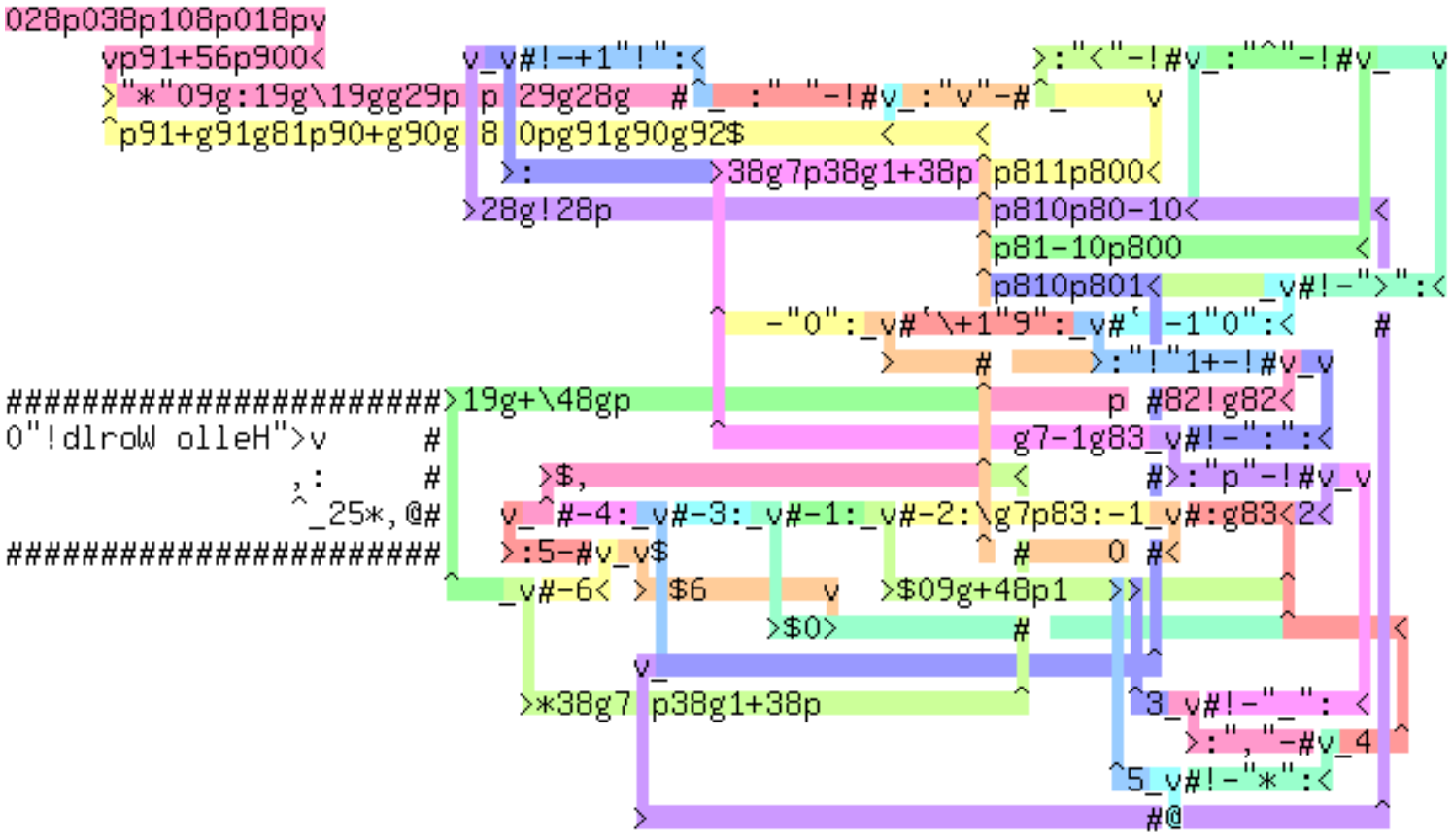
No more mini-languages: Autodiff in full-featured Python



David Duvenaud, Dougal Maclaurin, Matthew Johnson

Our awesome new world

- TensorFlow, Stan, Theano, Edward
- Only need to specify forward model
- Autodiff + inference / optimization done for you

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- a new language to learn

Autograd

github.com/HIPS/autograd

- differentiates native Python code
- handles most of Numpy + Scipy
- loops, branching, recursion, closures
- arrays, tuples, lists, dicts...
- derivatives of derivatives
- a one-function API!

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About 2000 downloads/month, used for...

- Population genetics simulations
- Inference libraries
- Protein folding simulations
- Material thermodynamics simulations
- Optimization on manifolds
- Neural Turing machine

Simple gradient code

```
anp.sinh.defvjp(lambda g, ans, vs, gvs, x : g * anp.cosh(x))
anp.cosh.defvjp(lambda g, ans, vs, gvs, x : g * anp.sinh(x))
anp.tanh.defvjp(lambda g, ans, vs, gvs, x : g / anp.cosh(x)**2)

anp.cross.defvjp(lambda g, ans, vs, gvs, a, b,
                       axisa=-1, axisb=-1, axisc=-1, axis=None :
                   anp.cross(b, g, axisb, axisc, axisa, axis), argnum=0)

def grad_sort(g, ans, vs, gvs, x, axis=-1,
                    kind='quicksort', order=None):
    sort_perm = anp.argsort(x, axis, kind, order)
    return unpermuter(g, sort_perm)
anp.sort.defvjp(grad_sort)
```

Most Numpy functions implemented

Complex & Fourier	Array	Misc	Linear Algebra	Stats
imag	atleast_1d	logsumexp	inv	std
conjugate	atleast_2d	where	norm	mean
angle	atleast_3d	einsum	det	var
real_if_close	full	sort	eigh	prod
real	repeat	partition	solve	sum
fabs	split	clip	trace	cumsum
fft	concatenate	outer	diag	norm
fftshift	roll	dot	tril	t
fft2	transpose	tensordot	triu	dirichlet
ifftn	reshape	rot90	cholesky	
ifftshift	squeeze			
ifft2	ravel			
ifft	expand_dims			

Autograd examples

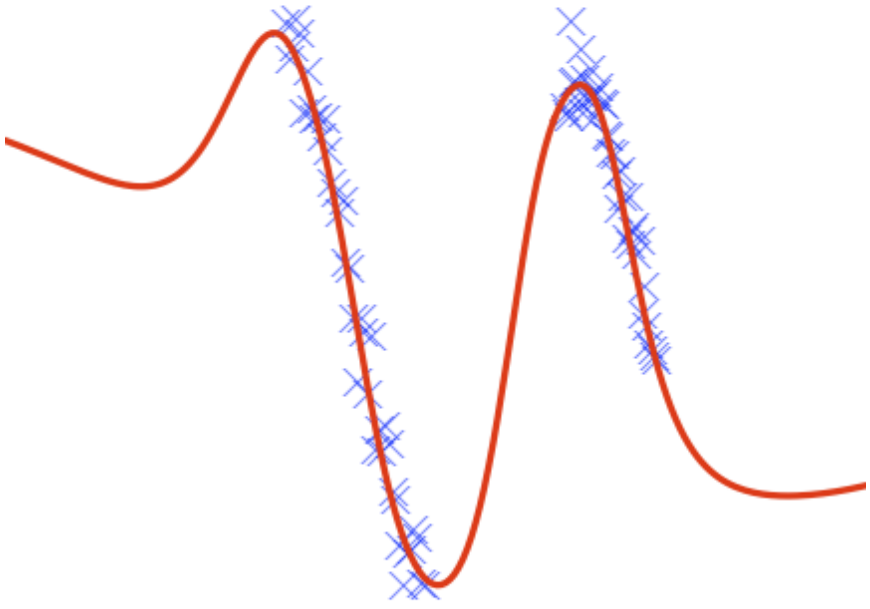
```
import autograd.numpy as np
from autograd import grad

def predict(weights, inputs):
    for W, b in weights:
        outputs = np.dot(inputs, W) + b
        inputs = np.tanh(outputs)
    return outputs

def init_params(scale, sizes):
    return [(npr.randn(nin, out) * scale,
             npr.randn(out) * scale)
            for nin, out in
            zip(sizes[:-1], sizes[1:])]

def logprob_func(weights, inputs, targets)
    preds = predict(weights, inputs)
    return np.sum((preds - targets)**2)

gradient_func = grad(logprob_func)
```



