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# Eidgenössische Technische Hochse Swiss Federal Institute of Technologie

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A dissertation submitted to ETH Zürich in partial fulfilment of the requirements for the degree of Master of Science ETH in Computer Science

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## **Declaration**

I David Yenicelik of ETH Zürich, being a candidate for the M.Sc. in Computer Science, hereby declare that this report and the work described in it are my own work, unaided except as may be specified below, and that the report does not contain material that has already been used to any substantial extent for a comparable purpose.

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# Abstract

Write a summary of the whole thing. Make sure it fits in one page.

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## Chapter 1

## Introduction

In Natural Language Processing (NLP), bilingual lexical induction (BLI) is a problem of inferring word-to-word mapping between two languages. While supervised BLI may be learned trivially from a dictionary, unsupervised BLI is highly non-trivial, and serves as a backbone to many unsupervised Neural Machine Translation (NMT) systems, without which the overall MT performance drastically drops [?] [?]. In addition to the unsupervised setting, words in a source language A often do not carry a one-to-one correspondence with words in a target language B. This further increases the difficulty of finding a (bijective) mapping between the two embedding spaces.

## 1.1 Scope of Work

Continuing where [?] left off, we want to investigate the performance of using normalising flows to model unsupervised lexicon matching between two probability distributions, which are defined by Gaussian embeddings, which have a one-to-one correspondence to tokens in the respective languages. We aim to use Gaussian embeddings for the robustness, and better integration into the probabilistic perspective. The method discussed in the paper also relies a lot on the (semi-)supervised loss component. We aim to investigate why this is the case, and would like to revise a method which is more robust

in the fully unsupervised setting.

Minimal goals: We propose the following steps on achieving this goal.

- 1. Replicate the algorithms and models which can generate through Gaussian embedding [?]. Do one sanity check by doing a sanity check on one of (SimLex / WordSim)
- 2. Replicate "Density matching for bilingual word embedding" to setup a baseline normalising flow between vector-word-embeddings [?].
- 3. Define loss between predicted and target embeddings (if change in definition is necessary)
- 4. Change the embeddings in point 2. to use Gaussian embeddings. Implement Loss functions found in point 3.

**Extended goals:** If the above points provide good performance, we would like to expand on the below points.

- 1. Implement a fully unsupervised extension by investigating the short-comings of [?].
- 2. Investigate "deeper" normalising flows than the linear flow in [?] as such [?] [?].

Contingency Plan / Further extended goals: Implementing the following points would allow for an applied perspective of this approach, showing that this methodology allows for more robust mapping, also outside the field of NLP.

- 1. Train embeddings for job-systems ESCO and AUGOV using Skip-Gram or Co-Occurence-matrix based. Do a sanity check.
- 2. Train Gaussian embeddings for job-systems. Do a sanity check.
- 3. Generate a small validation dataset between the European- and Australian job-system.

- 4. Setup some baseline algorithms based on NLP, graph-matching, colinear-PCA for matching as a non-NLP benchmark environment. Compare against above-proposed methods.
- 5. Find a normlising flow model to transform one job system into another.

## Chapter 2

# Background

We provide a short introduction to the background work.

## 2.1 Gaussian Embeddings

The problem that Gaussian Embeddings try to resolve is to find a suitable embedding space for entities, such as word-tokens.

In the context of NLP, the general goal is to map every word type w that is included in some dictionary  $\mathcal{D}$  in the latent space, such that w is close to all the context words c from another dictionary  $\mathcal{C}$ . This follws the (CITE THE RANDOM GUY) ASSUMPTION, which says that words in similar contexts have similar neighborhoods.

In an unsupervised context, we observe a sequence of words  $t(w)_i$  for each word-token  $w_i$  and their contexts  $c(w)_i$ .

We want to use some rank-based Energy, not absolute energy, such that we can make sure that neighboring words are pushed together, and two "random" words are pushed further away from each other.

