Today's Discussion

To date:

- Neural networks: what are they
- Backpropagation: efficient gradient computation
- Advanced training: conjugate gradient

Today:

- CG postscript: scaled conjugate gradients
- Adaptive architectures
- My favorite neural network learning environment
- Some applications

Scaled conjugate gradient algorithm

Basic idea: Replace line minimization:

$$E(\mathbf{w}_j + \alpha^* \mathbf{d}_j) \le E(\mathbf{w}_j + \alpha \mathbf{d}_j), \ \forall \eta.$$

with:

$$\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j}$$

Why #!@\$ are we doing this? Didn't we want to avoid computation of H ?

Conjugate gradient algorithm

- 1. Choose an initial weight vector \mathbf{w}_1 and let $\mathbf{d}_1 = -\mathbf{g}_1$.
- 2. Perform a line minimization along \mathbf{d}_i , such that:

 $E(\mathbf{w}_{j} + \alpha^{*}\mathbf{d}_{j}) \leq E(\mathbf{w}_{j} + \alpha\mathbf{d}_{j}), \forall \eta.$

- 3. Let $\mathbf{w}_{j+1} = \mathbf{w}_j + \alpha^* \mathbf{d}_j$.
- 4. Evaluate \mathbf{g}_{j+1} .

5. Let
$$\mathbf{d}_{j+1} = -\mathbf{g}_{j+1} + \beta_j \mathbf{d}_j$$
 where,

$$\beta_j = \frac{\mathbf{g}_{j+1}^T (\mathbf{g}_{j+1} - \mathbf{g}_j)}{\mathbf{g}_j^T \mathbf{g}_j}$$
(Polak-Ribiere)

6. Let j = j + 1 and go to step 2.

Scaled conjugate gradient algorithm

Well, yes but:

- Line minimization can be computationally expensive.
- Don't really have to compute **H**? Huh?

$$\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j}$$

A closer look at α_i

Don't have to compute H, only Hd_i.

Theorem:

- \mathbf{w}_0 = current *W*-dimensional weight vector,
- $\mathbf{g}(\mathbf{w}) = \nabla E(\mathbf{w})$ (gradient of *E* at some vector \mathbf{w}), and,
- \mathbf{H} = Hessian of *E* evaluated at \mathbf{w}_0 ,
- \mathbf{d} = arbitrary *W*-dimensional vector.

$$\mathbf{H}\mathbf{d} = \lim_{\varepsilon \to 0} \frac{\mathbf{g}(\mathbf{w}_0 + \varepsilon \mathbf{d}) - \mathbf{g}(\mathbf{w}_0 - \varepsilon \mathbf{d})}{2\varepsilon}$$

Computing Hd_i

$$\frac{\mathbf{g}(\mathbf{w}_0 + \varepsilon \mathbf{d}) - \mathbf{g}(\mathbf{w}_0 - \varepsilon \mathbf{d})}{2\varepsilon} \approx \frac{2\varepsilon \mathbf{H} \mathbf{d}}{2\varepsilon}$$
$$\frac{\mathbf{g}(\mathbf{w}_0 + \varepsilon \mathbf{d}) - \mathbf{g}(\mathbf{w}_0 - \varepsilon \mathbf{d})}{2\varepsilon} \approx \mathbf{H} \mathbf{d}$$

So:

$$\mathbf{Hd} = \lim_{\varepsilon \to 0} \frac{\mathbf{g}(\mathbf{w}_0 + \varepsilon \mathbf{d}) - \mathbf{g}(\mathbf{w}_0 - \varepsilon \mathbf{d})}{2\varepsilon}$$

 $\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j}$ now just requires two gradient evaluations...

