

Title of thesis

Second title line

Master Thesis of

Forename Surname

Department of Computer Science
Institute for Anthropomatics
and
FZI Research Center for Information Technology

Reviewer:	Prof. Dr.–Ing. J. M. Zöllner
Second reviewer:	Prof. Dr.–Ing. R. Dillmann
Advisor:	Dipl.–Inform. Max Mustermann

Research Period: XX. Monat 20XX – XX. Monat 2016

Title of thesis - Second title line

by
Forename Surname



Master thesis
in Monat 2016



Master thesis, FZI
Department of Computer Science, 2016
Gutachter: Prof. Dr.-Ing. J. M. Zöllner, Prof. Dr.-Ing. R. Dillmann

Affirmation

Ich versichere wahrheitsgemäß, die Arbeit selbstständig angefertigt, alle benutzten Hilfsmittel vollständig und genau angegeben und alles kenntlich gemacht zu haben, was aus Arbeiten anderer unverändert oder mit Abänderungen entnommen wurde.

Karlsruhe,
in Monat 2016

Forename Surname

Contents

1	Introduction	1
1.1	Motivation	1
1.2	Problem Statement	2
1.3	The Human Brain Project	2
1.4	Overview	2
2	Background	3
2.1	Probabilistic Graphical Models	3
2.1.1	Bayesian network	3
2.1.2	Markov Random Field	3
2.1.3	Energy-Based Models	4
2.1.4	Sampling	4
2.2	Neural Networks	5
2.2.1	Natural	5
2.2.2	Time discrete	7
2.2.3	Continuous	12
3	Related Work	15
3.1	Convolutional RBM	15
3.2	Sampling in SNNs	15
3.3	NN -> SNN conversion	16
3.4	eCD and Sampling Machines	16
4	Approach	19
4.1	Conversion	19
4.1.1	Conv DBNs	19
4.1.2	Conversion	19
4.2	eCD	21
4.2.1	Convolution	21
4.2.2	DBMs	22
5	Implementation	23
5.1	Discrete DBNs	23
5.2	Conversion	24
5.3	eCD	24

6 Experiments&Results	27
6.1 Datasets	27
6.1.1 1x4 Dataset	27
6.1.2 Strip Dataset	27
6.1.3 MNIST	27
6.2 Experiments	27
7 Conclusion and Outlook	29
7.1 Biological plausibility	29
A Some appendix	31
A.1 (if needed)	31
B List of Figures	35
C List of Tables	37
D Bibliography	39

1. Introduction

1.1. Motivation

In 2012 by winning the Imagenet Large Scale Visual Recognition Challenge 2012 convolutional neural network gained a big rise in popularity. Now they are becoming popular for their powerful abstraction mechanism in the fields of image and video classification and description and speech recognition. This can be contributed to compositional structure in which the world can be perceived and to their ability to extract high level features on spatial and/or temporal conditioned data.

Generating discriminative high level features extractors allows the system to dynamically adapt to the input data and work on various kinds of data. In addition the features extractors do not need to have any semantic representation and can be more complex to the manually build feature extractors. This also removes the labor-intensive and time consuming task of manually building feature extractors. Consequently they recently got adapted to solve robotic problems like grasp planning, drone navigation and autonomous driving.

One precursor of those are the deep belief networks (DBNs), which are built up of restricted boltzmann machines (RBMs). DBNs have shown excellent performance on image classification tasks in the early 2000s.

Compared to classical CNNs, DBNs allow recurrent connections and are trained in an unsupervised manner and do not need labeled data. They have been described as “probably the most biologically plausible learning algorithm for deep architectures we currently know”. DBNs can be used as generative model as well, which means they can sample data according to a learned distribution, e.g. find the most probable completion for a partially erased image.

Adding convolution to DBNs increased the performance of DBNs on image classification tasks, caught up to state of the art results and made the system more similar to the primate visual cortex than a standard RBM.

All those approaches use scalar values between neurons at discrete time slices to propagate information. This proposes some difficulties and is not biologically plausible, since biological neuron interleave linear and nonlinear operations, they communicate by stochastic binary values and are not synchronized. Spiking neural networks (SNNs), designed to simulate the communication between neurons with action potentials/ spikes, work in continuous time by design and do not suffer from the aforementioned limitations.

1.2. Problem Statement

To our best knowledge, up to today, there exists no system which utilizes the benefits of all those approaches. The main objective of this thesis is to realize such a spiking network, which integrates convolution and can be easily trained utilizing the RBM learning algorithm, to extract high level features. Two approaches are described. The first approach trains convolutional RBMs on discretized input data, to build up a DBN which is then converted to a SNN. The next approach directly works on continuous (event-driven) input-data and realizes a STDP learning algorithm with shared weights to train spiking convolutional RBMs directly. Both approaches will learn to extract high level features, which can be further used to classify an object or directly generate a grasp id.

1.3. The Human Brain Project

Heiko:

This thesis is under the scope of the research of the SP10 (sub-project 10) of the Human Brain Project. The Human Brain Project is an European Commission Future and Emerging Technologies Flagship and a large ten-year research project which aims to create a collaborative research infrastructure across national borders to progress the knowledge in neuroscience, computational neuroscience, medicine, scientific computation, and robotics. In the project over 120 institution from across Europe collaborate in 12 sub-projects.

The sub-project 10 of the Human Brain Project develops the Neurorobotics Platform which permits researchers to simulate robotic experiments with regard to neurorobotics. Apart from the platform, research is focused on the development of applications in robotics based on insights from neuroscience. One focus of the research group at the FZI Karlsruhe is the development of computational models for neurobiological inspired robotic grasping.

1.4. Overview

This thesis describes the approach and implementation of a spiking convolutional deep belief networks to extract high level features on continuous visual input. The thesis is structured as follows:

Chapter 2 introduces some background information, which is used in chapter 3 to describe state-of-the-art research used in this thesis. Chapter 4 will describe the different approaches to build such a spiking convolutional deep belief networks. In chapter 5 the different implementation steps and the architecture will be described. Chapter 6 outlines and compares the performance of the networks. Chapter 7 will conclude the gathered insight of this thesis, state its limitations and give suggestions for further improvements and research.

2. Background

2.1. Probabilistic Graphical Models

Neural networks can be seen as graphs

Using PGMs can be used to infer properties of some Neural Nets

PGMs are used to structure a model of the input and can be used for different task: Density estimation, Denoising, Missing value imputation, Sampling.