The above two problems can be regarded as finding an invertible transformation from one (embedding) space  $\mathcal{X} \in \mathbf{R}^d$  into another  $\hat{\mathcal{X}} \in \mathbf{R}^d$ . Normalising flows [?] [?] have proven to be a powerful tool in modeling such relations. As such, the aim of this project is to find a model based on normalising flows which is able to find an an invertible mapping f from  $\mathcal{X}$  to  $\hat{\mathcal{X}}$ . To keep the discussion focused, this thesis deals with the problem of finding a model for bilingual lexicon matching.

#### Word representations via Gaussian Embeddings

Gaussian Embeddings have been first proposed in the context of words, although prior work has been adapted in embedding matrix-rows into a mean and standard deviation (CAN CITE THE PEOPLE IN THE PAPER 2007, 2008). It is a continuous probabilistic relaxation of the otherwise common discrete point vectors. Each word is represented by a Gaussian distribution in high-dimensional space, allowing to better capture uncertainty and a representation and it's relationships. It can also express asymmetric relations more naturally than dot-products or cosine similarity, and enables for better-parametrized rule between decision boundaries.

Fitting Gaussian Mixture Models on embeddings have been done in order to apply Fisher kernels to entire documents.

Because this is an unsupervised learning task, we must use an energy function which is incorporated within a loss function that we try to minimize. The energy function describes (dis-)similarity between two items. The authors propose the following energy functions to derive a Gaussian word-embedding.

$$L_m(w, c_p, c_n) = \max(0, m - E_{\theta}(w, c_p) + E_{\theta}(w, c_n)$$
 (2.1)

Here, w is the word we want to sample,  $c_p$  is a "positive" context word, i.e. a word that co-occurs with the word w, and  $c_n$  is a "negative" context word, i.e. a word that does not co-occur with the word w. Usually the negative context words is sampled randomly from the corpus. The loss function reminds of a hinge-loss in logistic regression.

The authors propose two possible ways to learn the mean and variance of the Gaussian embeddings. They argue that the empirical covariance is not the

most effective method of deriving the words as Gaussian embeddings. This does not allow for inclusion between ellipsoids

Symmetric similarity: expected likelihood or probability product kernel We can use any kernel (which is symmetric by definition) to derive at an energy function. For two Gaussians f(x), g(x), the inner product is defined as:

$$E_{\theta}(w,c) = \int_{x \in \mathcal{R}^d} f(x)g(x)dx \tag{2.2}$$

$$= \int_{x \in \mathcal{R}^d} \mathcal{N}(x; \mu_w, \Sigma_w) \mathcal{N}(x; \mu_c, \Sigma_c) dx$$
 (2.3)

$$= \mathcal{N}(0; \mu_w - \mu_c, \Sigma_w + \Sigma_c) \tag{2.4}$$

For numerical feasibility and easy of differentiation, we usually maximize the  $\log E_{\theta}(w,c)$  for a given dataset with  $w \in \mathcal{W}, c \in \mathcal{C}$ . We will not go further in what the specific gradient of this log-energy is.

Asymmetric divergence: KL-Divergence We can use more directional supervision to exploit directional supervision, such as a knowledge graph.

Following energy-function is optimized:

$$-E(w_i, c_j) = D_{KL}(c_j || w_i)$$
(2.5)

$$= \int_{x \in \mathcal{R}^d} \mathcal{N}(x; \mu_{w_i}, \Sigma_{w_i}) \log \frac{\mathcal{N}(x; \mu_{c_j}, \Sigma_{c_j})}{\mathcal{N}(x; \mu_{w_i}, \Sigma_{w_i})} dx$$
 (2.6)

$$= \frac{1}{2} \left( \operatorname{tr} \left( \Sigma_i^{-1} \Sigma_j \right) + \left( \mu_i - \mu_j \right)^{\top} \Sigma_i^{-1} \left( \mu_i - \mu_j \right) - d - \log \frac{\det \left( \Sigma_j \right)}{\det \left( \Sigma_i \right)} \right)$$
(2.7)

Because of the loss function, this can entail information such as "y entails x" as a soft form of inclusion between two datasets (if KL divergence is used). If

a symmetric loss function is used, then this would most likely lead to overlap (IS THIS TRUE...???)