Computing Hd_{*i*}

$$\mathbf{Hd} = \lim_{\varepsilon \to 0} \frac{\mathbf{g}(\mathbf{w}_0 + \varepsilon \mathbf{d}) - \mathbf{g}(\mathbf{w}_0 - \varepsilon \mathbf{d})}{2\varepsilon}$$

First-order Taylor expansion of g(w) about w_0:

$$\mathbf{g}(\mathbf{w}) \approx \mathbf{g}(\mathbf{w}_0) + \mathbf{H}(\mathbf{w} - \mathbf{w}_0)$$

$$\frac{\mathbf{g}(\mathbf{w}_0 + \varepsilon \mathbf{d}) - \mathbf{g}(\mathbf{w}_0 - \varepsilon \mathbf{d})}{2\varepsilon} \approx \frac{[\mathbf{g}(\mathbf{w}_0) + \mathbf{H}(\varepsilon \mathbf{d})] - [\mathbf{g}(\mathbf{w}_0) - \mathbf{H}(\varepsilon \mathbf{d})]}{2\varepsilon}$$

New conjugate gradient algorithm

- 1. Choose an initial weight vector \mathbf{w}_1 and let $\mathbf{d}_1 = -\mathbf{g}_1$.
- 2. Compute α_i :

 $\alpha_{j} = -\mathbf{d}_{j}^{T}\mathbf{g}_{j}/\mathbf{d}_{j}^{T}\mathbf{H}\mathbf{d}_{j}, \forall \eta.$ 3. Let $\mathbf{w}_{j+1} = \mathbf{w}_{j} + \alpha_{j}\mathbf{d}_{j}.$ 4. Evaluate $\mathbf{g}_{j+1}.$ 5. Let $\mathbf{d}_{j+1} = -\mathbf{g}_{j+1} + \beta_{j}\mathbf{d}_{j}$ where, $\beta_{j} = \mathbf{g}_{j+1}^{T}(\mathbf{g}_{j+1} - \mathbf{g}_{j})/\mathbf{g}_{j}^{T}\mathbf{g}_{j}$ 6. Let j = j+1 and go to step 2. Any problems?

What about H < 0?

 $\alpha_j = -\mathbf{d}_j^T \mathbf{g}_j / \mathbf{d}_j^T \mathbf{H} \mathbf{d}_j$ might take uphill steps...

Idea:

- Replace **H** with $\mathbf{H} + \lambda \mathbf{I}$
- So:

$$\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j + \lambda \|\mathbf{d}_j\|^2}$$

What the #\$@! is this?

Model trust regions

Question: When should we "trust"

$$\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j}$$
?

Examining λ

$$\boldsymbol{\alpha}_{j} = \frac{-\mathbf{d}_{j}^{T}\mathbf{g}_{j}}{\mathbf{d}_{j}^{T}\mathbf{H}\mathbf{d}_{j} + \lambda \|\mathbf{d}_{j}\|^{2}}$$

- What is the meaning of λ being very large?
- What is the meaning of λ being very small (i.e. zero)?

Model trust regions

Question: When should we "trust"

$$\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j}?$$

- **1. H** is positive definite (denominator > 0)
- 2. Local quadratic assumption is good

Near a mountain, not a valley

Look at denominator of:

$$\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j + \lambda \|\mathbf{d}_j\|^2}$$
$$\delta = \mathbf{d}_j^T \mathbf{H} \mathbf{d}_j + \lambda \|\mathbf{d}_j\|^2$$

If $\delta < 0$, *increase* λ to make denominator positive.

How to increase λ ?

How about:

$$\lambda' = 2\left(\lambda - \frac{\delta}{\|\mathbf{d}_j\|^2}\right)$$

so that:

$$\delta' = \delta + (\lambda' - \lambda) \|\mathbf{d}_j\|^2$$
$$\delta' = \delta + \left[2\left(\lambda - \frac{\delta}{\|\mathbf{d}_j\|^2}\right) - \lambda\right] \|\mathbf{d}_j\|^2$$
$$\delta' = \delta - 2\delta + \lambda \|\mathbf{d}_j\|^2 = -\delta + \lambda \|\mathbf{d}_j\|^2$$

New effective denominator value

$$\lambda' = 2\left(\lambda - \frac{\delta}{\|\mathbf{d}_j\|^2}\right)$$
$$\delta' = -\delta + \lambda \|\mathbf{d}_j\|^2$$

So:

$$\delta' = -\left(\mathbf{d}_{j}^{T}\mathbf{H}\mathbf{d}_{j} + \lambda \|\mathbf{d}_{j}\|^{2}\right) + \lambda \|\mathbf{d}_{j}\|^{2}$$

 $\delta' = -\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j$ (what does this mean?)