2.1.1. Bayesian network

Bayesian networks or belief networks are directed acyclic graphs, in which random variables are represented by nodes and their causal dependencies are represented by (directed) edges/connections.

If there is a connection from Node X_i to X_j then X_i is referred to as a parent of X_j and, similarly, X_j is referred to as the child of X_i .

In addition to the DAG structure, which is often considered as the “qualitative” part of the model, one needs to specify the “quantitative” parameters of the model. The parameters are described in a manner which is consistent with a Markovian property, where the conditional probability distribution (CPD) at each node depends only on its parents. FORMULARS

A BN reflects a simple conditional independence statement. Namely that each variable is independent of its nondescendants in the graph given the state of its parents. This property is used to reduce, sometimes significantly, the number of parameters that are required to characterize the JPD of the variables. This reduction provides an efficient way to compute the posterior probabilities given the evidence. Such a reduction provides great benefits from inference, learning (parameter estimation), and computational perspective.

GRAPH EXAMPLE

Given some observed variables, the bayes net can be used to infer the most likely states of the other hidden variables. This process of computing the posterior distribution of variables given evidence is called probabilistic inference. A Bayesian network can thus be considered a mechanism for automatically applying Bayes’ theorem to complex problems, which is defined through the “qualitative” and “quantitative” parameters of the net.

2.1.2. Markov Random Field

In contrast to Bayesian networks, Markov random fields are undirected graphical models, in which random variables are represented by nodes and edges/ connections indicate conditional dependencies.

For all cliques in the graph a "Clique Potential" / "Likelihood" can be given, which indicates how likely it is for the variables in the clique to take given values.

A clique in a graph is a group of nodes which are in themselves fully connected, i.e. each node has each other node of the graph as a direct neighbour.

The clique potentials can be used to calculate an unnormalized probability distribution $p(x)$

FORMULAR

EXAMPLE

2.1.3. Energy-Based Models

One way to model the unnormalized probability distribution $p(x)$ is to use an Energy function $E(x)$:
Formular

This results in an energy-based model (EBM).

Since $\exp(a) \cdot \exp(b) = \exp(a+b)$, where a and b are the Energy of a subgraph/clique and the Energy of the complete graph is the sum of all cliques, it is apparent that EBMs are a sub category of MRF.

Because $\exp(z) > 0$, there is probability of each state if greater than zero.

2.1.4. Sampling

Sampling is concerned with the selection of a subset of individuals from within a statistical population to estimate characteristics of the whole population.

Sampling provides a flexible way to approximate many sums and integrals at reduced cost, which could only have been calculated using costly operations or were intractable altogether.

Graphical models also facilitate the task of drawing samples from a model.

Ancestral Sampling For directed graphical models is that a simple and efficient procedure called ancestral sampling can produce a sample from the joint distribution represented by the model. The basic idea is to sort the variables x_i in the graph into a topological ordering, so that for all i and j , j is greater than i if x_i is a parent of x_j . The variables can then be sampled in this order.

The topological sorting operation guarantees that the conditional distributions are valid and one can sample from them in order.

Markov Chain Monte Carlo If $p_{\text{model}}(x)$ is represented by an undirected model, Markov Chain Monte Carlo methods can be used. MCMC methods interpret the model as a Markov chain, and work best, when no state gets zero probability assigned by the model.

The basic idea is to begin a state x with some arbitrary value. Then for a (infinite) time x is repeatedly randomly updated using the by the model given transition distribution $T(x, x)$. Eventually x becomes a fair sample from $p(x)$ / from the stationary distribution of the Markov chain.

To get more than one sample, one can run more Markov chains in parallel, each initialized with a random starting state. Another method is to run just one Markov chain, run it from some burn in/ mixing time, which allows the Markov chain to reach its equilibrium, and then take samples at

different timesteps. Those approaches need the Markov chain to reach its equilibrium distribution, which is usually done, by letting it run for some burn in time. But there is no guaranty, that the Markov chain has settled in the given timespan. Another problem with the second approach is, that since it can be hard to escape probable states, and when not run for an infinite timespan, more likely states can be over represented and less likely states under represented, if they did not occur or over represented if they did occur and the number of samples it not big enough to scale the estimate.

Gibbs sampling Gibbs sampling is a commonly used MCMC algorithm. The basic idea in Gibbs sampling to perform the transition from one state to another in accordance with $T(x,x)$ is, to select a single variable x_i and sample it conditioned on its neighbours. Several variables can be sampled at the same time as long as they are conditionally independent given all of their neighbours.

2.2. Neural Networks

In this section at first the neuroscience foundations of Neural Networks will be explained. After that different models will be described, starting with models working at discrete time steps and then describing models, which work in continuous time.

2.2.1. Natural

Brain

The human brain is the main organ of the human central nervous system.

It weighs about 1.2 - 1.4 kg (2% of the total body weight) and consumes around 20 Watts (20% of the total human power consumption).

Its mainly to contribute for tasks as learning, memory, self-control, planing, reasoning, abstract thought, motor control, vision and language.

All these tasks can be attributed to different specialized regions in the brain.

It is mainly composed of neurons and glial cells and blood vessels.

While the neurons perform the computations, the other constituents are required for structural stabilization and energy support.

The human brains contains 10^{12} neurons and each single neuron is interconnected with 10^4 other neurons.

The resulting neural network with roughly 10^{15} neural connections are the main component for human intelligence.

Neuron

A neuron is the main processing unit in the brain.

A Neuron has a potential difference between the interior cell and ,separated by the cell membrane, its surroundings. Thus the potential difference is called the membrane potential.

Ion channels in the cell membrane allow positive and negatively charged ions to flow from the inside to the outside and from the outside to the inside. This can lead to an increase (depolarization) or decrease (hyper-polarization) of the membrane potential.

The ion flow is primarily determined by the charge difference at the membrane (electro static force E / reversal potential), the ion concentration differences (diffusion force/nernst potential) and ion pumps actively pumping ions across the membrane.

The membrane potential of a neuron at rest, with an equal influx and outflow of ions neutralizing each other, has usually a resting/equilibrium potential of -65 mV.

Each single neuron can be divided into three functional distinct parts: the dendrites, the soma and the synapses.

The dendrites are thin, complexly branching structures emerging from the cell body/ soma. They are access for signals of preceding neurons and forward and accumulate these signals to the cell body. These signals can either polarize or depolarize the part of the dendrites.

The cell body or soma which encompasses the nucleus, accumulates all signals/ depolarizations from the dendrites and if the accumulation succeeds are threshold, usually around -55 mV, a action potential/spike emerges.