Uncertainty calculation: In contrast to the empirical standard deviation as an uncertainty measure, we can now calculate the uncertainty of the inner product (i.e. the distribution  $P(z=x^Ty)$  using the following formula

$$\mu_z = \mu_x^T \mu_y \Sigma_z = \mu_x^T \Sigma_x \mu_x + \mu_y^T \Sigma_y \mu_y + \operatorname{tr}(\Sigma_x \Sigma_y)$$
 (2.8)

We then get an uncertainty bound, where c denotes the number of standard deviations away from the mean.

$$\mu_x^{\top} \mu_y \pm c \sqrt{\mu_x^{\top} \Sigma_x \mu_x + \mu_y^{\top} \Sigma_y \mu_y + \operatorname{tr}(\Sigma_x \Sigma_y)}$$
 (2.9)

We can learn the parameters  $\Sigma$  and  $\mu$  for each of these embeddings using a simple gradient-based approach, where we set hard constraints on

$$\|\mu_i\|_2 \le C, \forall i \tag{2.10}$$

$$mI < \Sigma_i < MI \tag{2.11}$$

The method shows competitive scores to the Skip-Gram model, although usually only with minor improvements depending on the benchmark-dataset.

## 2.2 General random information

For implicit generative models, we can measure the performance in two ways.

Assuming we have two probability density distributions, the true data-distribution p\* and the approximate posterior  $q_{\theta}$  whose goal is to mimick the original data distribution p\*.

One way is to apply the probability difference.

$$r_{\phi} = p * -q_{\theta} \tag{2.12}$$

This in general, however is not very intuitive, and does not easily allow to differentiate between mutliple differences. Loss functions that make use of this logic in clude the Max-Mean discrepancy, or moment-matching loss functions.

Another way it the probability ratio. This ratio is much more intuitive in that one can compare the performance of how different approximate posteriors match the true underlying data distribution.

$$r_{\phi} = \frac{p*}{q_{\theta}} \tag{2.13}$$

Examples include the f-diverence, class probability estimate and the Bregman divergence.

What do GANs usually do. We assume a simple vector z sampled from a feasibly-inferable distribution z.

$$z p(z) (2.14)$$

$$z p(z) (2.14)$$

$$x^{gan} = f_{\phi}(z) (2.15)$$

In general, we try to minimze an energy function, using some sort of reward R and action a taken in some state s. [THIS IS COMPLETE BULLSHIT]

$$\mathbf{E}_{\phi(a|s)}\left[R(s,a)\right] - KL\left[\pi_{\theta}(a|s)||p(a)\right] \tag{2.16}$$

## 2.3 Normalising Flows

In machine learning, we usually distinguish between fully observed models and latent variable models, where we have exclusively observed variables or both observed and unobserved variables.

In general, a generative model can be described as

$$p(x, z, \theta) = p(\theta) \sum_{i=1}^{N} p(x_i | z_i, \theta) \pi(z_i)$$
 (2.17)

And then the model is supplemented by a learning principle such as gradient-based learning (backpropagation), or structural learning (model selection using Monte-Carlo). In general, we want to estimate a probability density distribution by using a stochastic gradient estimator.

$$\nabla_{\phi} E_{q_{\phi}(z)} \left[ f_{\theta}(z) \right] = \nabla \int q_{\phi}(z) f_{\theta}(z) dz \qquad (2.18)$$

$$= \mathbf{E} \left[ \nabla f_{\theta}(g(\epsilon, \phi)) \right] \tag{2.19}$$

$$q(z') = q(z) \left| \det \frac{\delta f}{\delta z} \right|^{-1}$$
 (2.20)

We can now stack multiple bijective and invertible functions together to go from a complex distribution to a simple one (and vice-versa).