Goin' up? I'll show you...

Since the new denominator is:

$$\delta' = -\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j$$

the new value of α_j is:

$$\alpha_{j}' = \frac{-\mathbf{d}_{j}^{T}\mathbf{g}_{j}}{-\mathbf{d}_{j}^{T}\mathbf{H}\mathbf{d}_{j}} = \frac{\mathbf{d}_{j}^{T}\mathbf{g}_{j}}{\mathbf{d}_{j}^{T}\mathbf{H}\mathbf{d}_{j}}$$

$$\alpha_{j}' = -\alpha_{j}$$

$$\mathbf{H} > 0$$

$$\alpha_{j}'$$

$$\mathbf{H} < 0$$

Model trust regions

Question: When should we "trust"

$$\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j}?$$

- **1.** H is positive definite (denominator > 0)
- 2. Local quadratic assumption is good

How to test local quadratic assumption?

Check:

$$\Delta = \frac{E(\mathbf{w}_j) - E(\mathbf{w}_j + \alpha_j \mathbf{d}_j)}{E(\mathbf{w}_j) - E_Q(\mathbf{w}_j + \alpha_j \mathbf{d}_j)}$$

What's E_Q ?

$$E_{Q}(\mathbf{w}) = E(\mathbf{w}_{0}) + (\mathbf{w} - \mathbf{w}_{0})^{T}\mathbf{b} + \frac{1}{2}(\mathbf{w} - \mathbf{w}_{0})^{T}H(\mathbf{w} - \mathbf{w}_{0})$$

So:

$$E_{Q}(\mathbf{w}_{j} + \alpha_{j}\mathbf{d}_{j}) = E(\mathbf{w}_{j}) + \alpha_{j}\mathbf{d}_{j}^{T}\mathbf{g}_{j} + \frac{1}{2}\alpha_{j}^{2}\mathbf{d}_{j}^{T}\mathbf{H}\mathbf{d}_{j}$$

What does Δ tell us?

Local quadratic test

$$\Delta = \frac{E(\mathbf{w}_j) - E(\mathbf{w}_j + \alpha_j \mathbf{d}_j)}{E(\mathbf{w}_j) - E_Q(\mathbf{w}_j + \alpha_j \mathbf{d}_j)}$$

Adjustment of trust region:

- If $\Delta > 0.75$ then decrease λ (e.g. $\lambda = \lambda/2$)
- If $\Delta < 0.25$ then increase λ (e.g. $\lambda = 4\lambda$)
- Otherwise, leave λ unchanged

Scaled conjugate gradient algorithm (α_{j}, λ)

- **1. Compute** $\delta = \mathbf{d}_j^T \mathbf{H} \mathbf{d}_j + \lambda \|\mathbf{d}_j\|^2$.
- 2. If $\delta < 0$, set $\lambda = 2(\lambda \delta / \|\mathbf{d}_j\|^2)$.
- 3. Compute $\alpha_j = -\mathbf{d}_j^T \mathbf{g}_j / (\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j + \lambda \|\mathbf{d}_j\|^2)$.
- **4.** Compute Δ :

5.
$$\Delta = \frac{E(\mathbf{w}_j) - E(\mathbf{w}_j + \alpha_j \mathbf{d}_j)}{E(\mathbf{w}_j) - E_Q(\mathbf{w}_j + \alpha_j \mathbf{d}_j)}$$

6. If
$$\Delta > 0.75$$
, set $\lambda = \lambda/2$, else if $\Delta < 0.25$, set $\lambda = 4\lambda$.

Scaled conjugate gradient algorithm

- 1. Choose an initial weight vector \mathbf{w}_1 and let $\mathbf{d}_1 = -\mathbf{g}_1$.
- 2. Compute α_i , λ :

$$\alpha_j = \frac{-\mathbf{d}_j^T \mathbf{g}_j}{\mathbf{d}_j^T \mathbf{H} \mathbf{d}_j + \lambda \|\mathbf{d}_j\|^2}, \ \forall \eta$$