Structured gradients

```
print(grad(logprob)(init_params, inputs, targets))

[(array([[ -5.40710861, -14.13507334, -13.94789859,  28.6188964 ]]),
   array([-17.01486765, -28.33800594, -29.77875615,  49.78987454])),
 (array([[-71.47406027, -69.1771986 ,  -7.34756845, -17.96280387],
         [ 21.90645613,  22.01415812,   2.37750145,   5.81340489],
         [-39.37357205, -38.07711948,  -4.04245488,  -9.88483908],
         [-27.00357209, -24.79890695,  -2.56954539,  -6.28235645]]),
  array([-281.99906027, -278.86794587,  -29.90316231,  -73.12033635])),
  (array([[-410.89215947],
          [ 256.31407037],
          [ -31.39182332],
          [   6.89045123]]),
  array([-1933.60342748]))]
```

Structured gradients

```
def adam(grad, init_params, callback=None, num_iters=100,
         step_size=0.001, b1=0.9, b2=0.999, eps=10**-8):

    flattened_grad, unflatten, x = flatten_func(grad, init_params)

    m = np.zeros(len(x))
    v = np.zeros(len(x))
    for i in range(num_iters):
        g = flattened_grad(x, i)
        if callback: callback(unflatten(x), i, unflatten(g))
        m = (1 - b1) * g      + b1 * m  # First  moment estimate.
        v = (1 - b2) * (g**2) + b2 * v  # Second moment estimate.
        mhat = m / (1 - b1**(i + 1))    # Bias correction.
        vhat = v / (1 - b2**(i + 1))
        x = x - step_size*mhat/(np.sqrt(vhat) + eps)
    return unflatten(x)
```

How to code a Hessian-vector product?

```
def hvp(func):
    def vector_dot_grad(arg, vector):
        return np.dot(vector, grad(func)(arg))
    return grad(vector_dot_grad)
```

- hvp(f)(x, v) returns $\mathbf{v}^T \nabla_{\mathbf{x}} \nabla_{\mathbf{x}}^T f(\mathbf{x})$
- No explicit Hessian
- Can construct higher-order operators easily