(CITE REZENDE ET AL TUTORIAL ON DEEP GENERATIVE MODELS)

$$z_K = f_K \circ \dots \circ f_2 \circ f_1(z_0) \tag{2.21}$$

where the mass-preserving transformation of each function is

$$\log q_K(z_k) = \log q_0(z_0) - \sum_{k=1}^K \log \det \left| \frac{\delta f_k}{\delta z_k} \right|$$
 (2.22)

The loss function which we try to maximize as a product of this then becomes (through the use of expected stochastic gradients.

$$\mathcal{L} = \mathbf{E}_{q_0}(z_0) \left[ \log p(x, z_K) \right] - \mathbf{E}_{q_0}(z_0) \left[ \log p(x, z_K) \right]$$
 (2.23)

Where we used the reparametrization trick in the second equality. We can then use gradient-estimates to miminize the respective log-likelihood.

The REINFORCE (reparametrization trick) has been used in reinforcement learning

$$\nabla_{\theta} \mathbf{E}_{x \approx p_{\theta}(x)} [f(x)] \nabla_{\theta} p_{\theta}(x) = p_{\theta}(x) \nabla \nabla_{\theta} \log p_{\theta}(x)$$
 (2.24)

Rich families of posteriors can be formed.

Generally, if we have N variables, then 3 possible posteriors are (1) the fully connected graph which describes correlations between all variables, (2) a structured approximation model and (3) a fully factored model.

The posterior probabilities of these items respectively are:

1.

$$q^*(z|x) \propto p(x|z)p/z$$

2.

$$q(z) = \prod_{k} q_k(z_k | z_{j \neq k})$$

3.

$$q(z|x) = \prod_{k} q(z_k)$$

(Rezende 2015) A simple distribution is transformed into a more complex one by applying a sequence of invertible transformations until a desired level of complexity is attained.

In the above equation, the KL-divergence is the loss between the approximate posterior and the prior distribution (which acts as a regualizer). The second term, which is the expected log-likelihood is the reconstruction error.

Inference with normalizing flows provides a tighter, modified variational lower bound with additional terms that only add terms with linear time complexity.

in the asymptotic regime, this is able to recover the true posterioir distribution

Approximate posterior distribution of the latent variables  $q_{\phi}(z|x)$ 

$$\log p_{\theta}(x) = \log \int p_{\theta}(x|z)p(z)dz \tag{2.25}$$

$$= \log \int \frac{q_{\theta}(z|x)}{q_{\theta}(z|x)} p_{\theta}(x|z) p(z) dz$$
 (2.26)

$$= \log \int \frac{p(z)}{q_{\theta}(z|x)} p_{\theta}(x|z) q_{\theta}(z|x) dz$$
 (2.27)

$$= \mathbb{D}_{KL} \left[ q_{\phi}(z|x) || p(z) \right] + \mathbb{E}_{q} \left[ \log p_{\theta}(x|z) \right]$$
 (2.28)

$$\geq -\mathcal{F}(x) \tag{2.29}$$

TODO: WHERE DOES THE LAST  $q_{\theta}(z|x)$  disappear to? where  $-\mathbf{F}(x)$  is a free energy function, and where we used Jensen's inequality to obtain the final equation (through  $log E_q\left[\frac{p(x,z)}{q(z)}\right] \geq log E_q\left[\log\frac{p(x,z)}{q(z)}\right]$ ).  $p_{\theta}(x|z)$  is a likelihood function and p(z) is a prior over latent variables, and  $q_{\theta}(z|x)$  is a

(simple) model distribution with which we want to approximate p(x|z). item We try to minimize this loss

To train the model, we need to efficiently (1) compute the derivatives of the expected log-likelihood  $\nabla \phi \mathbb{E}_{q_{\phi}(z)}[p_{\theta}(x|z)]$  and (2) choose the richest, computationally-feasible approximate posterior distribution  $q(\dot{)}$ .

Using Monte Carlo gradient estimation and inference networks, which (when used together), they call *amortized variational inference*. TODO: TIMO SAID THAT THIS IS WRONGLY APPROXIMATING THE TRUE DISTRIBUTION? WHAT ABOUT PSI?