The spike is propagated across the axon and distributed to other subsequent (post-synaptic) neurons via the synapses, where they evoke (post-synaptic) potentials .

The axon is often covered by a fatty substance, the myelin, to regulate and improve the conductivity.

Synapses can be divided into two different categories: Electrical, which communicate with other neurons with electrical connections/synapses and chemical which chemical compounds.

Almost exclusively all synapses found in the brain are chemical.

Learning

Learning in the brain is associated with neural plasticity.

Neural plasticity describes the ability of the brain to change/reorganize the structure of brain, build new and alter connections.

Synaptic Plasticity describes the changes to the synaptic strength between single neurons.

Synaptic plasticity builds on the foundations, that neurons which are temporally and locally correlation, can lead to synaptic changes.

It can be seen as one of the important foundations of learning and memory.

It can be further divided in **short-term plasticity** which acts on a time scale one milliseconds to a minutes and **long-term plasticity**, lasting minutes or more.

Often the synapse strength change is contributed to a correlation between the firing time of the pre- and post-synaptic neurons.

One biological process LTP is based on, is Spike-timing-dependent plasticity (STDP), which is a abstract principle describing changes in the connection strength based on the relative timing of consecutive neurons.

"What fires together, wires together" A principle STDP is based on is the Hebbian theory which describes a simple mechanism for synaptic plasticity.

The Hebbian principle is often commonly summarized as "Cells that fire together, wire together".

Hebb originally stated it as follows: "When an axon of cell A is near enough to excite a cell B and repeatedly or persistently takes part in firing it, some growth process or metabolic change takes place in one or both cells such that A's efficiency, as one of the cells firing B, is increased."

Meaning, the more often A is active directly before B, the more likely A will have contributed to B's spike and the more causal B will become of A/ A and B become associated.

This can be mathematically generalized as: $\delta w = \theta_{x_i y}$

2.2.2. Time discrete

Artificial neural networks work/compute at discrete time slices, which can be interpreted as their tact. While it simplifies the neuron models a lot, it makes them exceptionally easy to handle on computers, which also work at a given tact rate.

Perceptron

The perceptron, also called Rosenblatt Perceptron, was invented in the late 1950s by Frank Rosenblatt. It was one of the first artificial neural networks, and can be seen as the foundation of most of the modern (deep) neural networks as well as linear discriminating classifiers.

Model The perceptron loosely models a neuron with a multi-dimensional input and a single output.

Be x the input and w the vector describing the synaptic weights. Each x_i is multiplied by it's synaptic weight w_i and then accumulated/summed up. $\sum x_i w_i = x \times w$.

If the sum exceed a threshold b the output y is set to 1 and 0 otherwise. $f = 1 \text{ if } w \times x + b > 0$.

Using the heaviside step function, the perceptron calculation rule can be rewritten as...

Decision Function This can be interpreted as a linear discriminating function, where w is a hyper plane in the data space diving data point which are classified as 1 or 0.

Perceptron Learning Given a set of datapoints X and their corresponding label Y and a learning rate θ .

For a x and y , their with eq... calculated output one update step can be described as

$$w = w + \delta w$$

, where

$$\delta w = \theta(d - y)x.$$

Thus the learning algorithm can be described as follows: 1. Initialize w randomly. 2. Select a data sample and calculate its output. 3. Calculate δw and update the weights.

This can be performed until all data samples are correctly classified i.e. $d = y$ or a predetermined number of iterations have been completed.

Multilayer-Perceptron

While the perceptron models a single neuron, the multi-layer perceptron can be seen as an extension modelling neural networks and by doing so overcoming the perceptrons disadvantage to only discriminate linearly.

Model architecture A MLP consists of multiple consecutive layers.

Each layer combines multiple perceptrons to map a multi-dimensional input to a multi-dimensional output.

A Layer is defined by: 1. The input dimension. 2. The number of perceptrons/ output dimension. 3. A Weight matrix W defining the weights between the synaptic connections 4. A activation function of the perceptrons.

The output of the layer is composed of the individual outputs of the perceptrons in the layer on the given input.

Thus be a perceptron the i th perceptron in the layer with the synaptic weights w_i its output d_i can be calculated using eq The output of the layer d is defined as $d = (d_1, \dots, d_n)$.

The weight matrix W can then be written as $W = (w_1, \dots, w_n)$.

This allows direct calculation rule for d : $d = f(Wx)$.

By using the output of a previous layer as input for the next layer, layers can be stacked up.

$$d_2 = f(W_2 d_1)$$

In this case there are only forward connections, so there are no cycles or loops (i.e. recurrent) connections in the network.

Activation functions For the choice of the activation function f there have been different functions, with different attributes proposed:

1. Step
2. Sigmoid
3. Softmax
3. Sign
4. Tanh
5. ReLU

Error functions To validate the quality of the model an error function/cost function is used. The cost function quantifies the performance of the model on a given task. The primary goal of a learning algorithm is to reduce the error on its task. Hereby it is important to note that the error function is often task dependent and rather independent on the learning algorithm used.

To compare the output o of the network to a label y of the input datum an error function or cost function is used $E(y, o)$.

Some of the more common are: 1. MSE 2. Cross entropy

Backpropagation The objective is to get weights/parameters which form (global) minimal point in the error function. One class of algorithms uses gradient descent to reduce the error and assign contributions to the weights and reach a minimal point.

In gradient descent for each parameter/weights its gradient is calculated, and by using the negative gradient direction each weight is updated.

Since MLPs have a nicely defined structure, gradient descent can be simplified using the chain rule to get an iterative procedure.

... gradient descent forumulars

Convolutinal Neural Networks

Convolutional neural networks (CNN) exploit spacial relations in the input, to regularize the complexity of the neural network by putting further constrains on the architecture of those nets, which make them easier to train and allow greater generalization with fewer training samples.

Instead of having fully connected nets, CNNs are only partially connected and use shared weights.

Convolution What is Conv and cross corelation

HBP Sem paper:

Convolution In the colvolution layer, the convolution operation with a learn- able kernel Matrix K is applied. An input given as an 3D tensor Y be composed of m 2D feature maps. Each feature map has the dimension $s \times t$. In the input layer, m is for example the number of color channels (3 in case of an RGB image), and s is the width and t the height of the input image. A discrete convolution with a (M, P, Q) filter matrix K_j at position (x, y) is then defined as :

Convolution In the colvolution layer, the convolution operation with a learn- able kernel Matrix K is applied. An input given as an 3D tensor Y be composed of m 2D feature maps. Each feature map has the dimension $s \times t$. In the input layer, m is for example the number of color channels (3 in case of an RGB image), and s is the width and t the height of the input image. A discrete convolution with a (M, P, Q) filter matrix K_j at position (x, y) is then defined as :

Architecture The most common architectures for convolutional nets for image classification are build up from stacks of alternating conv, pooling and normalization layers [17]. After those layers, fully connected layers are used as a classifier to assign labels to the extracted features (see Figure 7).