What about full Hessian?

```
def hessian(fun):
    return jacobian(jacobian(fun))
```

```
def project(vx, vy):
    # Project the velocity field to be approximately mass-conserving,
    # using a few iterations of Gauss-Seidel.
    p = np.zeros(vx.shape)
    h = 1.0/vx.shape[0]
    div = -0.5 * h * (np.roll(vx, -1, axis=0) - np.roll(vx, 1, axis=0)
                    + np.roll(vy, -1, axis=1) - np.roll(vy, 1, axis=1))
    for k in range(10):
        p = (div + np.roll(p, 1, axis=0) + np.roll(p, -1, axis=0)
                 + np.roll(p, 1, axis=1) + np.roll(p, -1, axis=1))/4.0
    vx -= 0.5*(np.roll(p, -1, axis=0) - np.roll(p, 1, axis=0))/h
    vy -= 0.5*(np.roll(p, -1, axis=1) - np.roll(p, 1, axis=1))/h
    return vx, vy

def advect(f, vx, vy):
    # Move field f according to x and y velocities (u and v)
    # using an implicit Euler integrator.
    rows, cols = f.shape
    cell_ys, cell_xs = np.meshgrid(np.arange(rows),
                                   np.arange(cols))
    center_xs = (cell_xs - vx).ravel()
    center_ys = (cell_ys - vy).ravel()

    # Compute indices of source cells.
    left_ix = np.floor(center_xs).astype(int)
    top_ix  = np.floor(center_ys).astype(int)
    rw = center_xs - left_ix
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    left_ix  = np.mod(left_ix,     rows)
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    flat_f = (1 - rw) * ((1 - bw)*f[left_ix,  top_ix] \
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    return np.reshape(flat_f, (rows, cols))

def simulate(vx, vy, smoke, num_time_steps):
    for t in range(num_time_steps):
        vx_updated = advect(vx, vx, vy)
        vy_updated = advect(vy, vx, vy)
        vx, vy = project(vx_updated, vy_updated)
        smoke = advect(smoke, vx, vy)
    return smoke, frame_list