- 3. Let $\mathbf{w}_{j+1} = \mathbf{w}_j + \alpha_j \mathbf{d}_j$.
- 4. Evaluate \mathbf{g}_{j+1} .
- 5. Let $\mathbf{d}_{j+1} = -\mathbf{g}_{j+1} + \beta_j \mathbf{d}_j$ where, $\beta_j = \mathbf{g}_{j+1}^T (\mathbf{g}_{j+1} - \mathbf{g}_j) / \mathbf{g}_j^T \mathbf{g}_j$ 6. Let j = j+1 and go to step 2.

Adaptive architectures

Standard learning:

- Select neural network architecture
- Train neural network
- If failure, go back to first step

Better approach:

• Adapt neural network architecture as function of training

Today's Discussion

To date:

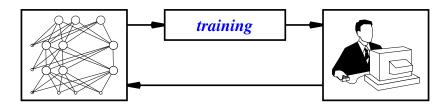
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Today:

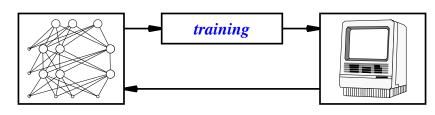
- CG postscript: scaled conjugate gradients
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- My favorite neural network learning environment
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Adaptive architectures

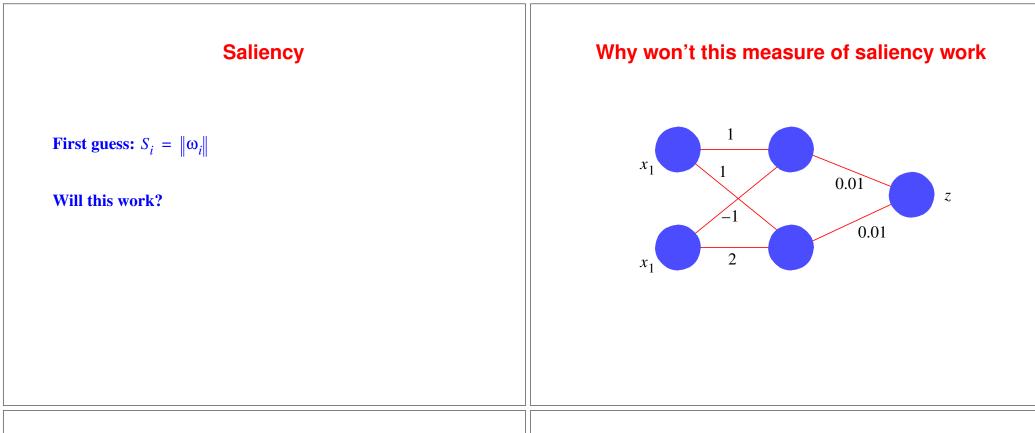
Standard learning:



Adaptive approach:



Adaptive architectures Pruning algorithms Basic idea: Problem: How do we do this? Start with really "big" network ٠ Eliminate "unimportant" weights/nodes Two main approaches: • Pruning (destructive algorithms) Retrain neural network • Growing (constructive algorithms) • **Advantages? Disadvantages? Problems? Pruning algorithms** Weight elimination schemes **Basic idea:** Idea: eliminate weights based on "saliency." Start with really "big" network • Eliminate "unimportant" weights/nodes **Definition:** saliency S_i = relative importance of weight ω_i • Retrain neural network **Any suggestions? Advantages?** (smaller final architectures) **Disadvantages?** (training cost of large network, retraining) **Problems?** (what is "unimportant?")



A better idea for saliency

Try to find relationship:

 $\delta E = \delta \mathbf{w}$

How can we do this?