Training $\Delta W = \sum \Delta w$

Hopfield Nets

Another models of time discrete neural networks with recurrent, like Hopfield Nets and Boltzmann machnies use binary units and have their origin in/ are similar to energy-based/graphical models.
(?)

Model Hopfield nets use binary units, meaning each unit/neuron can be either "on" (having the value 1) or "off" (having value -1 / 0).

A Hopfield net has recurrent connections with symmetric weights but no self connections: - $w_{ii} = 0$. - $w_{ij} = w_{ji}$.

The activation of a unit depends only on its neighbour unit and is performed using the following rule.

$$s_i = 1 \text{ if } \sum w_{ij}s_j - \theta_i \geq 0$$

The updates can be performed in two different way : - Asynchronous: One unit is updated at a time. The units are either chosen randomly or in a predefined order. - Synchronous: All units are updated at the same time (based on the previous values)

For each state there can be a "energy" assigned similar to energy based models. For a given state the energy of a Hopfield net is defined as:

$$E = -0.5 \sum w_{ij}s_i s_j + \sum \theta_i s_i$$

Properties / Advantages / Disadvantages Hopfield net as energy based model ($E = \dots \rightarrow p(s = 1) = \text{sgn}(E)$)

With each (asynchronous) update step the energy is thus guaranteed to stay the same or lower in value.

And the network will converge to a local minimum of the energy function, similar to a stable equilibrium state, where it can not escape from. Such a state can be called attractor state.

Hopfield nets can be trained as associative memory, where each stored pattern corresponds to an attractor state.

Such Hopfields net can perform pattern completion. But it can also end in a spurious state, an attractor state/local minimum, which was present in the training data.

Hopfield nets can be trained with a local Hebbian rule, i.e. for an update they only use information of neurons on either side of a connection:

$$w_{ij} = 1/n \sum^{patterns} s_i s_j$$

Boltzmann Machines

Boltzmann machines try to improve Hopfield nets by replacing the deterministic update rule with a stochastic one.

Model The units are similar to the Hopfield net.

The weights are similar.

The Energy is similar.

The Update rule can be written as: $p_{on} = 1/(1 + \exp(-E))$

A Boltzmann machine is an energy base model with the prob distribution defined by the Energy. Thus each state can be assigned a Energy which is directly indicates its probability.

To increase the capacity of the network latent variables are introduced. The units are divided into visible units which can be directly observed and hidden/latent units. With hidden units a

Boltzmann machine becomes a universal approximator probability mass functions over discrete variables.

E can then be written as $E(v, h)$

Each update step can be seen as gibbs sampling from the prob distribution.

Learning Rule Lowest energy for a trainig sample (= high probability)

Derivate E by weights $\Rightarrow p_{data} - p_{model}$

Nuding the variables/ weights to the data distribution

RBMs Reaching a eq state in BMs is difficult, since the hidden layers are fully connected (and conditioned on the visible data) still potentially infinite sampling steps have to be performed sequentially.

To evade the problem, a simple solution is to make the hidden units not depended on each other (and the visible not depended on each other), so all hidden units can be sampled independed and parallel to each other.

Thus a RBM is a Boltzmann machnine with two bipartite layers.

CD-Training Data distribution sampling easy. (postive)

Model distribution sampling hard, but short cut works (n sampling / up down steps \rightarrow CD n update)

Deep Belief Networks

A deep belief network (DBN) is a generative graphical model, or alternatively a type of deep neural network, composed of multiple layers of latent variables ("hidden units"), with connections between the layers but not between units within each layer.

A deep belief network can be build up by stacking up RBMs.

This does not only able to unsupervised train a deep belief net, but also improve the conditional distribution of the bottom RBM (Each time we replace the prior over the hidden units by a better prior).

In a deep belief net the two top layers have similar to a RBM symmetric weights while the symmetry is broken up between the lower layers to allow unsymmetric weights (Hinton Example).

Stacking Up RBMs To train a DBN , RBMs can be trained greedily and layerwise: 1. For each Layer train an RBM with CD to obtain the weight matrix W 2. Train the next RBM on top of the other one, but transform the input data using the previous RBMs (either sample hidden units or take the mean activations)

Fine-tuning The final DBN can be fine-tuned using the wake-sleep algorithm: 1. Do a stochastic bottom-up pass, and for each layer adjust the top-down weights, to be better at reconstructing the activation in the layer below.

2. Perform sampling steps in the top level RBM and adjust the weights with CD.
3. Do a stochastic top-down pass, and for each layer adjust the bottom-up weights, to be better at reconstructing the activation in the layer above.

Alternatively, if labels and an error function is given, back propagation can be performed to further fine tune the weights.

2.2.3. Continuous

While all the previous model did run at discrete time steps the next models will run in continuously which makes them more similar to natural neurons and neural networks.

Neuron Models

LIF The leaky integrate and fire neuron which phenomenological describes the membrane potential at the soma. It is one of the simplest and thus computationally most efficient, most important and popular spiking neuron model.

By discarding the different forms/shapes of the action potentials and reducing it the uniform spike events, the information is condensed to the precise spike times.

The model, described by linear equations, models the membrane potential integration due to spike input currents with a capacitor and introduces a leaky current with a resistor.

The model can be represented by a circuit of a single capacitor and a resistor with a battery.

FORMULAR

If the membrane potentiality exceeds a threshold a spike is emitted and the membrane potential is reset to its resting potential (and clamped to that for a given refractory period / time).

FORMULAR

Input: CUBA vs COBA

The spike can be modelled by different functions, among which the most commonly used are:

1. The alpha function
2. The rectangular function.

With those, the leaky integrate-and-fire neuron can be described by the following equations:

FORMULARS

Due to its simplifications the LIF model can not capture some natural observed behaviour such as long-lasting refractoriness or adaptation.

Hodgkin-Huxley The Hodgkin-Huxley model tries to improve some of the limitations of the LIF model.

The model explicitly allows to model different ion channels.