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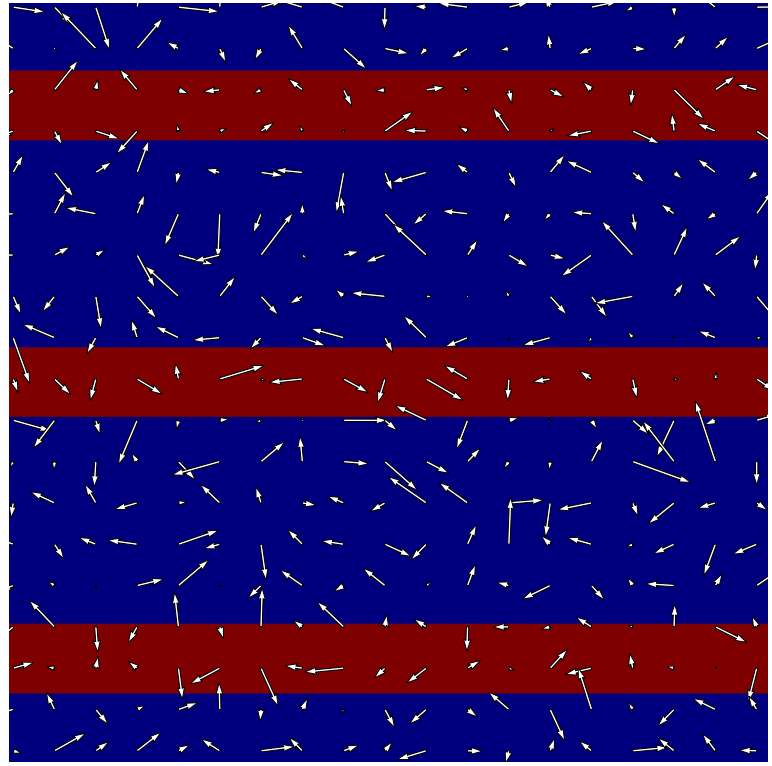

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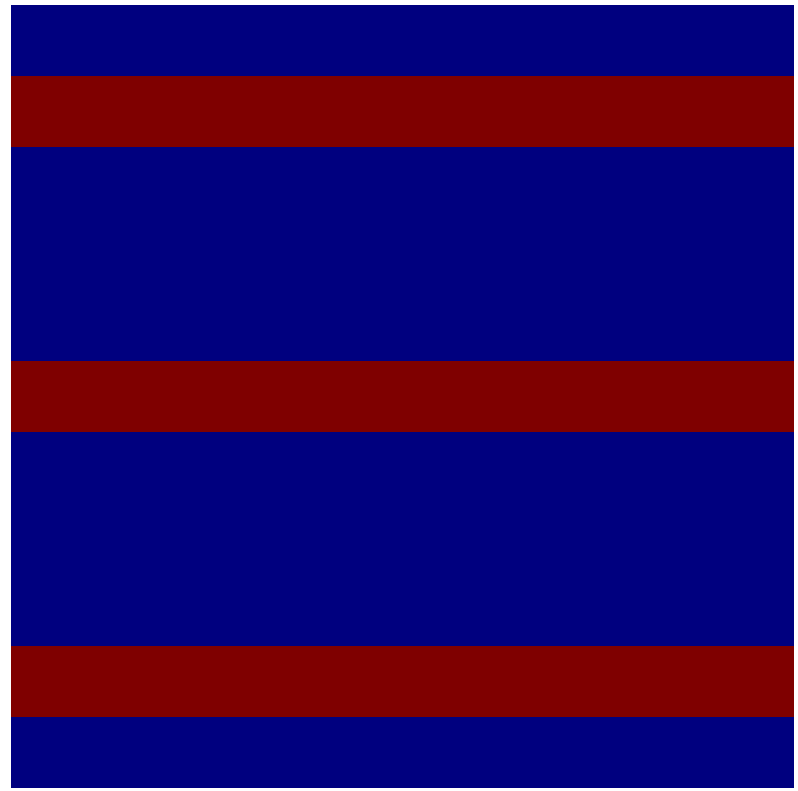

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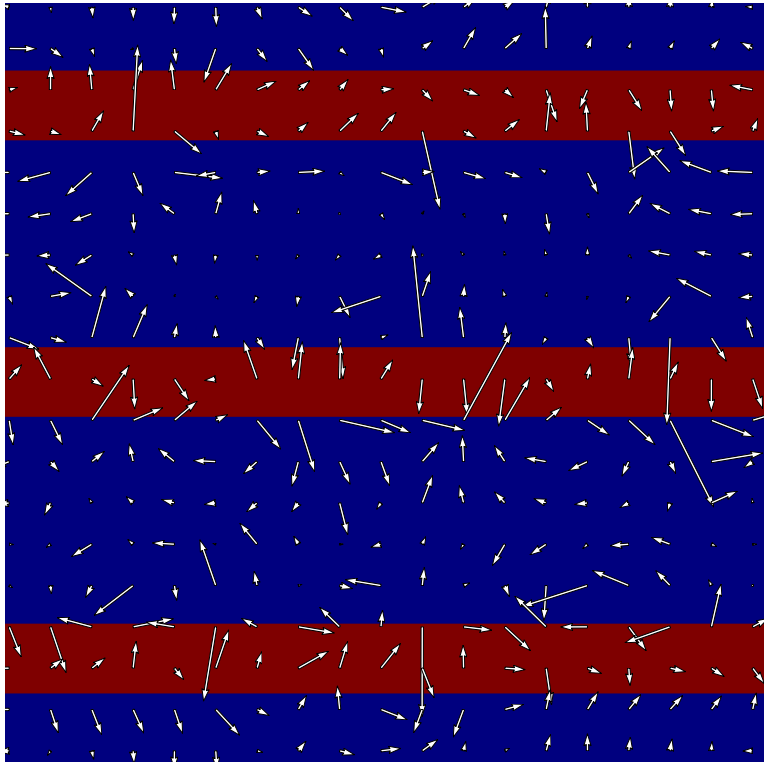

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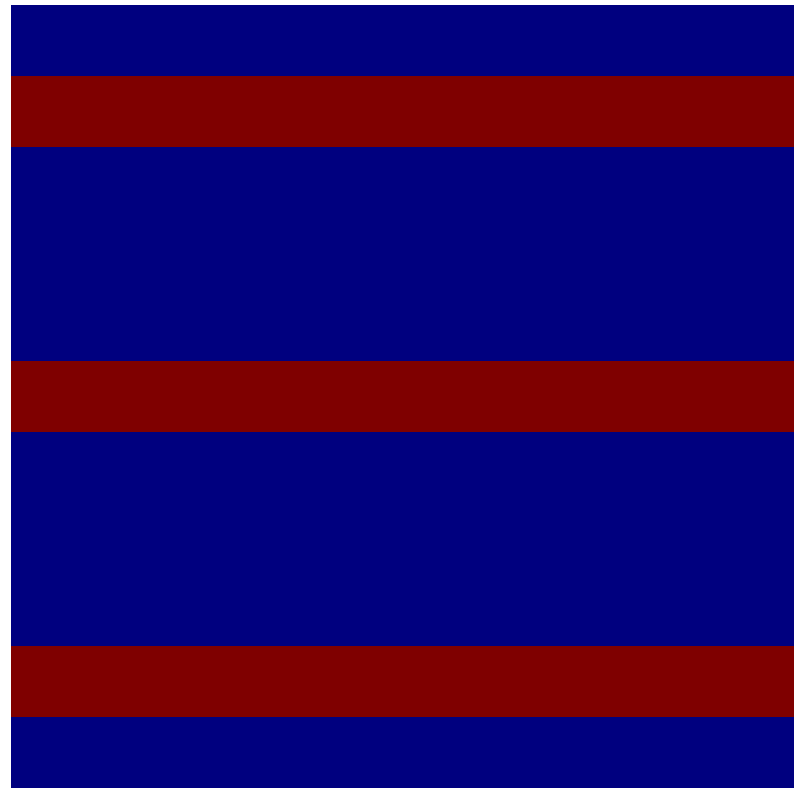
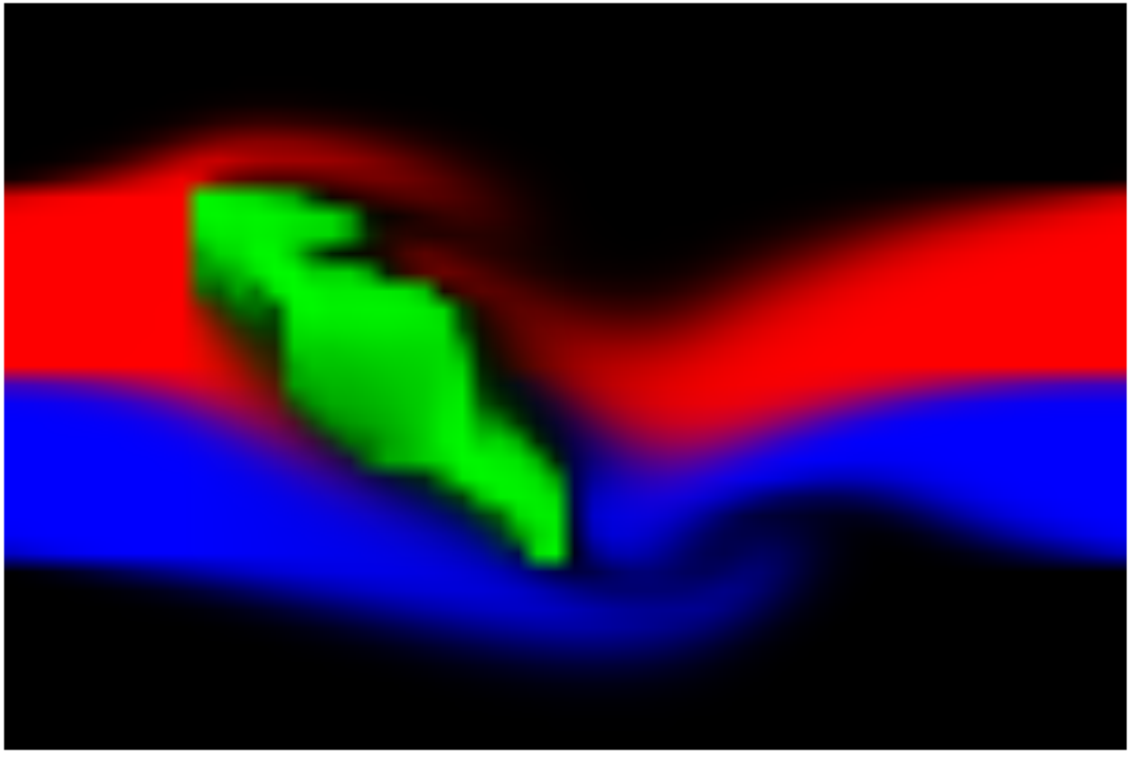
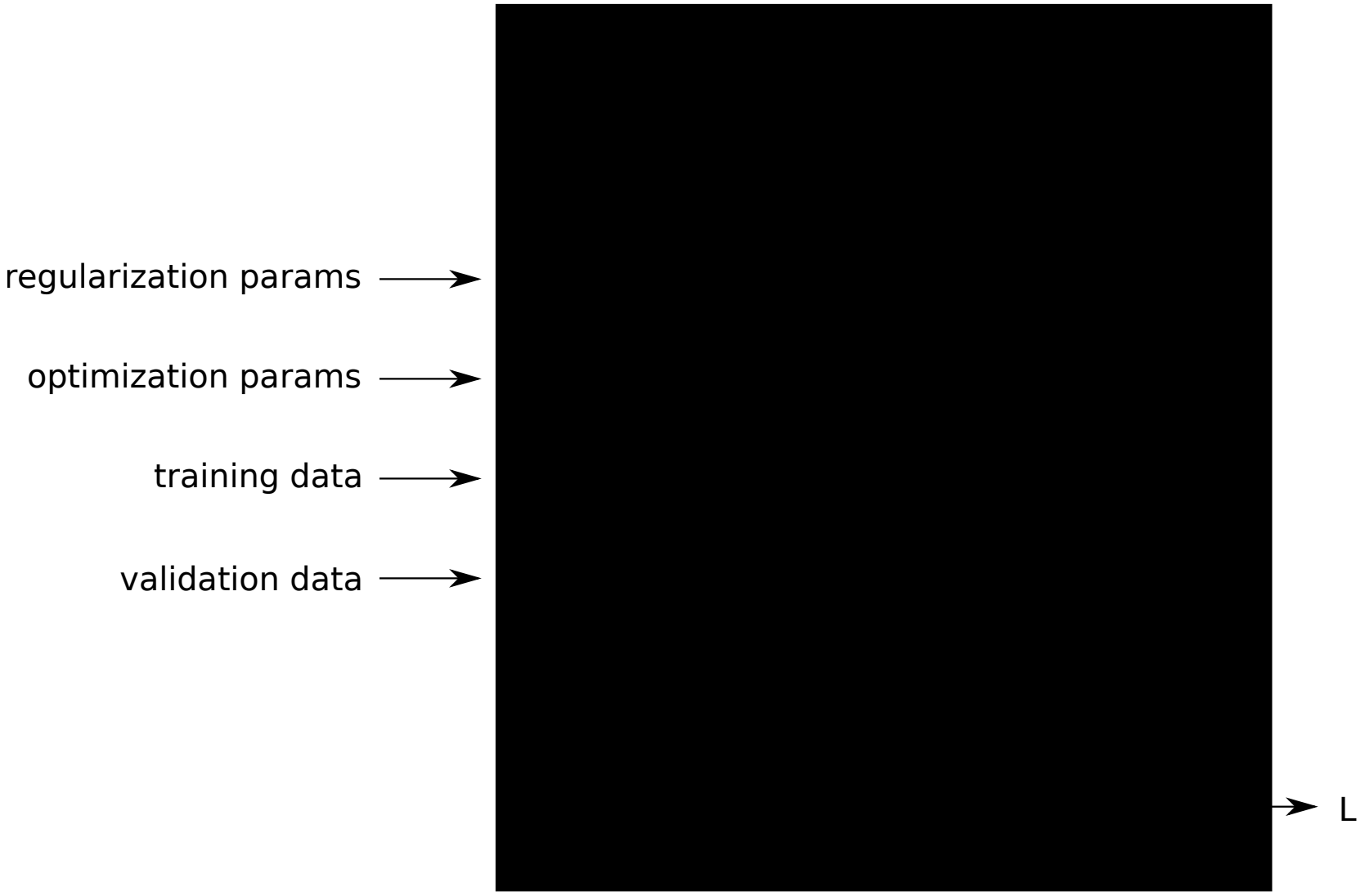


More fun with fluid simulations



Can optimize any objective!

Can we optimize optimization itself?



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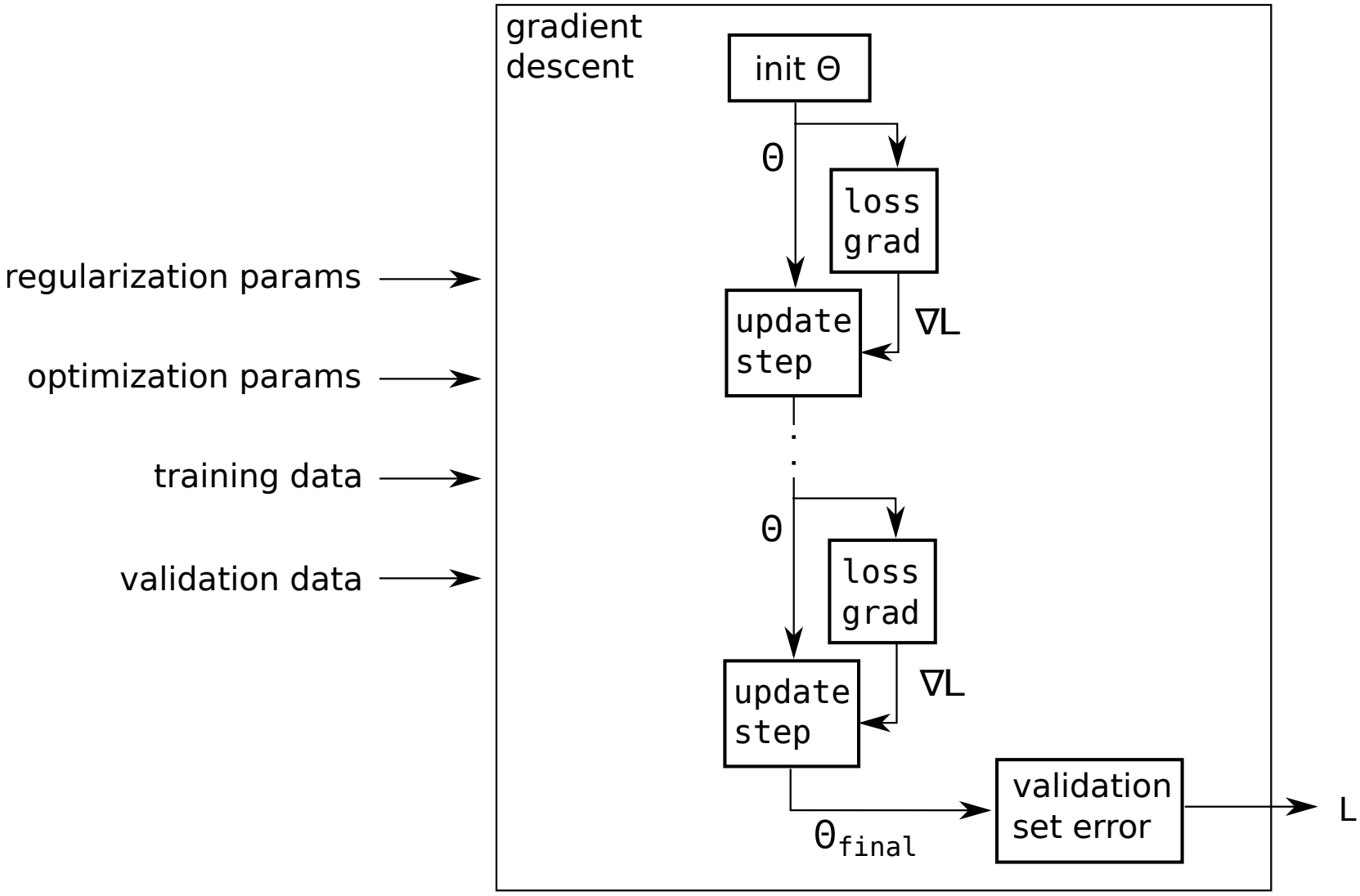
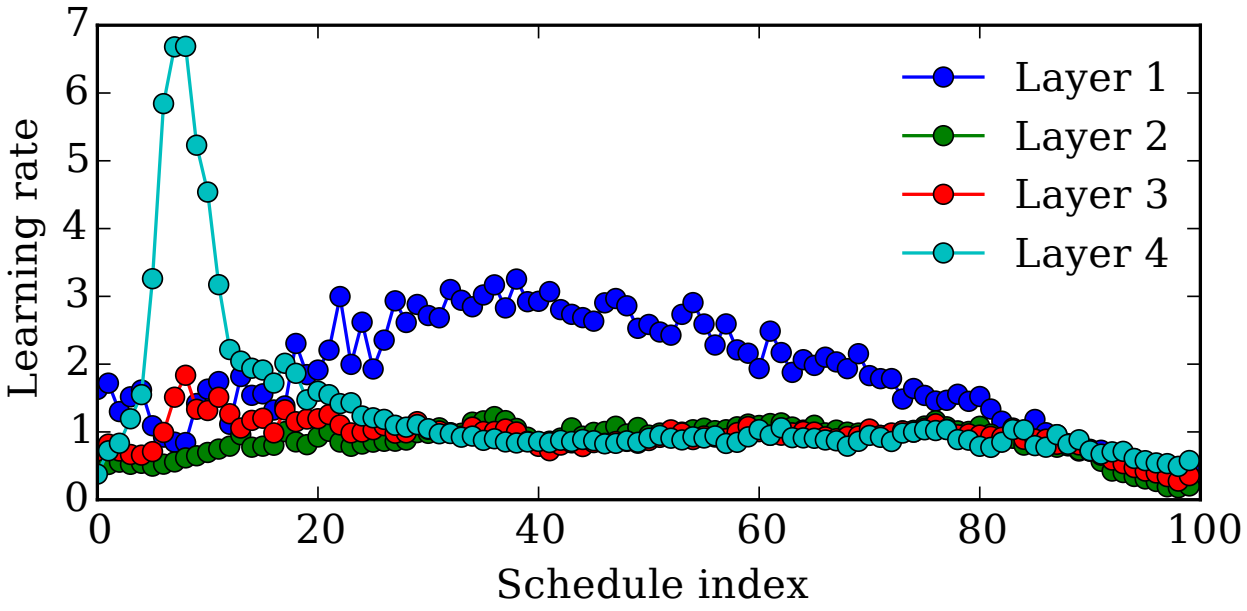
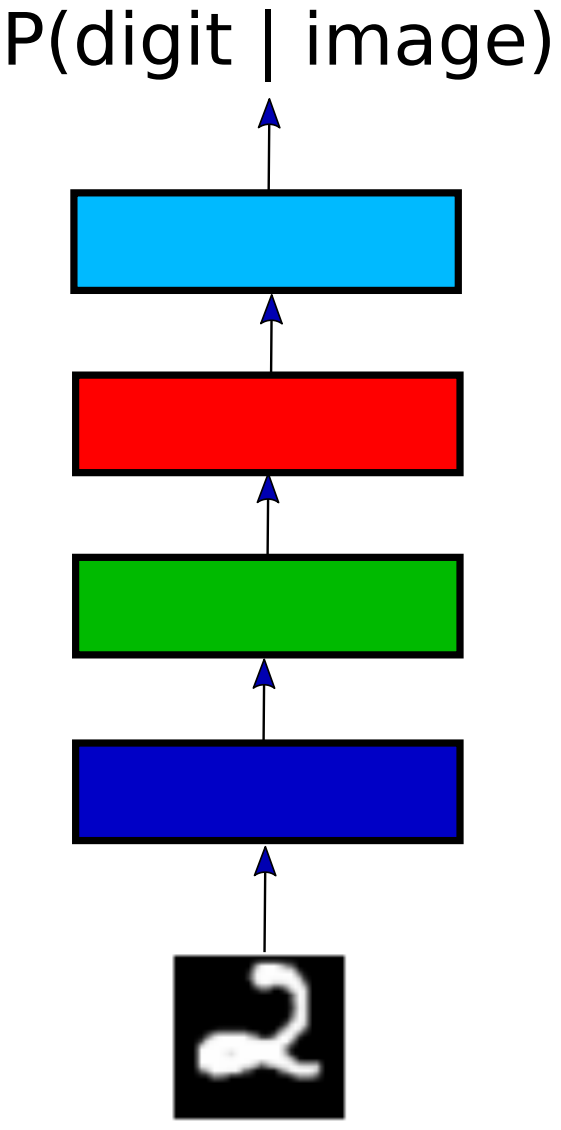


Optimized training schedules

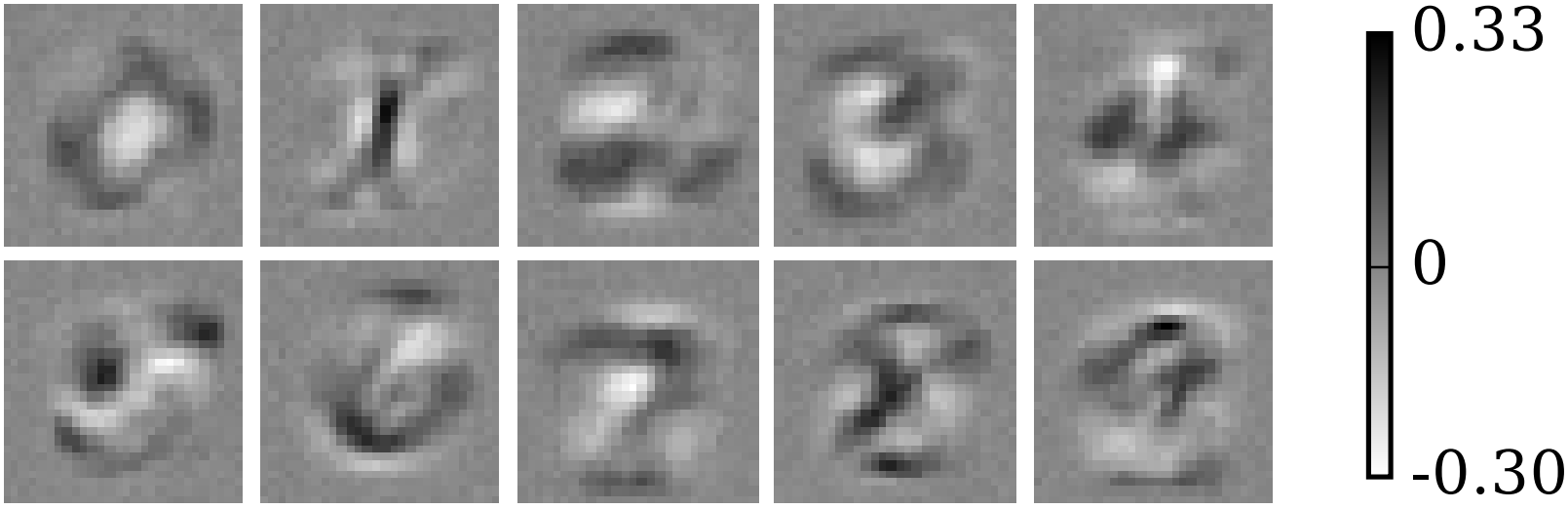




What else could we optimize?

Optimizing training data

- Training set of size 10 with fixed labels on MNIST