• Brute force:

 $\delta E_i = \left\| E(\mathbf{w}) - E(\mathbf{w} + \delta \mathbf{w}_i) \right\|$

 $\delta \mathbf{w}_i = [0, ..., 0, -\omega_i, 0, ..., 0] (problems?)$

More on saliency

Use ol' reliable: 2nd order Taylor approximation

$$E(\mathbf{w}) = E(\mathbf{w}_0) + (\mathbf{w} - \mathbf{w}_0)^T \nabla E(\mathbf{w}_0) + \frac{1}{2} (\mathbf{w} - \mathbf{w}_0)^T H(\mathbf{w} - \mathbf{w}_0)$$

Now:

$$\delta \mathbf{w} = \mathbf{w}_1 - \mathbf{w}_0 \text{ (what are } \mathbf{w}_1 \text{ and } \mathbf{w}_0 \text{?)}$$

$$\delta E = E(\mathbf{w}_1) - E(\mathbf{w}_0) = \delta \mathbf{w}^T \nabla E(\mathbf{w}_0) + \frac{1}{2} \delta \mathbf{w}^T \mathbf{H} \delta \mathbf{w}$$

Can we simply this?

Optimal Brain Damage

$$\delta E = \frac{1}{2} \delta \mathbf{w}^T \mathbf{H} \delta \mathbf{w}$$

1. Idea: assume Hessian is diagonal

$$\delta E = \frac{1}{2} \sum_{i} H_{ii} \delta \omega_i^2$$

2. Resulting saliency:

$$S_i = \frac{H_{ii}\omega_i^2}{2}$$

- 3. Eliminate weights with smallest saliency
- 4. Retrain remaining weights

Optimal Brain Surgery

- Smarter idea: don't assume Hessian is diagonal
- Eliminate need for retraining

Now, assume you want to remove weight ω_i :

We want to minimize,

$$\delta E = \frac{1}{2} \delta \mathbf{w}^T \mathbf{H} \delta \mathbf{w}$$

subject to constraint

 $\delta \omega_i = -\omega_i (why?)$

Optimal Brain Surgery

Use Lagrange multipliers:

- We can minimize f(x) subject to constraint g(x) = 0 by minimizing L = f(x) + λg(x)
- $\lambda = \text{Lagrange multiplier}$

For our case:

$$f(\mathbf{x}) = \delta E = \frac{1}{2} \delta \mathbf{w}^T \mathbf{H} \delta \mathbf{w}$$

 $g(\mathbf{x}) = \delta \omega_i + \omega_i$

Optimal Brain Surgery

Minimize:

$$L = \frac{1}{2} \delta \mathbf{w}^T \mathbf{H} \delta \mathbf{w} + \lambda (\delta \omega_i + \omega_i)$$

•••

Solution:

$$\delta \mathbf{w} = -\frac{\omega_i}{[\mathbf{H}^{-1}]_{ii}} \mathbf{H}^{-1} \mathbf{u}_i$$

 $\delta E_i = \frac{1}{2} \frac{\omega_i^2}{[\mathbf{H}^{-1}]_{ii}}$ (what's the problem?)

Optimal Brain Surgery

1. Evaluate the inverse Hessian \mathbf{H}^{-1} .

2. Evaluate:

$$\delta E_i = \frac{1}{2} \frac{\omega_i^2}{[\mathbf{H}^{-1}]_{ii}}$$

- **3. Eliminate weight** ω_i , $\delta E_i < \delta E_j$, $i \neq j$.
- 4. Update all weights (no retraining)

$$\delta \mathbf{w} = -\frac{\omega_i}{[\mathbf{H}^{-1}]_{ii}} \mathbf{H}^{-1} \mathbf{u}_i$$

Pruning algorithms: key issues

- Large network to small network
- Need definition of saliency
- May need retraining step

Big problem: *lots of wasted training effort*

Node elimination scheme

Idea: Node pruning — need saliency of node, not weight

Define:

$$z_j = \gamma \left(\alpha_j \sum_i \omega_{ij} z_i \right)$$
 (output of unit j with addition of α_j)

Then:

$$s_j = E(\alpha_j = 1) - E(\alpha_j = 0)$$

 $s_j \approx \partial E / \partial \alpha_j |_{j=1}$

Growing algorithms

Basic idea:

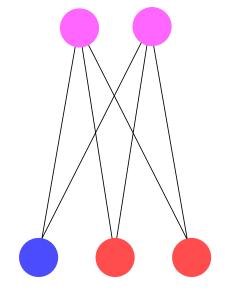
- Start with really small network
- Add hidden units as required

Advantages? Disadvantages? Problems?

