Notes on Recurrent Models of Visual Attention (2015)

We present a novel recurrent neural network model that is capable of extracting information from an image or video by adaptively selecting a sequence of regions or locations and only processing the selected regions at high resolution.

Glossary: -

A RNN that can be trained using reinforcement learning methods to learn task-specific policies.

Main aim is to reduce cost of computation when finding (images) or tracking (video) a object. Computational complexity is linear in the number of pixels. Compare this to the human visual system that "extracts" areas of "importance" and gives them more attention.¹ This reduces complexity of the task (where do I look?) and ignores noise (whats relevant for me?). Both number of parameters and amount of computation can be controlled independently of the size of input.

Litterature from neuroscience relies on saliency² and general attention.³ Should be expanded.

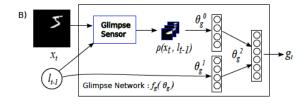
Authors frame it as a control problem by implementing a RNN that processes input (frames) sequentially by attending to different parts of the frame and incrementally combines this information to build a dynamic interal representation. The next location is chosen at time t based on past information $X_{\leftarrow t}$ and the policy. The procedure uses backpropagation to train the neural network components and policy gradient.

Previous work has focussed on sliding window paradigmes for cascades,^{4,5} branch and bound approach⁶ or proposition of candidate windows based on their likelihood to contain objects^{7,8} or

salience as mentioned above. These methods prioritise "interesting" regions of the image but fail to integrate information across fixations. Others have used a sequential decision tasks. 9–14

Essentially this is a POMDP. At each t the agent gets input $x_t = \{x_{\leftarrow t}, x_t\}$ through a bandwidth limited sensor \mathcal{M} (How is the sensor **implemented?**) that senses the (hidden) environmental state s_t (full image or current frame and past information). If $x_t = s_t$ the agent has "full information". The agent decides where to place the sensor (attention) and can change the true (hidden) state of the environment by taking action a_t . Since $\mathcal{M}: x_t \to s_t$ is only partial the agent needs to build a internal representation $p(s_t, x_t | x_{\leftarrow t})$ using past information effectively. At each t the agent gets a scalar reward r_t that depends on the actions "and can be delayed" (What is meant by delay?). The agent chooses the policy π_i that maximises $\sum r_t$.

 $\mathcal{M}(x_t, l_{t-1})$ where l_{t-1} is last location of sensor. Region around l is high-res but progressively gets lower res further from l. This is low the dimensionality of s_t is reduced to glimpse g_t . This is used in f_g (What is this subscript?) that produces a vector $g_t = f_g(x_t, l_{t-1}; \theta_g)$ (What does; mean here?) where $\theta_g = \{\theta_g^1, \theta_g^2, \theta_g^3\}$ (Is this raised to power 1 2 3?).



The internal (hidden state) $h_t = f_h(h_{t-1}, g_t; \theta_h)$ or core network summarises past information in h_{t-1} and must therefore be Markovian. The external input is g_t . At each t the agent performs two actions: Deploy sensor a_s and a environmental specific action a_t that depends on the task. Locations are chosen from a distribution parameterised by the location network $f_l(h_t; \theta_l)$ where $l_t \sim p(\cdot|f_l(h_t; \theta_l))$

and environment action $a_t \sim p(\cdot|f_a(h_t;\theta_a))$. For classification they use softmax and the exact formulations of the dynamic environment depends (as said) on the task (motor control, joystick, etc.). The model can be augmented to a cost-sensitive classisifier by adding negative reward for each additional glimpse. As said, the agent maxmises $\sum r_t$ or alternatively $\sum \gamma r_t$ where gamma is a discount factor. E.g. $r_T = 1$ if the object is classified correctly after T time $r_T = 0$ other ways.

Thus the agent needs to learn a stochastic policy $\pi((l_t, a_t)|s_{\leftarrow t}; \theta)$ where θ maps the history of past interactions with the environment $s_{\leftarrow t} = x_1, a_1, l_1, \dots, x_{t-1}, l_{t-1}, a_{t-1}, x_t$ to a distribution over actions for the current time step, subject to the constraint of the sensor. (What the f*** does that mean?)

Parameters of interest are given by the glimpse network, core network (figure below) and action network $\theta = \{\theta_g, \theta_h, \theta_a\}$ that are learned by maximising total (**Expected?**) reward. The policies of the agent in combination with environment gives a distribution over possible interaction sequences $\phi_{1:N}$. Maximise under

$$J(\theta) = \mathbb{E}_{p(\phi_{1:T};\theta)} \left[\sum_{t=1}^{T} r_t \right] = \mathbb{E}_{p(\phi_{1:T};\theta)} \left[R \right] \quad (1)$$

where $p(\phi_{1:T}; \theta)$ depends on policy.

Maximising $J(\cdot)$ is nontrivial as the expectation is over all possible high-dimensional interactions between policy and (unknown) environment. Applying gradient decent on a RL algorithm gives

$$\nabla_{\theta} J = \sum_{t=1}^{T} \mathbb{E}_{p(\phi_{1:T};\theta)} \left[\nabla_{\theta} log \pi(u_t | \phi_{1:T}; \theta) \cdot R \right]$$

$$\approx \frac{1}{M} \sum_{t=1}^{M} \sum_{i=1,t=1}^{M,T} \nabla_{\theta} log \pi(u_t^i | \phi_{1:T}^i; \theta) \cdot R^i$$
(2)

where each ϕ^i denotes a intertaction sequence obtained by running the current agent for i = 1...M episodes of T length. u_t denotes action at time t. For each run θ is adjusted to such that the log probability of actions that lead to high cumulative reward is increased, while

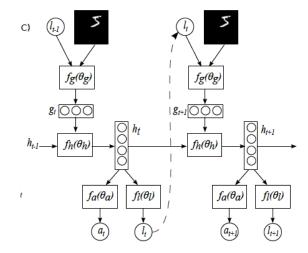
actions that produce low cumulative reward is reduced. $\nabla_{\theta} log \pi(u_t^i | \phi_{1:T}^i; \theta)$ is solved by back-propagation.

Equation 2 is an unbiased estimate of the gradient but may have high variance. It is common to solve this by considering

$$\frac{1}{M} \sum_{i=1,t=1}^{M,T} \nabla_{\theta} log \pi(u_t^i | \phi_{1:T}^i; \theta) \cdot (R^i - b_t)$$
 (3)

where R^i is total reward following action u_t^i and b_t is a baseline that depends on $\phi_{1:T}^i$ (via. h_t^i) but not on the action itself u_t^i (Kind of cryptic. How is this baseline calculated?). Taking expectations to equation 3 is equal to equation 2 but may have lower variance (Why?). Natural to select $b_t = \mathbb{E}(R_t)$ this baseline is simply a value function. Equation 3 results in an algorithm that ascribes higher log-probability to actions that were followed by larger-than-expected rewards (essentially prediction error) and vice versa for smaller rewards. The baseline is learned by reducing squared error between R_t^i and b_t .

If we have labeled data we can train on $log\pi(a_t^*|\phi_{1:T}^i;\theta)$ where a_t^* denotes labeled ground truth.



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