# Value Function Approximation

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## Dealing with imperfect value functions

$$||V - TV||_{\infty} \le \epsilon \to ||V - V^*||_{\infty} \le \frac{\epsilon}{1 - \gamma}$$

- How reassuring is this?
- Does this worst case hold in practice?

### Fitted value iteration (model-based)

- Assume:
  - Very large state space can't represent the value function as a vector
  - Generic machine learning "fit" operator that fits a continuous function based upon a set of training points
- Fitted VI algorithm:
  - Randomly initialize approximate value function V<sub>0</sub>
  - i=0
  - Repeat until done\*
    - Sample states S=s1...sm
    - Fit  $V_{i+1}$  on  $TV_i(s^1)...TV_i(s^m)$ .  $\leftarrow$  T is the Bellman operator
    - i=i+1
- Shorthand: V<sub>i+1</sub>=fit(TV<sub>i</sub>)
- How do we define "done"?

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### How to compute TV(s) in approximate VI

- Challenges:
  - V is not a vector, but some other representation
  - TV involves an expectation over next states, next states which may not be in original sample set S, i.e. off-sample extrapolation is likely required
- If number of next states is large and/or no model is available
  - Sample next states too
  - · Evaluate expected next state value by Monte Carlo
    - Generate many next states for each state
    - Possible if model/simulator can be easily reset
  - Or, train function approximator on r+γV(s') for each sampled s' (let function approximator do the averaging for you)

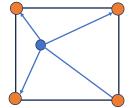
#### Properties of Fitted VI (FVI) – part I

- Properties of FVI depend upon properties of Fit function
- Recall that Bellman operator "T" is a contraction in max norm, i.e.,  $||V_1 V_2||_{\infty} < \epsilon \rightarrow ||TV_1 TV_2||_{\infty} < \gamma \epsilon$ ,  $0 \le \gamma < 1$
- If two operators, F and G are contractions (i.e. for any value function FV and GV are contractions) then F(GV) is a contraction
- Non-expansion: If H is a non-expansion in max norm, then:  $||V_1-V_2||_{\infty}<\epsilon \rightarrow ||HV_1-HV_2||_{\infty}\leq \gamma\epsilon$
- If one of F or G is a non-expansion in max norm, and the other is a contraction, the F(GV) is a contraction

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### Properties of Fitted VI (FVI) - part II

- It follows from previous slide that if Fit is a non-expansion in max norm, then fitted VI is a contraction in max norm
- What choices of Fit are non-expansions?
- Most common examples are averagers, e.g., interpolation
- Fitted VI with interpolation:
  - Pick S=s<sup>1</sup>...s<sup>m</sup> to be a grid of points
  - Implementing Fit:
    - For points in S, store TV(s) exactly
    - For points outside of S, use a distance weighted average of nearest neighbors



## Properties of Fitted VI with averagers

- It converges!
- But to what?
- Suppose  $\varepsilon$  = largest approximation error introduced at any iteration
- Total error is bounded by  $\varepsilon/(1-\gamma)$

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## Is this good news?

- Good news:
  - Convergence yay!
  - In some cases it may be possible to estimate  $\epsilon$
- Bad news:
  - Averagers do not scale well
  - Keeping ε small requires dense S
  - Achieving dense S is exponentially expensive in dimension of space

#### **Beyond Averagers**

- Moving beyond averagers requires more powerful function approximation
- Linear approximation is more powerful that averagers because it can extrapolate beyond points in S=s<sup>1</sup>...s<sup>m</sup> (observe that by design, any point not in s<sup>1</sup>...s<sup>m</sup> has value > min(V(s<sup>1</sup>)...V(s<sup>m</sup>)) and < max(V(s<sup>1</sup>)...V(s<sup>m</sup>))
- Non-linear approximation (e.g. neural networks) is more powerful than linear approximation

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#### Linear Value Function Approximation

- |S| typically quite large
- Pick linearly **independent** features  $\Phi = (\phi_1 ... \phi_k)$  (basis functions)
- Desire weights **w**=w<sub>1</sub>...w<sub>k</sub>, s.t.

$$V^*(s) \approx \hat{V}(s) = \sum_{i=1}^k w_i \phi_i(s)$$

 $\hat{V} = \Phi \mathbf{w}$ 

W is a kx1 column vector Ф is an mxk matrix (m is number of states sampled)

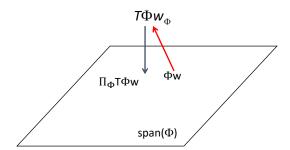
### Why is linear regression so important?