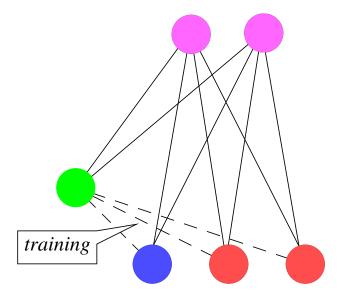
Growing algorithmsCascadeBasic idea:• Start with really small network• Add hidden units as required

Advantages? (reduced training cost, optimized networks) Disadvantages? (?) Problems? (arrangement of added weights/nodes)

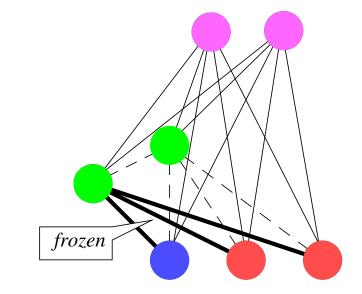
Cascade growing: initial network

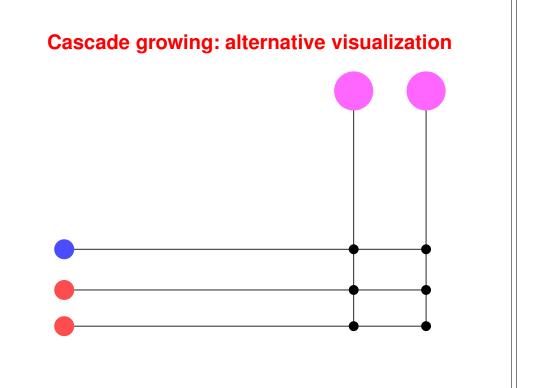


Cascade growing: first hidden unit

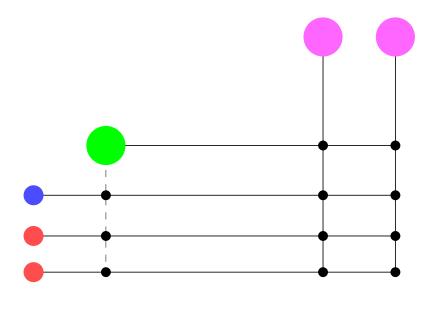


Cascade growing: second hidden unit

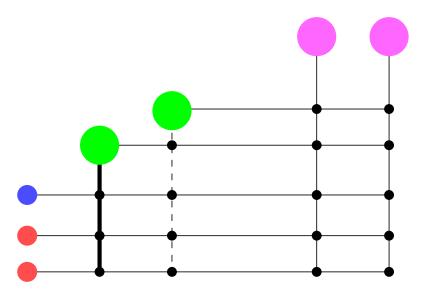




Cascade growing: alternative visualization

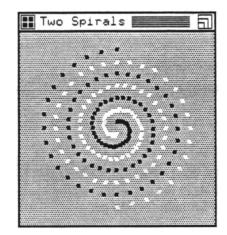


Cascade growing: alternative visualization

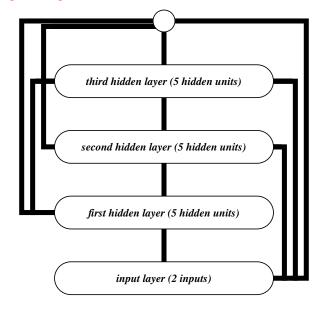


Cascade neural networks

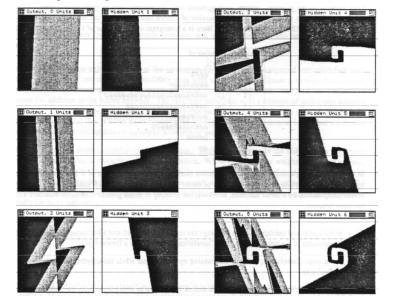
Do you ever need deeply nested structure?