The first model described by Hodgkin-Huxley introduced three different ion channels, namely sodium, potassium and a leak current of Cl^- ions, which they discovered in their experiments on axon of a squid.

Each channel is described by a resistor with a battery.

The Hodgkin-Huxley model with three ion channels can be described by the following equations :

FORMULARS

where the gating variables are described by :

FORMULARS

This model can be extended/generalized to cope more three ion channels with their dynamics, to better match the characteristics/biophysics of different neurons in the brain:

FORMULARS

While this allows to model to predict and simulate various effects observed in the brain, like frequency adaptation, with high accuracy, the Hodgkin-Huxley model is more computationally expensive.

Poisson A Poisson neuron, produces stochastic firing according to a Poisson process.

The firing rate λ or rate function $\lambda(t)$ determines the dynamics of the homogeneous or inhomogeneous Poisson process and thus of the spike times.

The probability of a spike is given by:

FORMULAR

where the occurrence of a spike is in depended off previous spikes.

Since the spikes are produced by a Poisson process, the expected number of spikes for an interval is given by :

*FORMULAR***Neural Coding**

Spikes are the main way information in the brain is transmitted/exchanged.

This information can be encoded in different ways.

Three ways, which were observed in the brain, are rate coding, temporal coding and population coding.

In rate coding the information is purely in the spike frequency.

In temporal coding the information is in addition transmitted via the different points in time at which spikes arrive.

In population coding the information is distributed over the joint activity of several neurons.

Learning

Learning in the brain describes the generalized term, how information in the brain is stored (in contrast to learning a task memory etc is also considered learning).

Most models of learning algorithms in spiking neural networks build on changes in the synaptic weights/ strength between to neurons to store information.

For changes to be seen as learning only time-spans lasting minutes to days or more are considered, which are covered by LTP (long-term potentiating) and LTD (long-term depression).

The models most commonly used are inspired on the research and discoveries of Hebb and his Hebb principle.

STDP Spike time depended plasticity poses such an learning algorithm which was inspired by research on single neurons with artificially induced current.

The experiments indicated a increase of the synaptic weight if a post-synaptic spike occurred in strong temporal vicinity after a pre-synaptic spike and a decrease of the synaptic weight if a post-synaptic spike occurred in strong temporal vicinity before a pre-synaptic spike. If the further apart the spike times of the different neurons were, the weaker was the effect on the spike.

This lead to the following update rules:

FORMULARs

Symetric STDP

3. Related Work

3.1. Convolutional RBM

The convolutional RBM was invented more or less at the same time by Bengio and Lee.

In similarity to CNN it can be seen as the advancement of energy based model to compositional data.

Describing images in terms of spatially local features needs fewer parameters, generalizes better, and offers re-usability as identical local features can be extracted from different locations of an image. cRBM have shared weights and are not fully connected.

For propagating information up can be seen as the convolution/cross correlation with a filter matrix. The down propagation part uses the flipped kernel (since $w_{ij} = w_{ji}$) (with some padding).

FORMULARS

Images.

Learning is similar to normal RBMs with CD.

Lee proposes an softmax based probabilistic max pooling to introduce sparseness in the hidden activations (for a neighbourhood), on which a pooling layer can be stacked.

This can also be achieved by setting the bias to a negative value.

3.2. Sampling in SNNs

One new encoding of information in the brain has indicated by stochastic neural transitions in the brain, stochastic inference.

The first framework, which allowed MCMC sampling with spiking neurons was introduced by Buesing.

One simplification which can be made is a discretization of the time in time slices (see Buesing for the generalization to continuous time).

Buesing proposes a stochastic neuron model which activates with a probability proportional to the input.

If the neuron spikes its state is set to $z=1$ and it stays in the $z=1$ state for its refractory period $z=0$.

A neuron which is not in refractory period is given the state $z=0$.

Thus for a given time step the state of the network is defined by the state of the individual neurons.

IMAGE

To get the markovian properties ($p(z|z)$) back, a auxiliary variable is introduced which indicates the time a neuron has to stay in the refractory period.

It can be shown, that a update in a such networks can be seen as a gibbs sampling step and therefore the network performs gibbs sampling.

Experiments show, as $t \rightarrow \infty$, the network in fact is able approximates a boltzmann distribution.

Replacing the rectangular PSP with a more biological plausible alpha shaped PSP deteriorates the performance, but is still reasonable well.

Petrovici improved the model by replacing the stochastic neuron model by LIFs a more common and biologically inspired neuron model.

He proved under high frequency (poisson) noise, which leads to a high conductance state of the membrane potential and therefore to a low membrane time constant, the membrane potential can be seen assumed to be normal distributed.

FORMULARS

This leads to a stochastic membrane potential as well as to stochastic firing if the membrane potential crosses the threshold.

Consequently he defines the probability of a neuron to spike as the number of spikes in a given time period in relation to the total number of spikes which could have occurred.

FORMULAR + Example

Here it is important to note, the given the HCS state the time from a neuron to go from the resting potential to the spiking threshold can be neglected (important !!!).

This allows the neuron to show a firing behavior which can be matched by a logistic function.

By normalizing the weights, the LIF neurons can so perform neural sampling similar to Buesing.

3.3. NN -> SNN conversion

This allows a quite simple transformation of a (offline) trained boltzmann machine to neural networks, where the weight have to be scaled *FORMULAR* and the previous described LIF neurons can be used.

Connor uses a different approach, where he instead of approximating sigmoid units by LIF neurons, uses the sigert neuron description to implement units in the RBM which activate similary to LIF neurons. Such a trained net can be directly transfered to a SNN without any adaptation.

There also have been different approaches to transform a CNN to a SNN (see seminar paper).

3.4. eCD and Sampling Machines

Different approaches have been proposed to train a rate based RBM, where the first was probably by Hinton. They use multiple binary stochastic input units of the same input to simulate a rate based input.

The next approaches make use of the synaptic sampling described in the previous chapter.

One approach is the evtCD by Daniel Neil which works in continuous time with spiking networks and a STDP variant. He simulates the positive and negative phase of the CD by simply

unrolling the RBM (with shared weights). But this approach only allows a certain number of CD steps and is due to the weight synchronization not very plausible.

The more sophisticated approach which also uses STDP was proposed by Neftci. He uses bidirectional synapses, between a visible and hidden layer of LIF neurons. In their approach the training time is divided into four parts: 1. The data signal is applied and the system is allowed to model the data distribution 2. Positive STDP is used to get $v_i h_j$ -data (with stdp) and is added to the weights (positive phase) 3. The data signal is removed and the system is allowed to model the model distribution 4. Negative STDP is used to get $v_i h_j$ -model (with stdp) and is subtracted from the synaptic weights.

$$w = w + w_{pos} - w_{neg}$$

4. Approach

We evaluate two different approaches to get convolutional deep belief networks.