- Started from blank images

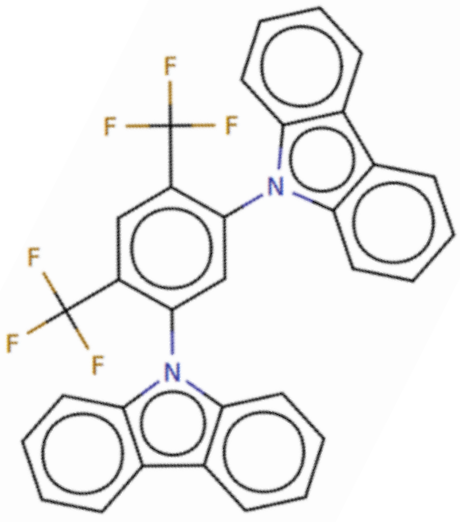


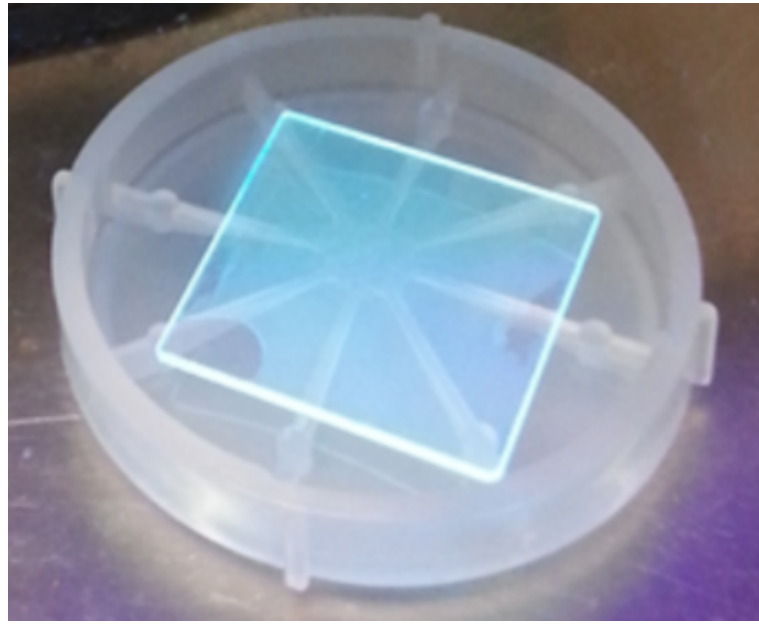
Maclaurin, Duvenaud & Adams, 2015

github.com/HIPS/hypergrad

Machine learning for virtual screening

Predicting properties of molecules from examples



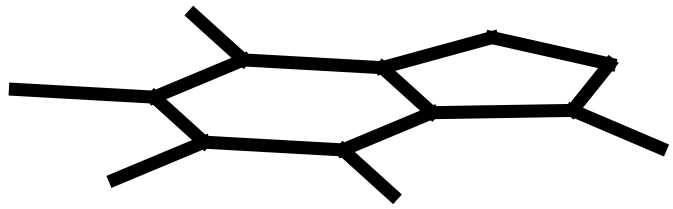


- Problem: molecular graphs can be any size and shape
- How to turn any graph into a fixed-size vector?

Convolution on graphs

Efficient approximation
(Glem *et al.*, 2006):

- Initialize nodes with atom type
- Combine each node with neighbors
- Use softmax to index



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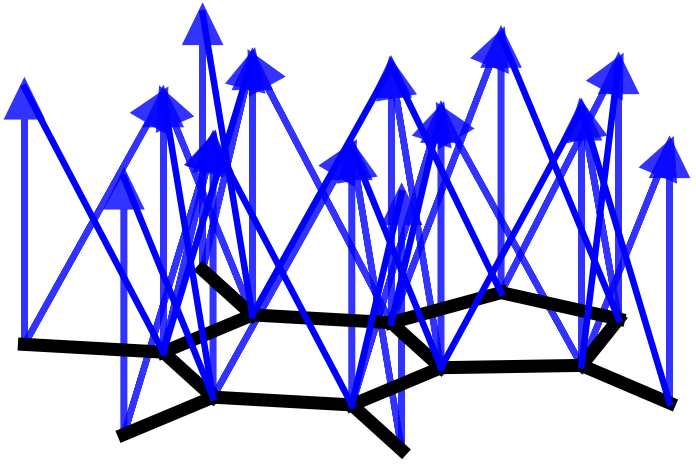

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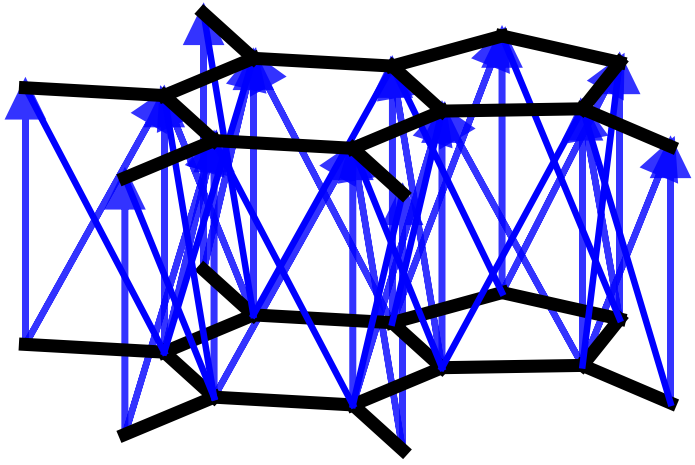

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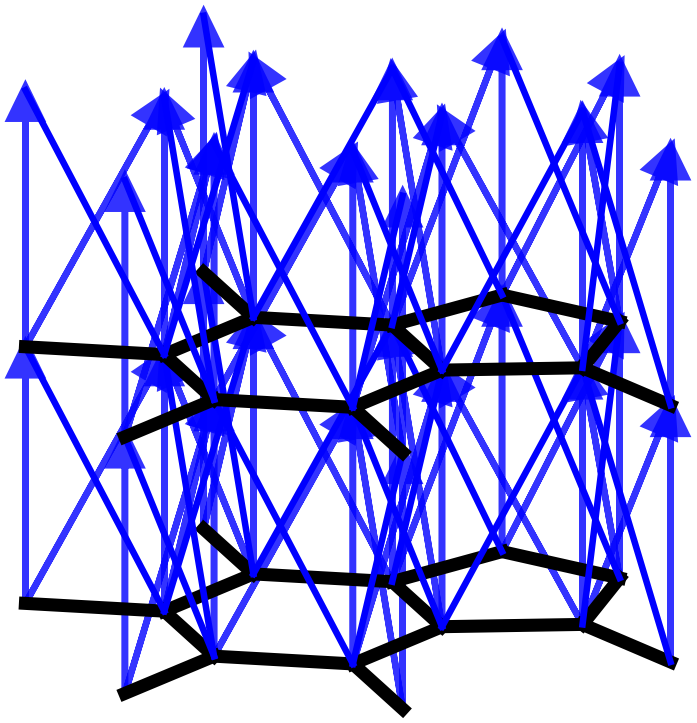