For stochastic backpropagation, we make use of two techniques:

• Reparameterization: The latent variable is reparametrized in terms of a known base distribution and a differentiable transofmration. As an example, if  $q_{\phi}(z)$  is a Gaussian distribution  $\mathbf{N}(z|\mu, \sigma^2)$  with trainable parameters  $\phi = \mu, \sigma^2$ , then we can reparametrize the variable z as

$$z \sim \mathbf{N}(z|\mu, \sigma^2) \iff z = \sigma + \sigma\epsilon, \epsilon \sim \mathbf{N}(0, 1)$$
 (2.30)

• Backpropagation with Monte Carlo. We backpropagated w.r.t. the parameters  $\phi$  of the variational distribution using a Monte Carlo approximation. This is our samples batch.

$$\nabla \phi \mathbb{E}_{q_{\phi}(z)} \left[ f_{\theta}(z) \right] \iff \nabla \phi \mathbb{E}_{\mathbf{N}(\epsilon|0,1)} \left[ \nabla_{\phi} f_{\theta}(\mu + \sigma \epsilon) \right]$$
 (2.31)

where we have simply rewritten z using the reparametrization trick.

For continuous latent variables, it has the lowest variance among competing estimators [CITEEE].

Normalizing flows are most often used in the context of inference network.

An inference network learns an inverse map from observations to latent variables. The simplest latent variable is a Gaussian distribution.

Rezende 2015 propose a *Deep Latent Gaussian Model*, which consists of a hierarchy of L layers of Gaussian latent variables  $z_l$  for layer l.

$$p(x, z_1, \dots, z_L) = p(x|f_0(z_1)) \prod_{l=1}^L p(z_l|f_l(z_{l+1}))$$
 (2.32)

We induce priors over each latent variable  $p(z_l) = \mathbf{N}(0, I)$ , and the observation distribution  $p_{\theta}(x|z)$  is any distribution that is conditioned on  $z_1$  and is parametrized by a neural network.

This model is very general, and captures includes other models, such as factor analysis and PCA; non-linear factor analysis, and non-linear Gaussian belif networks as special cases (Rezende 2014, cited in Rezende 2015)

We know that  $\mathbf{D}_{KL}[q||p] = 0$  is achived when  $q_{\phi}(z|x) = p_{\theta}(z|x)$  (i.e. the approximate q matches the true posterior distribution). IS THIS THE ONLY TIME THIS CAN BE ACHIEVED?

A normalising flow captures more flexibel distributions.

Summary normalising flows:

A normalizing flow is a set of basic rules for transofmration of densiities, considers an invertible, smooth mapping

$$f: \mathbf{R}^d \to \mathbf{R}^d$$

where the input dimensionality  $d_0$  and output dimensionalities  $d_L$  match. The inverse is denoted  $f^{-1} = g$ , i.e. the composition  $g \circ f(z) = z$ . When we want to transform a random variable z' = f(z), we receive the following distribution:

This stems from the Identity that the probability mass must be conversed amongst operations:

$$p_x(x) = \frac{p_z(z)V_Z}{V_X} \tag{2.33}$$

$$z' = f(z) \tag{2.34}$$

$$\frac{z'}{q(z')} = \frac{f(z)}{q(z)} \tag{2.35}$$

#### WHERE DOES THE PRIME COME FROM

We can then construct arbitrarily complex densities by composing several simple maps and successively applying the above mass-preserving operation.

$$z_K = f_K \circ \dots \circ f_2 \circ f_1(z_0) \tag{2.36}$$

We then apply the log-function to have tractable probability functions.

