Data Mining II
Optimization & Parameter Tuning

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Why Parameter Tuning?

• What we have seen so far
  – many learning algorithms for classification, regression, ...

• Many of those have parameters
  – k and distance function for k nearest neighbors
  – splitting and pruning options in decision tree learning
  – hidden layers in neural networks
  – C, gamma, and kernel function for SVMs
  – ...

• But what is their effect?
  – hard to tell in general
  – rules of thumb are rare
Parameter Tuning – a Naive Approach

• You probably know that approach from the exercises
  1. run classification/regression algorithm
  2. look at the results (e.g., accuracy, RMSE, …)
  3. choose different parameter settings, go to 1

• Questions:
  • when to stop?
  • how to select the next parameter setting to test?
Parameter Tuning – Avoid Overfitting!

• Recap overfitting:
  – classifiers may overadapt to training data
  – the same holds for parameter settings

• Possible danger:
  – finding parameters that work well on the training set
  – but not on the test set

• Remedy:
  – train / test / validation split
Parameter Tuning – Avoid Overfitting!

- Parameter option: pruning (yes/no)
Parameter Tuning – Avoid Overfitting!

- Real example: train a local polynomial regression model
  - Parameter to tune: find the optimal maximum degree of the polynomial

- Tuning with proper validation: degree = 3
Parameter Tuning – Avoid Overfitting!

• Real example: train a local polynomial regression model
  – Parameter to tune: find the optimal maximum degree of the polynomial

• Tuning overfitting: degree = 9
Parameter Tuning: Brute Force

• Try all parameter combinations that exist

• Consider, e.g., a k-NN classifier
  – try 30 different distance measures
  – try all k from 1 to 1,000
  – use weighting or not
    → 60,000 runs of k-NN

→ we need a better strategy than brute force!
Parameter tuning is an optimization problem
Finding optimal values for N variables
Properties of the problem:
  - the underlying model is unknown
    • i.e., we do not know changing a variable will influence the results
  - we can tell how good a solution is when we see it
    • i.e., by running a classifier with the given parameter set
  - but looking at each solution is costly
    • e.g., 60,000 runs of k-NN

Such problems occur quite frequently
Intermezzo: Beyond Parameter Tuning

• Related problem:
  – feature subset selection
  – cf. Data Mining 2, first lecture

• Given n features, brute force requires $2^n$ evaluations
  – for 20 features, that is already one million
    $\rightarrow$ ten million with cross validation
Intermezzo: Beyond Parameter Tuning

- **Knapsack problem**
  - given a maximum weight you can carry
  - and a set of items with different weight and monetary value
  - pack those items that maximize the monetary value

- **Problem is NP hard**
  - i.e., deterministic algorithms require an exponential amount of time
  - Note: feature subset selection for N features requires $2^n$ evaluations
Intermezzo: Beyond Parameter Tuning

- Many optimization problems are NP hard
  - Routing problems (Traveling Salesman Problem)
  - Integer factorization
    - hard enough to be used for cryptography
  - Resource use optimization
    - e.g., minimizing cutoff waste
  - Chip design
    - minimizing chip sizes
Intermezzo: Beyond Parameter Tuning

http://xkcd.com/287/
Parameter Tuning: Brute Force

• Properties of Brute Force search
  – guaranteed to find the best parameter setting
  – too slow in most practical cases

• Grid Search
  – performs a brute force search
  – with equal-width steps on non-discrete numerical attributes (e.g., 10, 20, 30, .., 100)

• Parameters with a wide range (e.g., 0.0001 to 1,000,000)
  – with ten equal-width steps, the first step would be 1,000
  – but what if the optimum is around 0.1?
  – logarithmic steps may perform better
Parameter Tuning: Heuristics

• Properties of Brute Force search
  – guaranteed to find the best parameter setting
  – too slow in most practical cases

• Needed:
  – solutions that take less time/computation
  – and *often* find the best parameter setting
  – or find a *near-optimal* parameter setting
Beyond Brute Force

https://xkcd.com/399/
Parameter Tuning: One After Another

- Given $n$ parameters with $m$ degrees of freedom
  - brute force takes $m^n$ runs of the base classifier

- Simple tweak:
  1. start with default settings
  2. try all options for the first parameter
     2a. fix best setting for first parameter
  3. try all options for the second parameter
     3a. fix best setting for second parameter
  4. ...