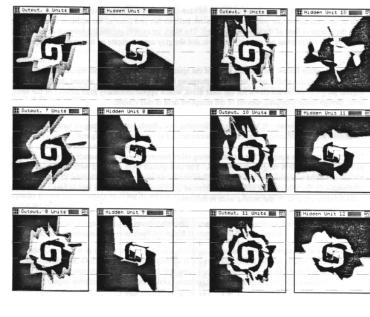
Two-spiral problem: best fixed architecture



Two-spiral problem: cascade architecture



Two-spiral problem: cascade architecture



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NN environment that rocks...

Two problems with traditional neural networks:

- *Fixed* architecture
 - Difficult to guess "appropriate" architecture
 - Functional complexity requirements can vary widely
- *Slow* learning algorithms (e.g. backprop, quickprop)

My neural network approach:

- *Flexible* architecture
 - Cascade neural networks
 - Variable activation functions
- *Fast* learning algorithm (e.g. NDEKF)

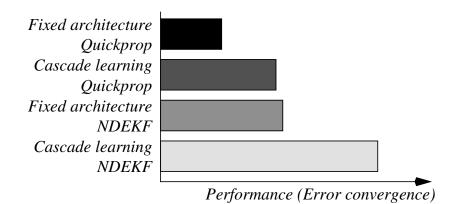
Cascade neural networks with nodedecoupled extended Kalman filtering (NDEKF)

Types of problems investigated:

- Continuous function approximation
- Dynamic system modeling

Cascade learning and NDEKF combine to result in better error convergence.

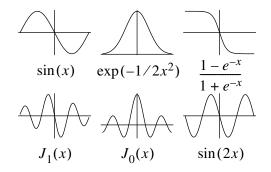
A sneak peak at results



Additional flexibility: variable activations

Cascade neural networks already offer great flexibility... However, why restrict candidate activation functions?

- Sigmoidal activation functions may not offer best results.
- Sinusoidals and/or others may be more appropriate:



Additional flexibility: variable activations

For continuous mapping problems:

- Variable networks converge to better minima.
- Sinusoidal networks about same as variable networks.

Better learning: extended Kalman filtering

View neural network training problem as system identification problem.

- Let weights of neural network represent *state* of nonlinear dynamic system.
- Let neural network be that nonlinear system.

Extended Kalman filter training:

- <u>Advantage</u>: Explicitly accounts for pairwise interdependence of weights with conditional error covariance matrix.
- <u>*Disadvantage*</u>: $O(W^2)$ computational complexity, where *W* is number of weights in network.

Decoupled extended Kalman filtering

Key insight:

- Some weights are more interdependent than others.
- Group weights into groups.
- Ignore interdependence between groups of weights (block diagonalize conditional error covariance matrix).

Even better idea: Group weights by node!

Node-decoupled extended Kalman filtering

Key insight: Decouple (group) weights by node: Natural formulation for cascade learning

- One weight group for current hidden unit
- One additional weight group for each output unit



- Matrix operations reduce to vector operations.
- Computational complexity reduces to $O\left(\sum W_i^2\right)$

Computational complexity					
NDEKF requires inversion of an $m \times m$ matrix, ($m =$ number of outputs)					
Cascade learning with NDEKF typically requires less than 10 epochs/hidden unit.					
• Several orders of magnitude less than backprop or quickprop approaches.					
 Computational complexity similar to fixed-architecture networks trained with NDEKF. 					
Experimental studies Four learning approaches:					
Symbol	Explanation				
Fq	fixed-architecture training with quickprop				
Cq	cascade-network training with quickprop				
Fk	fixed-architecture training with NDEKF				

cascade-network training with NDEKF

Ck

Ratio of computational cost between a cascade/NDEKF epoch and an equivalent fixed-architecture/backprop epoch (for few outputs):

Computational complexity

- Example: for 400 inputs and 20 hidden units ratio is less than 100.
- Example: for 20 or less inputs, ratio is less than 10.