Probability mass must be conserved. From the change of variable formula, we know that taking infinitesimal c:hanges toward the direction of the final probability distribution, we get

$$\int z'q(z')dz' = \int f(z)q(z)dz$$
 (2.37)

$$q(z') = q(z) \ln \left| \det \frac{\delta f(z)}{\delta z'} \right|$$
 (2.38)

There are different ways to use normalising flows:

Although computing the above gradient is expensive  $(O(n^3))$  complexity), we can use the Matrix determinant lemma, and construct out mapping matrices in a special way as follows (https://blog.evjang.com/2018/01/nf1.htm)

$$det(A) = det(L + V \cdot D \cdot V^{T})$$
(2.39)

$$= det(L)(1 + (VD^{\frac{1}{2}})^{T}(L^{-1}D^{\frac{1}{2}}V))$$
 (2.40)

$$W = PL(U + diag(s)) \tag{2.41}$$

Then the cost of computing det(W) becomes O(c) instead of  $O(c^3)$ . (CITING GLOW Glow: Generative Flow with Invertible 1 x 1 Convolutions)

where we have used the Matrix Determinant Lemma for the second equality, with L is a lower triangular matrix, and D is a diagonal matrix.

subsectionExtensions on Normalising Flows

(https://www.shakirm.com/slides/DeepGenModelsTutorial.pdf)

Ever since Normalising flows were proposed, a number of modified versions were presented by [CITE].

Each one calculates the consecutive flow based on a different composition of the previous latent variables  $z_k$ .

#### **Planar Flow**

The planar flow is a simple sequential flow with invertible piecewise linear transformations.

$$z_k = z_{k-1} + uh(w^T z_{k-1} + b) (2.42)$$

### Real NVP (2017)

The input for each step is split into two parts. The intuition behind this is similar to a residual network, which includes logic that allows the gradient to flow back easier.

$$y_{1:d} = z_{k-1,1:d} (2.43)$$

$$y_{d+1:D} = t(z_{k-1,1:d}) + z_{d+1:D} \odot \exp(s(z_{k-1,1:d}))$$
(2.44)

Here, t and s are arbitrary functions, such as deep neural networks or simple linear transforms.

Provides a set of stably invertible and learnable transformations.

This results in an exact log-likelihood calculation.

In the paper, they for the networks s and t, they use residual network with batchnorm and weight normalization.

#### Inverse AR Flow (2017)

According to the authors, this method scales well to high dimensional latent spaces. Each transformation inside the flow is based on an autoregressive neural network.

At each timestep, all prior variables are taken into consideration.

$$z_k = \frac{z_{k-1} - \mu_k(z_{k}, x)}{\sigma_k(z_{k}, x)}$$
 (2.45)

#### Non-Linear Independent Components Estimation (NICE)

Another volume-preserving flow is the NICE flow. First, we arbitrarily partition the latent vector into two components  $z = (z_A, z_B)$ .

$$f(z) = (z_A, z_B + h_\lambda(z_A))$$
 (2.46)

$$g(z') = (z'_A, z'_B + h_\lambda(z'_A)) \tag{2.47}$$

Here,  $h_{\lambda}$  can be chosen in such a way that h is a deep neural network with parameters  $\lambda$ 

LOL, COULD WE JUST CREATE A NEW ONE WITH MORE PROPERTIES??

### Multiple

Normalising flows [?], [?] are a statistical technique where a series of invertible transformations  $f_t$  are applied to a simple distribution  $z_0 \sim q_0(z)$ , to yield increasingly complex distributions  $z_t = f_t(z_{t-1})$ , s.t. the last iterate  $z_T$  has the desired and more flexible distribution. As long as we can efficiently compute the Jacobian determinant of the transformation bijection  $f_t$ , we can both (1) evaluate the density of our data (by applying an inverse transformation and computing the density in the base distribution), and (2) sample from our complex distribution (by sampling from the base distribution and applying the forward transformation) These can be used for classification and clustering [?], [variational inference tasks [?] such as image-generation [?]], enriching the posterior (and prior!) [?], and density estimation [?].

### Glow: Generative Flows with Invertible 1x1 Convolutions (2018)

We use 1x1 convolutional layers as a generalization of the permutation in sequential flows with lower triangular learnable rotation matrices (assuming that the number of input and output channels are equivalent).