- This reduces the runtime to $n \times m$
  - i.e., no longer exponential!
  - but we may miss the best solution
Parameter Tuning: Interaction Effects

• Interaction effects make parameter tuning hard
  – i.e., changing one parameter may change the optimal settings for another one

• Example: two parameters p and q, each with values 0,1, and 2
  – the table depicts classification accuracy

<table>
<thead>
<tr>
<th></th>
<th>p=0</th>
<th>p=1</th>
<th>p=2</th>
</tr>
</thead>
<tbody>
<tr>
<td>q=0</td>
<td>0.5</td>
<td>0.4</td>
<td>0.1</td>
</tr>
<tr>
<td>q=1</td>
<td>0.4</td>
<td>0.3</td>
<td>0.2</td>
</tr>
<tr>
<td>q=2</td>
<td>0.1</td>
<td>0.2</td>
<td>0.7</td>
</tr>
</tbody>
</table>
Parameter Tuning: Interaction Effects

- If we try to optimize one parameter by another (first \( p \), then \( q \))
  - we end at \( p=0, q=0 \) in six out of nine cases
  - on average, we investigate 2.3 solutions

<table>
<thead>
<tr>
<th></th>
<th>( p=0 )</th>
<th>( p=1 )</th>
<th>( p=2 )</th>
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<tbody>
<tr>
<td>( q=0 )</td>
<td>0.5</td>
<td>0.4</td>
<td>0.1</td>
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<tr>
<td>( q=1 )</td>
<td>0.4</td>
<td>0.3</td>
<td>0.2</td>
</tr>
<tr>
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<td>0.1</td>
<td>0.2</td>
<td>0.7</td>
</tr>
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Hill-Climbing Search

- a.k.a. greedy local search
- always search in the direction of the steepest ascend
  - "Like climbing Everest in thick fog with amnesia"

```plaintext
function HILL-CLIMBING(problem) returns a state that is a local maximum
inputs: problem, a problem
local variables: current, a node
               neighbor, a node

current ← MAKE-NODE(INITIAL-STATE[problem])
loop do
    neighbor ← a highest-valued successor of current
    if VALUE[neighbor] ≤ VALUE[current] then return STATE[current]
    current ← neighbor
```
Hill-Climbing Search

- Problem: depending on initial state, one can get stuck in local maxima
Hill Climbing Search

- Given our previous problem
  - we end up at the optimum in three out of nine cases
  - but the local optimum \((p=0, q=0)\) is reached in six out of nine cases!
  - on average, we investigate 2.1 solutions

<table>
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Variations of Hill Climbing Search

- **Stochastic hill climbing**
  - random selection among the uphill moves
  - the selection probability can vary with the steepness of the uphill move

- **First-choice hill climbing**
  - generating successors randomly until a better one is found, then pick that one

- **Random-restart hill climbing**
  - run hill climbing with different seeds
  - tries to avoid getting stuck in local maxima
Local Beam Search

- Keep track of k states rather than just one
- Start with k randomly generated states
- At each iteration, all the successors of all k states are generated
- Select the k best successors from the complete list and repeat
Simulated Annealing

- Escape local maxima by allowing “bad” moves
  - Idea: but gradually decrease their size and frequency
- Origin: metallurgical annealing