Experimental studies

Key questions:

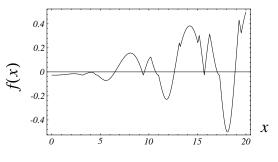
- Do we improve learning using NDEKF by going from fixed-architecture networks to cascade-type learning?
- Do we improve cascade learning by switching from quickprop (simple training) to NDEKF?
- Are any of more advanced methods (*Cq*, *Fk*, *Ck*) an improvement over baseline *Fq* (fixed-architecture/ quickprop) training method?

Five learning problems

Problem (A): smooth, continuous FA

 $f_1(x, y, z) = z\sin(\pi y) + x$ $f_2(x, y, z) = z^2 + \cos(\pi x y) - y^2$

Problem (B): nonsmooth, continuous FA



Learning results (avg. RMS error)

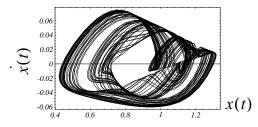
	Ck	Fk	Cq	Fq
(A)	42.1 (4.2)	127.1 (37.3)	94.5 (6.2)	N/A
(B)	7.4 (2.0)	12.4 (3.2)	14.5 (4.0)	65.0 (18.2)
(C)	15.6 (1.5)	20.7 (4.8)	29.9 (2.0)	N/A
(D)	4.6 (0.6)	10.2 (4.0)	9.4 (2.7)	16.7 (2.2)
(E)	42.0 (5.9)	60.5 (3.1)	72.6 (16.3)	90.3 (8.3)

Five learning problems

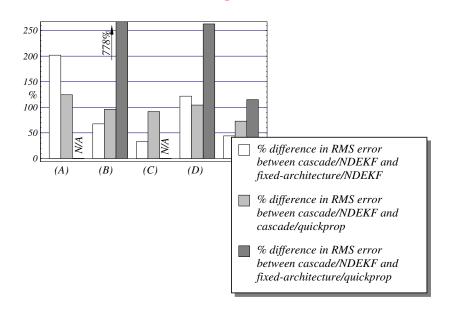
Problem (C): deterministic dynamic system

$$\begin{split} u(k+1) &= f[u(k), u(k-1), u(k-2), x(k), x(k-1)] \\ f[x_1, x_2, x_3, x_4, x_5] &= \frac{x_1 x_2 x_3 x_5 (x_3-1) + x_4}{1 + x_3^2 + x_2^2} \end{split}$$

Problems (D) & (E): *chaotic Mackey-Glass dynamic system* (*t*+6) and (*t*+84)

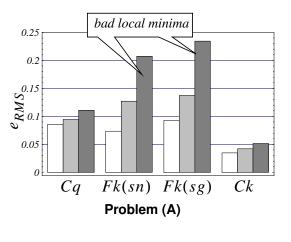


Learning results



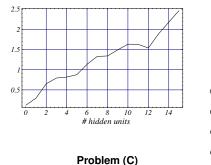
Why is Ck better than Fk?

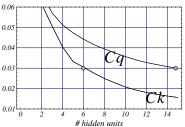
"NDEKF at times requires a small amount of redundancy in network in terms of total number of nodes in order to avoid poor local minima..." — [Puskorius & Feldkamp, 1991]



Why is *Ck* better than *Cq*?

As hidden units are added in cascade learning, NDEKF is better equipped to handle increasingly correlated weights to new hidden units.





Cascade/NDEKF advantages/disadvantages

- Cascade learning and NDEKF complement each other well.
- Cascade learning minimizes the potentially detrimental effect of node-decoupling.
- Cascade learning minimizes the problem of poor local minima in NDEKF.
- NDEKF better handles the increased correlation of weights as the number of hidden units increases in cascade learning.
- NDEKF requires no learning parameter tuning.
- Cascade/NDEKF converges efficiently to better local minima than either cascade or NDEKF by themselves.
- *Disadvantage:* computationally efficient with few outputs.

Today's Discussion

To date:

- Neural networks: what are they
- Backpropagation: efficient gradient computation
- Advanced training: conjugate gradient

Today:

- CG postscript: scaled conjugate gradients
- Adaptive architectures
- My favorite neural network learning environment
- Some applications