TODO: Convolution operators and reduction to permutation operators

$$\log \left| \det \left( \frac{d \operatorname{conv2D}(h; W)}{dh} \right) \right| = h \cdot w \cdot \log |\det(W)|$$
 (2.48)

(read the paper a bit more before citing this, but cite this: https://arxiv.org/pdf/1901.08624.pdf)

The authors cite strong qualitative experimental results in generating faces, thus empirically proving that higher dimensional flows  $(R^d)$  are effective and learn a continuous latent space.

### Flow++ (2019)

The Flow++ model describes three problems with existing methods.

- 1. uniform noise is a suboptimal dequantization which hurts training loss and generalization (i.e. training points are mapped to discrete points in space which have the full probability mass)
- 2. the current models are not expressive enough
- 3. convolutional layers in coupling are not powerful enough

The authors respectively propose three ways to

- 1. variational flow-based quantization (WUT?)
- 2. logistic mixture CDF coupling flows (WUT?)
- 3. self-attention in conditioning of networks

The coupling layers can be described as

$$y_1 = x_1 (2.49)$$

$$y_2 = x_2 \cdot exp(a_{\theta}(x_1)) + b_{\theta}(x_1)$$
 (2.50)

where the functions  $a_{\theta}$  and  $b_{\theta}$  are learnable functions (possibly normalizing flows which ). These have to be invertible affine transformations

It is the first model which starts to close the gap between autoregressive models and flow-based models.

# Chapter 3

## Related Work

As outlined by (https://ruder.io/cross-lingual-embeddings/) the task of finding cross-lingual embedding models fall within one of the four categories:

- 1. Monolingual mapping: Here we train the embedding for each language separately, and the try to find a mapping from one space into another. The learned mapping is usually linear.
- 2. Pseudo-cross-lingual: Here, a pseudo-language is created by mixing corpus from different languages. The cross-lingual context is used to learn the shared embedding space. [To what extent is this applicable?]
- 3. Cross-lingual training: parallel corpus used to train combined embeddings. Similar words will be mapped close to each other in same embedding space.
- 4. Joint optimization: Optimizing a combination of cross-lingual and monolingual losses.

For parallel data, we can have different types of data:

- 1. Word-aligned data
- 2. Sentence-aligned data
- 3. Document aligned data

- 4. Lexicon (translations between multiple language-tables)
- 5. No parallel data (only monolingual resources)

These are the ways to train word-embeddings (amongst others):

- word2vec variants
- skip-gram with negative sampling (SGNS)
- continuous bag-of-words (CBOW)
- GloVe matrix factorization approach

Learning a linear projection from monolingual embeddings and parallel data.

Projection via CCA (Canonical component analysis). A transformative matrix is learned for each language, the mapped space is the same space across different languages. Using 80% of projection vectors and using top k components generally yield highest performance.

Normalisation and orthogonal transport. Product similarity measure (correlation) can be seen as Wasserstein distance (because it measures how much data points are away from each other)  $c^T c$ . Vectors are normalised during training with unit vector. Because unit length, we can solve this by an orthogonal matrix. (Does any information go away? Maybe it's good to regualrize even further so that we know that the mapping is indeed possible?)

Max-margin intrudercs. Linear least squares leads to the hubness problem. Words tend to appear as nearest neighbors of many other words (-¿ perhaps some non-random sampling is best here? to move away). Use a ranking loss (sounds like optimal transport again...) Ranking loss is also used in wasserstein context.

Alignment based projection: Count number of times each word in source language is aligned with each word in target language document / sentence.

Adversarial autoencoder Autoencoder learns to re-create the embedding, while discriminator learns to maximize the projection

Orthogonal transformation, normalisation, mean centering Relation between the constraints is not clear. Starting with basic optimization objective, they add regualizers and loss functions as intuitively make sense. -¿ Dot-products need to be maintained after the mapping (as much as possible) constraint (i dont like)-¿ equal contribution to objective done by normalizing the embedding..-¿ two randomly selected words should in general not be similar

**Pseudo-Cross-Label** Captures interactions between words in different languages. Mostly skipped because brain capacity...

Joing training with 2 x 2000 dimensional input (one for each job-system)... can then translate each one in one latent space to another. How do we find parallel logic here, however..?