- Averagers interpolate (weak, resource hungry approximation)
- Regression Extrapolates (potentially more powerful)
- Linear regression = special case of most other methods
  - Neural networks
  - Kernel methods
- If regression fails, not much optimism on other methods

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#### Linear Fixed Point

•  $\Pi_{\Phi}$ V=projection of V into span( $\Phi$ )



• If we converge, we have:  $\Pi_{\Phi} T \Phi_{W} = \Phi_{W}$ 

### Example: Stability Problem [Tsitsiklis & Van Roy 1996]

**Problem:** Convergence not guaranteed



No rewards,  $\gamma = 0.9$ : V\* = 0

Consider linear approx. w/ single feature  $\phi$  with weight w.

$$\hat{V}(s) = w \cdot \phi(s)$$
 Optimal w = 0  
since V\*=0

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#### **Example: Stability Problem**



From V<sup>i</sup> perform projected iteration for each state

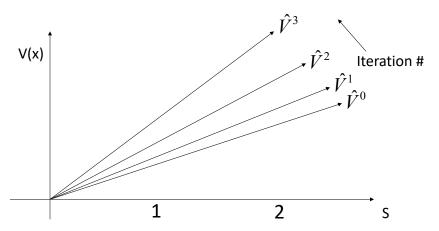
$$T[\hat{V}^i](s_1) = \gamma \hat{V}^i(s_2) = 1.8w^i$$
$$T[\hat{V}^i](s_2) = \gamma \hat{V}^i(s_2) = 1.8w^i$$

Can't be represented in our space so find wi+1 that gives least-squares approx. to exact backup

After some math we can get:  $\mathbf{w}^{i+1} = 1.2 \mathbf{w}^{i}$ 

What does this mean?

#### **Example: Stability Problem**



Each iteration of approximation makes things worse! Even for this simple problem fitted VI diverges.

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## Van Roy's Result

- Bellman operator *fixed policy* is a contraction in the weighted L<sub>2</sub> norm
- Weights come from the stationary distribution of P
- Linear regression in the weighted L<sub>2</sub> norm is non expansive in the weighted L<sub>2</sub> norm
- Understanding this:
  - Weighted norm redefines distance function so that different dimensions in the original space have different importance
  - Equivalent scaling the dimensions of the space
- Combined Regression-Bellman operator is a contraction!

To what does it converge?

$$\|V^{\pi} - \widehat{V}^{\pi}\|_{2,\rho} \le \frac{1}{\sqrt{1-\kappa^2}} \|V^{\pi} - \Pi V^{\pi}\|_{2,\rho}$$

- $\rho$  is the stationary distribution of  $\text{P}_\pi$
- $\kappa$  is the effective contraction rate (< $\gamma$ )

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#### Q-iteration: Generalization of Value Iteration

- $\forall s, a : Q(s, a) \leftarrow R(s, a) + \gamma \Sigma'_s P(s'|s, a) V(s')$
- $V(s') = \max_{a'} Q(s', a')$
- Q-iteration has similar convergence properties to value iteration

#### Application to stopping

- What about optimization?
- How to think about Bellman operator with max
  - Define T\*Q as the Q-iteration operator
  - $T_Q^*$  is a contraction is Max Norm
- Is T\*<sub>Q</sub> a non-expansion in weighted L<sub>2</sub>?
- No. 🙁
- But... It is non-expansion if max is always done with a constant
- Optimal stopping: Should I continue or stop and receive a payout?

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### Financial application

- Want to assign a price to an asset with following properties:
  - Can be held by owner for an arbitrary amount of time
  - Can cash out at some future time and receive a state-dependent reward
- Want to compute present value of this asset
- Features:
  - · Variables relevant to immediate value of asset
  - · Variables relevant to future value of the asset
- Supposedly used by some financial institutions to price assets

#### Perspective: Is weighted L<sub>2</sub> reasonable?

- In many ways more reasonable than Max norm
  - Worst case over entire state space hard to evaluate
  - Sampling methods can never provide guarantees without additional assumptions
- How do you achieve weighted L<sub>2</sub> in practice? (Sample from "real world" states)
- Weighted L<sub>2</sub> gives lower weight to less frequently occurring states
  - · Common cases get the most weight
  - Rare events may be wrong but that is forgivable(?)

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### Q-iteration in general

- What if Fit is a neural network?
- Linear value function approximation is a special case of this
- (Lack of) convergence guarantees from linear VFA apply to neural networks, but...
- If approximation error introduced at each step can be bounded by a constant, then overall approximation error is low
- Is this a reasonable assumption?

## Properties of approximate VI methods

- Convergence not guaranteed except in special cases
- Success has traditionally required very carefully chosen features and/or dense coverage to achieve low error
- Deep learning, which "automatically" learns feature representations, and uses massive amounts of samples has partially overcome this challenge