensor)) = prim::TupleConstruct(%hy.1, %60)
  return (%62)
with prim::FusionGroup 1 = graph(%13 : Float(8, 10),
      %39: Tensor,
     %44 : Tensor):
 %45 : Float(8, 10), %46 : Float(8, 10), %47 : Float(8, 10), %48 : Float(8, 10) = prim::ConstantChunk[chunks=4, dim=1
 %40 : Float(8, 10), %41 : Float(8, 10), %42 : Float(8, 10), %43 : Float(8, 10) = prim::ConstantChunk[chunks=4, dim=1
 %37 : int = prim::Constant[value=1]() # ../eq.py:24:63
 %38 : Float(8, 10) = aten::add(%45, %40, %37)
 %34 : Float(8, 10) = aten::add(%46, %41, %37)
 %30 : Float(8, 10) = aten::add(%47, %42, %37)
 %26 : Float(8, 10) = aten::add(%48, %43, %37)
 %ingate.3 : Float(8, 10) = aten::sigmoid(%38) # ../eg.py:26:17
 %forgetgate.3 : Float(8, 10) = aten::sigmoid(%34) # ../eg.py:27:21
 cellgate.3 : Float(8, 10) = aten::tanh(%30) # ../eg.py:28:19
  \text{%outgate.3}: Float(8, 10) = aten::sigmoid(%26) # ../eg.py:29:18
  return (%hy.1, %cy.1)
```

(4) Remove redundant guards, and use profiled properties to apply fusion.

O Profile-guided optimization

Possibilities

If we know tensors are constant, we can pre-multiplying weights to remove batch norms or load weights into grams on accelerators.

If we know that the bool that is input to an if-statement is almost always true, we can eliminate the other branch from the optimized code.



WHAT'S NEXT?

TORCHSCRIPT AS A PLATFORM



QUANTIZATION

Model quantization done safely and automatically using JIT transformations.



MOBILE

A lightweight interpreter that can run on-device.



BACKENDS

Support for lowering models to static graph compilers, like TVM, Glow, XLA.



TRY IT

AND GIVE US FEEDBACK!



TUTORIALS

pytorch.org/tutorials

Introduction to TorchScript: https://pytorch.org/tutorials/beginner/ Intro_to_TorchScript_tutorial.html

Loading a TorchScript model in C++: https://pytorch.org/tutorials/advanced/ cpp_export.html



DOCUMENTATION

TorchScript reference: https://pytorch.org/docs/master/jit.html



FEEDBACK

"jit" label on github: https://github.com/pytorch/pytorch/issues? q=is%3Aissue+is%3Aopen+label%3Ajit