The first one is the most classical one, where a offline/ in discrete time trained DBN is transferred to the spiking domain.

The second approach trains the RBMs event based with STDP directly in the spiking domain.

4.1. Conversion

The conversion can be roughly described in two steps: 1. Train the RBMs offline and build up an DBN 2. Convert the DBN to the spiking domain

4.1.1. Conv DBNs

To train convolutional DBNs we proceed similar to Lee.

At first the RBMs are trained greedily with CD.

After we stop training the first RBM, we convert the dataset, by sampling of the hidden layer of the trained RBM (one sampling/forward-pass step), into a new dataset.

This can be simply performed by multiplying the dataset with the trained weights, using the sigmoid activation function, and performing Bernoulli sampling on the activations.

By converting a data sample, we hope to extract features/ the most important structure of the data sample.

On this converted dataset the next RBM is trained with CD.

To get a measurement of the feature quality we integrate label information and perform classification with the DBN.

Our approach is similar to Hinton. The last layer of the DBN is a fully connected RBM with input consisting of the last converted dataset as well as the label. The RBM is also trained with CD to associate the label, with the converted dataset/features.

To evaluate the performance, we transform the data sample, and to the top RBM we only input the converted data sample with no label input. We then let the RBM predict a label by performing up and downward passes, with the data input fixed.

4.1.2. Conversion

For the conversion we have three different variations

Conversion as CNN One way to convert the DBN to the spiking domain, is by interpreting it as a pretrained CNN (we do not perform any commonly used gradient descent fine tuning to

get comparable results with only CD trained models, but fine tuning could further improve the performance).

For the conversion, we proceed similar to Cao and Diehl. They use avg pooling and relu functions, to get a similar architecture as SNNs. But in contrast, we don't use any pooling, since for RBM there is no simple way to integrate avg pooling (what Cao and Diehl use), and for spiking CNN there is no simple way to integrate max pooling. We also use the sigmoid function, since RBMs are commonly trained with sigmoid activations (even so Hinton proposed relu for rbms as well) and the "activation" function of rate based LIF neurons with a refractory period matches the sigmoid function more closely.

DBN layer is replaced by a layer of LIF neurons. The connections are replaced by (directed) synapses, with the weights of the DBN synapses scaled with a constant factor, to get similar activations.

Conversion with COBA LIF Another way to convert the DBN to the spiking domain, is by interpreting it as a directed graphical model, a sigmoid belief network, and perform ancestral sampling.

This approach is heavily based on the synaptic sampling theory, i.e. it uses spiking neurons to perform sampling.

The sampling can be either performed by conductance base LIF neurons or current base LIF neuron as described by Petrovici.

For the COBA neurons, we choose a biological plausible neuron model (see parameters in table). The high conductance and increased mean potential (gaussian distributed) and thus a firing probability of .5, is achieved by using high frequency Poisson generated (inhibitory and excitatory) spikes to bring the neuron to a high conductance state (HCS).

This neuron model has an input-current/spikes to firing frequency mapping/ transfer function which approximates a sigmoid function.

Image

The PSP are chosen to have an alpha shape instead of an t_{rec} long rectangular PSP, which according to Buesing may introduce some discrepancies when performing sampling in comparison to Gibbs sampling.

IMAGEvglrbsamplings

The DBN is converted by simply converting each layer to a layer of CUBA LIFs with poisson noise, and the connections are transformed to synapses with the weights scaled to achieve a function similar to the sigmoid function.

Consequently the DBN simply performs ancestral sampling with the data sample as evidences and the label as inferred state.

Conversion with CUBA LIF To reduce the computational expenses of the COBA models, they can be replaced by less computationally complex CUBA models.

Their model parameters are chosen to simulate a HCS state.

Therefore the membrane time constant as well as the membrane conductance are reduced and a static input current is inserted.

By adding high frequency poisson noise, a sigmoid shaped frequency mapping/ transfer function is achieved.

Imagesigmoid/boltzmann

The DBN conversion is similar to the COBA case, where the COBA neurons are now replaced by the adapted CUBA LIFs.

4.2. eCD

Another approach tries to learn the conv spiking DBN online with an STDP based learning rule.

The main idea of the learning rule is adapted from Neftci.

We also adapt this LIF neuron model.

FROMLAR

A similar STDP rule is used, but we extended the model with a learning rate. This rule can be presented as an iterative rule as follows:

FORMULAR

The training can basically be divided in 4 phases.

While this poses similarities to pCD since the activity of the hidden layer of the previous step is used as starting state for the next step, we extend the model by a 5th phase between the two data samples, where the model is "flushed" thus enabling normal CD.

STDPCURVE

4.2.1. Convolution

We implement convolution with local receptive field and sharing weights between neuron through weight synchronization.

The receptive fields have the same size and each receptive field is connected to a neuron.

The receptive fields are overlapping and next to each other, with the output neuron having the same topology as their input regions.

The receptive fields cover partial regions of the input data, where the weight of a neuron covering a certain region can be seen as a (flipped) convolution filter over the partial input data.

Since the weights of different neurons with different feature maps are shared, the neural activity can be seen as a convolution over the input data with a (flipped) filter of the size of the receptive fields, where the weights are the shared weights from the receptive fields to the neurons. This can be also called the feature map of the convolution.

By using more layers of neurons over the same receptive field, where each layer shares one set of weights, more feature maps can be generated.

IMAGES

Since each weight has their local STDP based update rule (eCD), we have to find a way to synchronize/share weights between all neurons in a layer.

To keep the weights in neuron in a layer the same, we perform a weight synchronization step at discrete time steps, since updating all weights after a single update did not show any promising results.

Thus the synchronization at a time step can be described by the following rule:

$$w_1 = w + \delta w_1 \quad w_2 = w + \delta w_2$$

\implies

$$w' = 1/2(w_1 + w_2) = 1/2(w + \delta w_1 + w + \delta w_2) = w + 1/2(\delta w_1 + \delta w_2)$$

Thus similar to conv RBMs we can simply take the mean of the weight changes and apply it to all the weights, which is equal to just using the average of the new weights.