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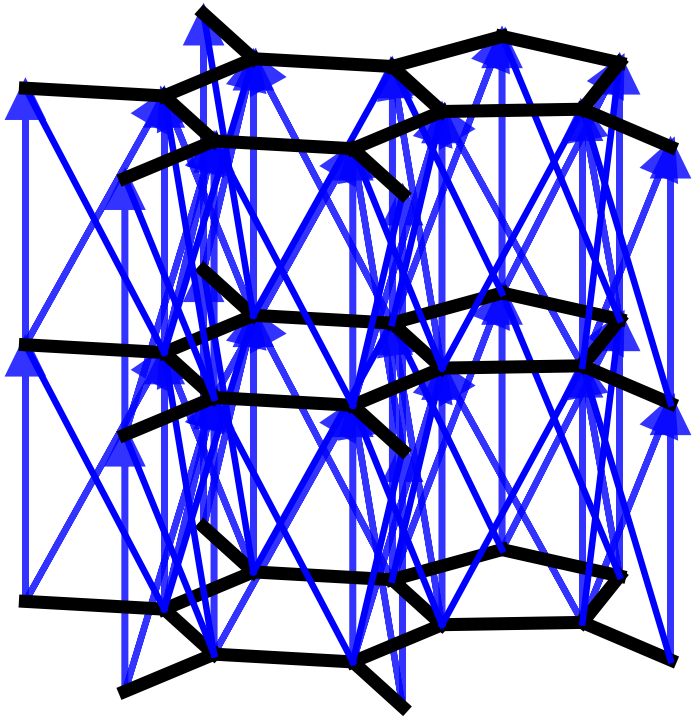

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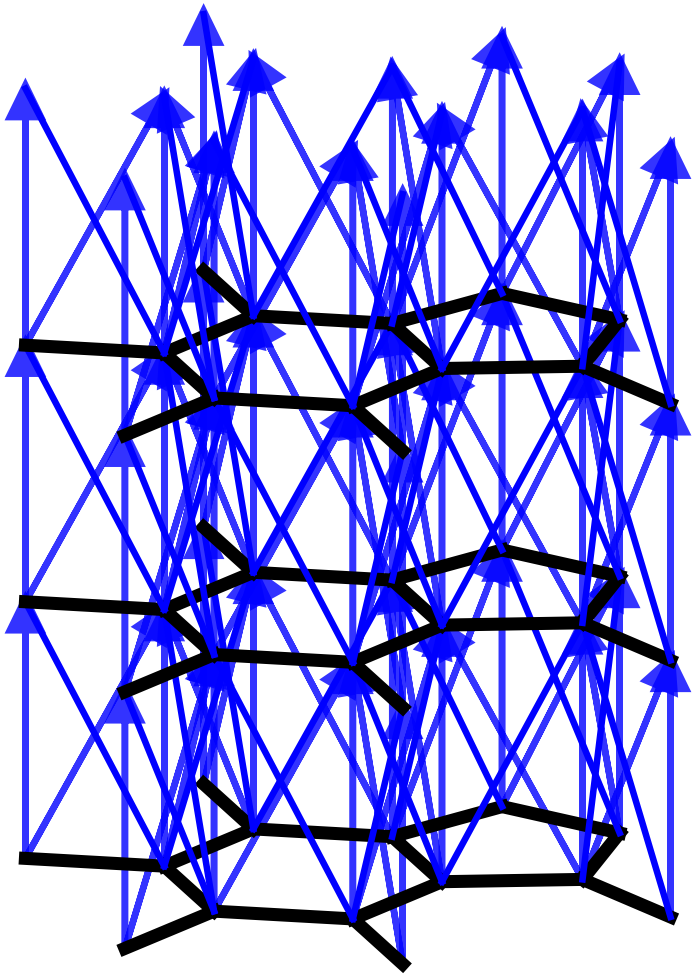

Convolution on graphs

Efficient approximation (Glem *et al.*, 2006):

- Initialize nodes with atom type
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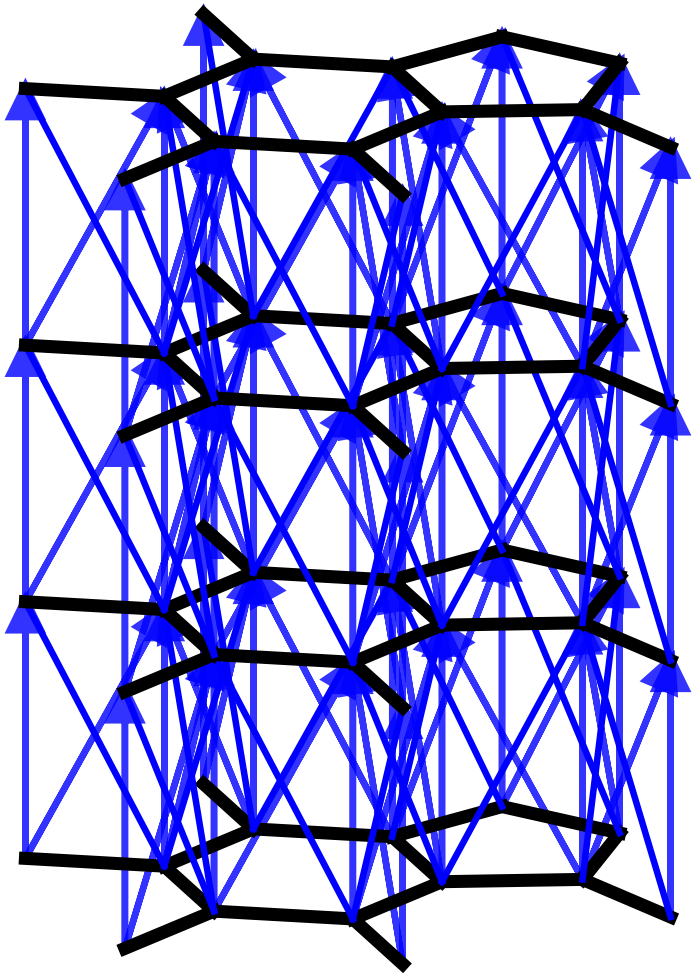
- Use softmax to index



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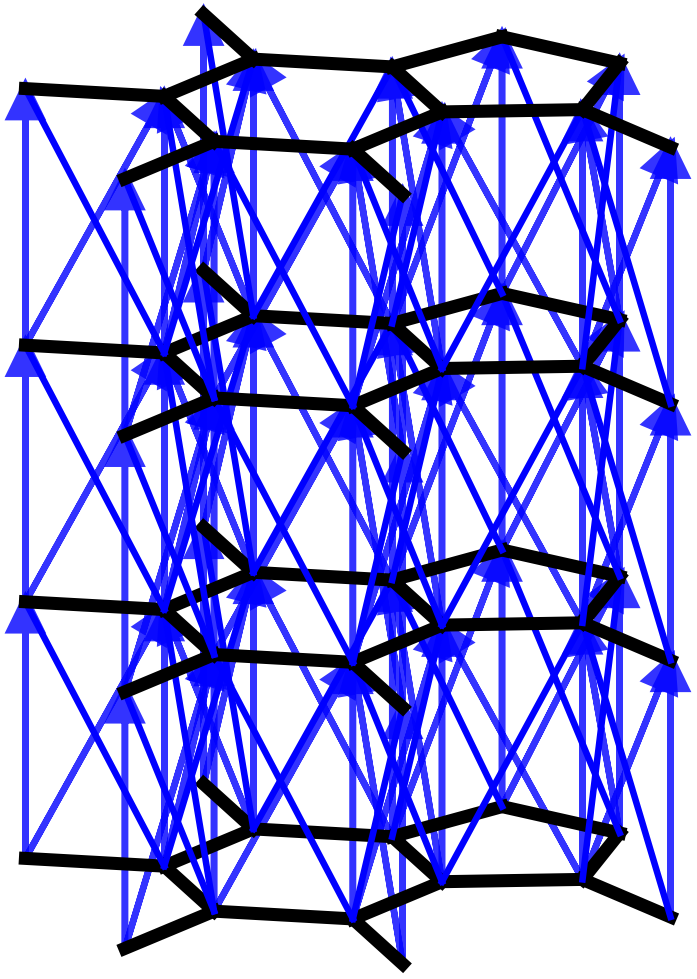

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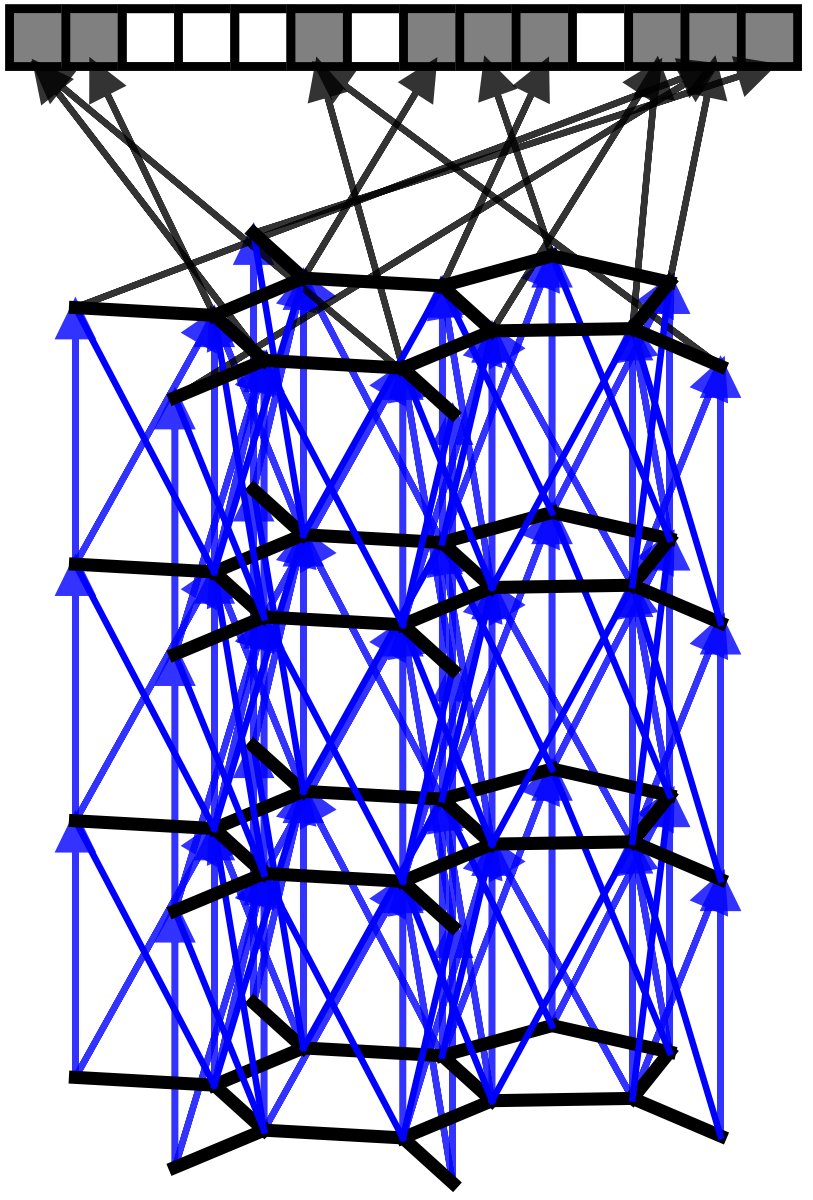
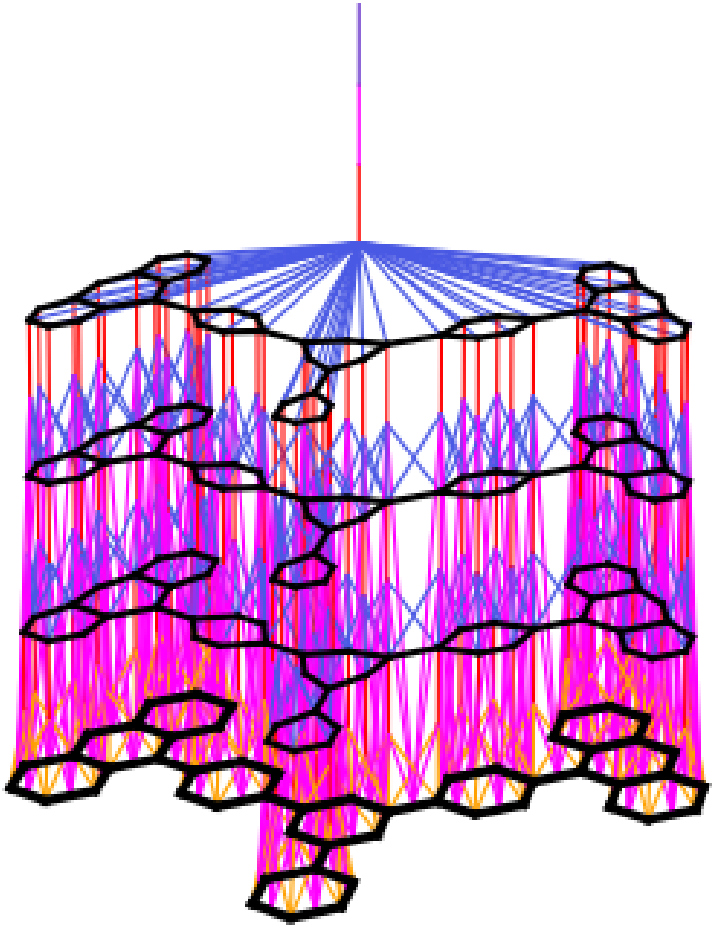
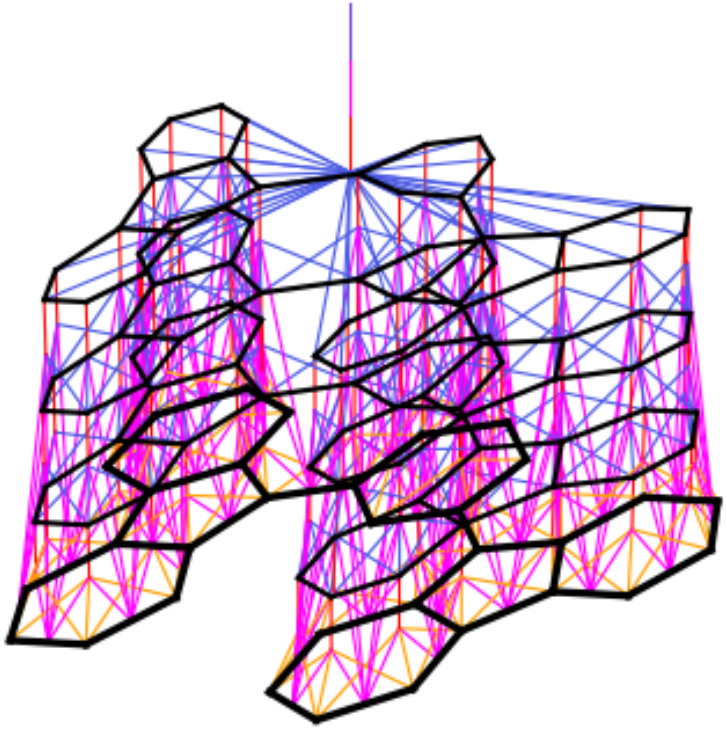


Each input yields a different computational flow





But what about inference?

Stan also provides inference routines...

But what about inference?

Stan also provides inference routines...

Ryan Adams @ryan_p_adams · 7 Nov 2015

@DavidDuvenaud