- Bouncing ball analogy:
  - Shaking hard (= high temperature)
  - Shaking less (= lower the temperature)
- If $T$ decreases slowly enough, best state is reached
Simulated Annealing

function SIMULATED-ANNEALING( problem, schedule) return a solution state
input: problem, a problem
        schedule, a mapping from time to temperature
local variables: current, a node.
                next, a node.
                T, a “temperature” controlling the probability of downward steps

current ← MAKE-NODE(INITIAL-STATE[problem])
for t ← 1 to ∞ do
    T ← schedule[t]
    if T = 0 then return current
    next ← a randomly selected successor of current
    ∆E ← VALUE[next] - VALUE[current]
    if ∆E > 0 then current ← next
    else current ← next only with probability $e^{\Delta E / T}$
Genetic Algorithms

• Inspired by *evolution*

• Overall idea:
  – use a population of individuals (solutions)
  – create new individuals by crossover
  – introduce random mutations
  – from each generation, keep only the best solutions ("survival of the fittest")

• Developed in the 1970s

• John H. Holland:
  – Standard Genetic Algorithm (SGA)

Charles Darwin (1809-1882)
Genetic Algorithms

• Basic ingredients:
  – individuals: the solutions
    • parameter tuning: a parameter setting
  – a fitness function
    • parameter tuning: performance of a parameter setting (i.e., run learner with those parameters)
  – a crossover method
    • parameter tuning: create a new setting from two others
  – a mutation method
    • parameter tuning: change one parameter
  – survivor selection
SGA Reproduction Cycle

1. Select parents for the mating pool
   (size of mating pool = population size)
2. Shuffle the mating pool
3. For each consecutive pair apply crossover with probability $p_c$, otherwise copy parents
4. For each offspring apply mutation
   (bit-flip with probability $p_m$ independently for each bit)
5. Replace the whole population with the resulting offspring
SGA Operators: 1-point crossover

- Choose a random point on the two parents
- Split parents at this crossover point
- Create children by exchanging tails
- $P_c$ typically in range $(0.6, 0.9)$
SGA Operators: Mutation

- Alter each gene independently with a probability $p_m$
- $p_m$ is called the mutation rate
  - Typically between $1/pop\_size$ and $1/chromosome\_length$

<table>
<thead>
<tr>
<th>parent</th>
<th>1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1</th>
</tr>
</thead>
<tbody>
<tr>
<td>child</td>
<td>0 1 0 0 1 0 1 1 0 0 0 1 0 1 1 0 0 1</td>
</tr>
</tbody>
</table>
SGA Operators: Selection

- Main idea: better individuals get higher chance
  - Chances proportional to fitness
  - Implementation: roulette wheel technique
    » Assign to each individual a part of the roulette wheel
    » Spin the wheel n times to select n individuals

<table>
<thead>
<tr>
<th></th>
<th>fitness(A)</th>
<th>fitness(B)</th>
<th>fitness(C)</th>
</tr>
</thead>
<tbody>
<tr>
<td>A</td>
<td>3</td>
<td></td>
<td></td>
</tr>
<tr>
<td>B</td>
<td>1</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>C</td>
<td>2</td>
<td></td>
<td>2</td>
</tr>
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</table>

3/6 = 50% 1/6 = 17% 2/6 = 33%
Crossover OR Mutation?

• Decade long debate: which one is better / necessary ...

• Answer (at least, rather wide agreement):
  – it depends on the problem, but
  – in general, it is good to have both
  – both have another role
  – mutation-only-EA is possible, crossover-only-EA would not work
Crossover OR Mutation? (cont’d)

- Exploration: Discovering promising areas in the search space, i.e. gaining information on the problem
- Exploitation: Optimising within a promising area, i.e. using information
- There is co-operation AND competition between them
  - Crossover is explorative, it makes a *big* jump to an area somewhere “in between” two (parent) areas
  - Mutation is exploitative, it creates random *small* diversions, thereby staying near (in the area of) the parent
Crossover OR Mutation? (cont'd)