In addition we introduce fixed negative connections between neuron in the top/ hidden layers (biological plausible \rightarrow more sparse King). This removes one advantage of the RBMs since the hidden layers are no longer independent, which makes it harder to sample from the true distributions, but since the network continuously performs sampling steps, the approximation appears to be good enough, if the weights are not too strong and prevent a change to a different mode.

By connecting neurons to neurons on a similar position in another feature map, appears to make the features more discriminative and less correlated. Also this poses some similarities to adding a negative (here: structures) bias to the hidden units, which has shown to result in better features (see Norouzi M)

An intuitive interpretation, is that if one feature reacts to a certain input it will be highly active and prevent the others from being active as well and thus prevent them from learning the same.

4.2.2. DBMs

To build up a spiking DBN we train the BMs layer wise and forward the input of the previous layer to the next layer.

To be exact the spiking DBN is a mixture between a DBN and a DBM, since the single BMs are still bidirectional connected and only the hidden layer is forwarded in a directed manner.

IMAGE

Converting a RBM to a directed BN did not show any good results, since the activation of the hidden unit turned out to be important for generating a good estimate of the data distribution.

Due to some top-down influences, when stacking a new BM on the trained BM, the hidden distribution gets distorted and unfit input for the new BM to be trained on.

In our case we simply forward, the hidden activations to the input layer of a new two layered BM. An approach similar to Salakhutdinov's DBMs with doubling the weights and composing the models in the end may sound promising and could be tried as well.

5. Implementation

5.1. Discrete DBNs

The DBNs were implemented in the Theano-Framework, to utilize the computational power of the GPU. The implementation of the single RBMs is adapted from the official Theano RBM (REF).

An upward pass is performed as follows: As an input we take a 4D tensor in the form of (batch size, input channels, input height, input width), and for the upwards pass convolve it with a filter of the shape (number of filters, input channel, filter height, filter width) with a stride of 1. This results in feature maps of the shape (batch size, number of filters, input height - filter height + 1, input width - filter width + 1). A constant c is added to the activation as bias. On those feature maps a sigmoid function is applied. Afterwards the activations are sampled with a Bernoulli distribution (FRMLA), which results in the activation of the hidden units.

The downward pass on the activation of the hidden units is performed similarly: As input we take the activation of the hidden unit, and convolve it with the 180 degree flipped kernel, where the first and second dimensions are swapped. Afterwards again a sigmoid function is applied and accordingly sampled to get a new visible layer activation.

One Gibbs sampling step consists of one upwards and a consecutive downward pass.

The free energy of the model is calculated using formula (TODO ref) on output of the convolution on the visible units.

After performing one Gibbs sampling step to get the distribution of the data (positive phase), we perform $k-1$ additional Gibbs sampling steps to get an estimate of the model distribution (negative phase).

As error function we define the difference between the free energy of the positive phase and the negative phase (TODO ref formula). We use Theano's auto-differentiation to determine the gradient and perform gradient descent with a learning rate lr and a weight decay of x .

FORMULA

We train the RBM for a given number of epochs, with a batch size smaller than the dataset, thus performing stochastic gradient descent.

To build up the DBN each RBM was trained greedily. After training one RBM the entire dataset was converted by doing one upwards pass in the trained RBM and using the activation of the hidden units as the new input data for the next RBM.

PSEUDOCODE

The last layer consists of a fully connected RBM where the input data consists of the converted data set concatenated with a one-hot encoding of the label.

This is the first time the label is used to train an RBM and does not affect the extracted features of the DBN up to this point.

5.2. Conversion

To simulate the converted DBNs we use the pyNN framework with Nest as spiking network simulator.

To simulate the CNN and DBN we use LIF with CUBA and COBA units neurons, respectively. The parameters can be seen in Table x(@TODO table).

Each unit in the DBN is represented by a neuron.

A layer in the DBN is represented by a (not interconnected) neuron population.

If a unit in the DBN is connected to one in a higher layer with a weight w_{ij} , a synapse is added with a the between the two neurons representing the units with the scaled weight $w_{ij} * scaleFactor$

The input is transformed to a Poisson distributed spikes, where the rate of the Poisson process is given by the original data value/ image intensities scaled with some constant offset to generate more frequent spikes.

The input is directly fed into the bottom layer neurons with a fixed weight w_{in}

The network is run for a runtime of t .

To get the classification results, the spikes in the label layer are counted to determine the label we take the index of the neuron with the max activation:

$$\operatorname{argmax}_a(x)$$

5.3. eCD

The eCD online learning was implemented in PyNN with Nest as well as Brian simulator, but due to the simulation speed we chose Brian for most of our experiments.

As neuron type we again chose a exp COBA LIF neuron with high frequency noise.

Each input element gets transformed to Poisson distributed spikes, where the rate is proportional to the input value. E.g. for a 100 dimensional data, 100 Poisson spike generators with different weights are used.

Each RBM consist of a visible and hidden layer. The visible consist of one neuron population with # input data dim of neurons representing the input, the data population. In addition a second neuron population represeting the label input, the label population, can be added to the visible units. The hidden layer is an neuron population of the size nmbr filters x (input height - filter height + 1) x (input width - filter width + 1).

The data population is sparsely connected to the hidden layer with synchronized weights, while the label population is fully connected to the hidden layer.

IMAGE

In the hidden layer we connect each neuron to all other neurons at in a square at same position in a different feature map with a inhibitory synapse with a fixed negative weight.

IMAGE

The training of the spiking RBM is performed with eCD:

In the first phase (Data burn-in), we input layer is induced with a strong negative current and the data input is fed as spikes with a high synaptic weight, so that the visible layer is unaffected

from any spikes in the top layer and only spikes in accordance to the input data. The hidden layer is unaffected. The STDP learning flag is set to $g = 0$, so no learning is allowed

In the second phase (Data distribution), now the STDP learning flag is set to $g = 1$ so positive learning is allowed. This should drive the weights to represent the data distribution.

In the third phase (Model burn-in), we set the data input to the input layer to zero and remove the induced negative current, to let it reach the model distribution. In this phase the learning is disabled, setting the STDP learning flag to $g = 0$

In the fourth phase (Model distribution), the STDP learning flag is set to $g = -1$, enabling only negative (un-)learning. This will unlearn the model distribution.

In the optional fifth phase (Flush), we induce a strong current to the visible and hidden layer to remove all activity and allow a fresh start in the first phase.

IMAGE

The training of one data sample is set to last $n * t_{ref}$ cycles.

The weight synchronization is set at discrete time points, after n training samples. This is performed by taking the mean over all the weights, which are to be the shared. In addition a small weight decay is introduced:

$$w = \text{mean}(w) - \text{decay} - \text{rate} * \text{mean}(w)$$

Each RBM is trained with m data samples.