```
def elbo(p, lp, D, N):
 v=exp(p[D:])
 s=randn(N,D)*sqrt(v)+p[:D]
 return mvn.entropy(0, diag(v))+mean(lp(s))
gf = grad(elbo)
```

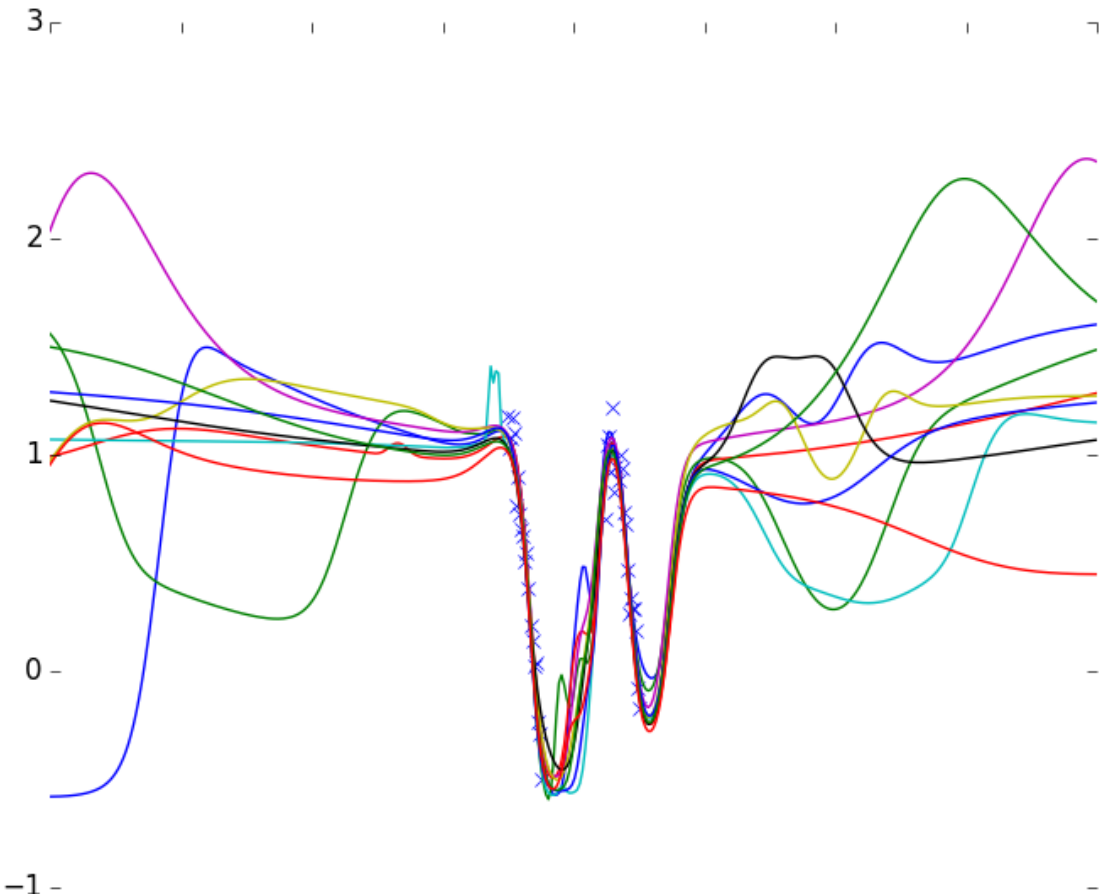
1 7 22 •••

... which are a tiny amount of code with autodiff!

Show Bayesian Neural Network

Our awesome new world

- You can autodiff directly in high-level languages
- Makes data-dependent graphs, higher-order ops, small code easy
- Frees our minds



Collaborators

github.com/HIPS/autograd

Dougal Maclaurin

Matthew Johnson

Ryan Adams

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Thanks!