- Recall the solution space example from Hill Climbing
  - crossover makes big jumps
  - mutation makes small steps
Crossover OR Mutation? (cont’d)

• Only crossover can combine information from two parents
• Only mutation can introduce new information (alleles)
• To hit the optimum you often need a ‘lucky’ mutation
Genetic Feature Subset Selection

- Feature Subset Selection
  - can also be solved by Genetic Programming

- Individuals: feature subsets
- Representation: binary
  - 1 = feature is included
  - 0 = feature is not included
- Fitness: classification performance
- Crossover: combine selections of two subsets
- Mutation: flip bits
Selecting a Learner

• So far, we have looked at finding good parameters for a learner
  – the learner was always fixed

• A similar problem is selecting a learner for the task at hand
• Again, we could go with search
• Another approach is meta learning
Selecting a Learner by Meta Learning

• Meta Learning
  – i.e., learning about learning

• Goal: learn how well a learner will perform on a given dataset
  – features: dataset characteristics, learning algorithm
  – prediction target: accuracy, RMSE, ...
Selecting a Learner by Meta Learning

• Used in the *Automatic System Construction* extension
• regression trained on
  – 90 datasets
  – 54 features

• Examples for features
  – number of instances/attributes
  – fraction of nominal/numerical attributes
  – min/max/average entropy of attributes
  – skewness of classes
  – ...

*Automatic System Construction* extension
Selecting a Learner by Meta Learning

• Used in the *Automatic System Construction* extension
...and now for something completely different.

- Recap: search heuristics are good for problems where...
  - finding an optimal solution is difficult
  - evaluating a solution candidate is easy
  - the search space of possible solutions is large
- Possible solution: genetic programming
- We have encountered such problems quite frequently
- Example: learning an optimal decision tree from data
Genetic Decision Tree Learning

• e.g., GAIT (Fu et al., 2003)
  – also the source of the pictures on the following slides
• Population: candidate decision trees
  – initialization: e.g., trained on small subsets of data
• Create new decision trees by means of
  – crossover
  – mutation
• Fitness function: e.g., accuracy
Genetic Decision Tree Learning

- Crossover:
Genetic Decision Tree Learning

- Mutation:

Subtree-to-subtree Mutation

Subtree-to-leaf Mutation
Genetic Decision Tree Learning

- Feasibility Check:
Combination of GP with other Learning Methods

• Rule Learning ("Learning Classifier Systems"), since late 70s
  – Population: set of rule sets (!)
  – Crossover: combining rules from two sets
  – Mutation: changing a rule

• Artificial Neural Networks
  – Easiest solution: fixed network layout
  – The network is then represented as an ordered set (vector) of weights e.g., \([0.8, 0.2, 0.5, 0.1, 0.1, 0.2]\)
  – Crossover and mutation are straight forward
  – Variant: AutoMLP
    • Searches for best combination of hidden layers and learning rate
Parameter Optimization vs. Pruning

- Architecture of a neural network can be seen as parameters
  - How many hidden layers? Which size?

- Pruning approaches: train large network, then start eliminating connections

Han et al. (2015): Learning both Weights and Connections for Efficient Neural Network
Wrap-Up

• Parameter tuning is important
  – many learning methods work poorly with standard parameters
  – often no global optimum, dataset dependent

• Parameter tuning has a large search space
  – trying all combinations is infeasible
  – interaction effects do not allow for one-by-one tuning
Wrap-Up

• Heuristic Methods
  – Hill climbing with variations
  – Beam search
  – Simulated Annealing
  – Genetic Programming

• Other uses of genetic programming
  – Feature subset selection
  – Model fitting
Next Steps DMC

• Register your team now
  – by sending an e-mail to Nico

• Task is out on Thursday
  – have a look as soon as possible
  – familiarize yourselves with the task and data

• We will have a first round of presentations **next week**
  – details will be announced

• Also next week
  – decision on evaluation metrics & protocol
Questions?