After a RBM is trained, the next RBM will be trained on the output of the previous RBM. Therefore the a connection with a strong synaptic weight is established from the hidden layer of the bottom RBM to the top RBM.

The original input data is still fed to the bottom RBM while the training data for the top RBM is now the activations of hidden layer in bottom RBM.

At test time the data is fed to the bottom RBM and the network is run for a fixed timespan t_{test} . There is no external input forwarded the label layer and the number of spikes in the label are counted. The index of the label with the most spikes represents the predicted data class.

6. Experiments&Results

6.1. Datasets

We evaluate our models on three different datasets. The dataset size is primarily limited by the computational resources, such as memory and computation time.

6.1.1. 1x4 Dataset

The simplest dataset we use a 1 dimensional binary dataset of size 4, where either the first or the last element is set to 1.

6.1.2. Strip Dataset

We generate a 10x10 pixel noisy stripe dataset, with three different oriented stripes, horizontal, diagonal, vertical.

This could represent a similar object to a pen recorded with a event-based camera, and result in an grasp id.

In the easiest version of this dataset, the stripes always occur on the same places, with some random noise.

An more complex version of the datasets contains the stripes randomly distributed across the whole image.

The dataset can be either binary or continuous.

6.1.3. MNIST

We also evaluate the models on the MNIST dataset.

The MNIST dataset consists of 60000 28x28 pixel gray images of the numbers 0-9.

We evaluate our models on the normal MNIST dataset, as well as a to dvs events converted version of MNIST.

6.2. Experiments

We orient our experiments primarily on the strip dataset, due to time and computational constrains.

After evaluating our models on the strip dataset, we look on the performance on other datasets for comparison and generalization.

6.2.1. Convolution vs no Convolution

6.2.2. Lateral connections

6.2.3. Hidden sparsity/ learning the data distribution

7. Conclusion and Outlook

7.1. Biological plausibility

Studies by Hubel and Wiesel suggest, certain neurons respond so similar features at different position of in visible field.

This suggest similar synaptic weights.

One explanation of this could be shared weights similar to CNNs.

Since the weight updates in the brain are primarily believed to be local, there is no know principle to keep weights between different neurons in the brain synchronized.

So while the trained structure, with receptive fields and similar weights is quite plausible, the training procedure here is not.

A more plausible way in the brain to get similar weights, is due to the similarity of the inputs, e.g. if two receptive fields get quite similar input, their weights will probably converge to the same target values.

While this requires all receptive fields to be presented with the whole data, presenting each field with some part of the data but updating all fields with a combined update can be more computational effective. This could be a principle CNN utilize to get biological plausible result, while performing completely biological plausible updates.

Another biological not completely plausible part of our presented system are the bidirectional synapses.

This in turn could be easily translated to two directional synapses, with some weight synchronization. While in this case local updates are sufficient, to keep the weights similar (e.g. applying a similar learning rule to both weights), and research on discrete NNs has shown some automatic weight synchronization in Autoencoders (Bengio), where is no biological proof.

The STDP flag determining either completely positive or negative does not appear to be plausible as well.

Even so this system has many constrains, it might could be counted among one of the more biological plausible deep learning architectures.

A. Some appendix

A.1. (if needed)

Liste der noch zu erledigenden Punkte

B. List of Figures

C. List of Tables

D. Bibliography

- [1] G. Berndes, M. Hoohwijk, and R. van den Broek. The contribution of biomass in the future global energy supply: a review of 17 studies. *Biomass and Bioenergy* 25, pages 1–28, 2003.
- [2] W. Feng, H. J. van der Kooi, and J. de S. Arons. Phase equilibrium for biomass conversion processes in subcritical and supercritical water. *Chem. Eng. J.* 98, pages 105–113, 2003.
- [3] X. H. Hao, L. Guo, X. M. Mao, and H. J. Chen. Hydrogen production from glucose as a model compound of biomass gasified in supercritical water. *Int. J. Hydrogen Energy*, pages 53–64, 2003.
- [4] R. A. H. III. The age of energy gases. *Int. J. Hydrogen Energy* 27, pages 1–9, 2002.
- [5] I.-G. Lee, M.-S. Kim, and S.-K. Ihm. Gasification of glucose in supercritical water. *Ind. Eng. Chem. Res.* 41, pages 1182–1185, 2002.
- [6] Y. Matsumura and T. Minowa. Biomass gasification in near- and super-critical water: Status and prospects. *Biomass and Bioenergy* 29, pages 269–292, 2005.
- [7] M. Osada, T. Sato, M. Watanabe, T. Adschiri, and K. Arai. Low temperature catalytic gasification of lignin and cellulose with a ruthenium catalyst in supercritical water. *Energy & Fuels* 18, pages 327–333, 2004.
- [8] I. T. and Y. Matsumura. Gasification of cellulose, xylan and lignin mixtures in supercritical water. *Ind. Eng. Chem. Res.* 40, pages 5469–5474, 2001.
- [9] S. T, S. Kurosawa, R. L. Smith, T. Adschiri, and K. Arai. Water gas shift reaction kinetics under noncatalytic conditions in supercritical water. *J. Sup. Fluids* 29, pages 113–119, 2004.
- [10] M. O. T. Sato, M. Watanabe, M. Shirai, and K. Arai. Gasification of alkylphenols with supported noble metal catalyst in supercritical water. *Ind. Eng. Chem. Res.* 42, pages 4277–4282, 2003.
- [11] J. D. Taylor, C. M. Herdman, B. C. Wu, K. Wally, and S. F. Rice. Hydrogen production in a compact supercritical water reformer. *Int. J. of Hyd. Ene.* 28-11, pages 1171–1178, 2003.
- [12] T. Yoshida, Y. Oshima, and Y. Matsumura. Gasification of biomass model compounds and real biomass in supercritical water. *Biomass&Bioenergy* 26, pages 71–78, 2004.
- [13] Y. Yoshida, K. Dowaki, Y. Matsumura, R. Matsuhashi, D. Li, H. Ishitani, and H. Komiyama. *Comprehensive comparison of efficiency and CO2 emissions between biomass energy conversion technologies-position of supercritical water gasification in biomass technologies*, *Biomass and Bioenergy*. 2003.

- [14] D. Yu, M. Aihara, and M. J. Antal. Hydrogen production by steam reforming glucose in supercritical water. *Energy & Fuels*, 7, pages 574–577, 1993.