(-¿ Perhaps it makes sense to draw all the literature in the context of wassetstein, to do a probabilistic and principled analysis from there on.)

Observation (FRAGE: Frequency-Agnostic Word Representation): It is a feature that most-frequent words are embedded to different locations than non-most-embedded features. Xing et al argue there is a mismatch between the way it is learned, and the projection and loss-function. Tries to find a uniform theory across the three components.

If we use monolingual resources only, the common approach is

I focus on work which solely includes as input the bilingual word-embeddings. Because I want to devise an algorithm which can map from any "matrix column" to any other "matrix column" (i.e. token to token, or item to item), I will not include papers which take as input NLP sentences / corpora used to build the embeddings.

This work deals with some summary on how unsupervised bilingual lexicon matching was achieved in past papers.

I will give a short summary of papers sorted by year of appearance, and what additional contributions and observation were added since the last iteration.

Some terminiology before we start

**online data** implies that some sort of parallel corpora is available

**offline data** implies that we use pre-trained monolingual embeddings to arrive at a toen-translation task

The general goal of bilingual lexicon matching is to learn a shared embedding space where words possessing similar meanings are projected to nearby points.

Most existing methods focus on minimizing the distance between the two token-datasets using some (variation) of Wassterstein, Jensen-Shannon divergence, etc.

What are datasets that we can use for benchmarking

• MUSE dataset (as used in Zhou 2019 et al

•

Instead of mapping one system into another, we can also allow for joint-training of the word-embeddings.

(This resource is great!!! https://ruder.io/cross-lingual-embeddings/)

### 3.1 Discrete Methods

#### 3.1.1 Loss in Translation

Finding a non-orthogonal mapping is more effective, especially with distance dictionaries

Orthogonality preserves distances between word-pairs.

Euclidean closest nearest neighbor is taken as the translation pair.

Hubs are words that appear too frequently close to other words. This retrieval criterion suffers from the hubness problem.

Removing the loss creates a discrepancy between inference and training.

(-¿ Why should a loss-function be symmetric, not be symmetric?)

(-¿ especially with chinese. perhaps the non-orthogonality captures the ambigious tokens..?)

From the original problem of finding an orthogonal projection, we assimilate the auxiliary problem of

the CSLS criterion looks as follows:

$$CSLS(x,y) = -2\cos(x,y) + \frac{1}{k} \sum_{\mathbf{y}' \in \mathcal{N}_Y(\mathbf{x})} \cos(\mathbf{x}, \mathbf{y}') + \frac{1}{k} \sum_{\mathbf{x}' \in \mathcal{N}_X(\mathbf{y})} \cos(\mathbf{x}', \mathbf{y})$$
(3.1)

We then use the objective function of a relaxed

$$\min_{\mathbf{W} \in \mathcal{O}_d} \frac{1}{n} \sum_{i=1}^n -2\mathbf{x}_i^{\mathsf{T}} \mathbf{W}^{\mathsf{T}} \mathbf{y}_i \tag{3.2}$$

$$+\frac{1}{k} \sum_{\mathbf{y}_j \in \mathcal{N}_Y(\mathbf{W}\mathbf{x}_i)} \mathbf{x}_i^{\top} \mathbf{W}^{\top} \mathbf{y}_j$$
 (3.3)

$$+\frac{1}{k} \sum_{\mathbf{W} \mathbf{x}_i \in \mathcal{N}_X(\mathbf{y}_i)} \mathbf{x}_j^{\mathsf{T}} \mathbf{W}^{\mathsf{T}} \mathbf{y}_i \tag{3.4}$$

### 3.1.2 MUSE Facebook using a Min-Max objective

Mikolov et. al learn a mapping between the two sets of Y and X:

$$W^{\star} = \operatorname{argmin}_{W \in M_d(\mathbf{R})} \|WX - Y\|_F \tag{3.5}$$

Translation of any source token s to a target token t is defined as

$$t = \operatorname{argmax}_{t} \cos(Wx_{s}, y_{t})$$

(i.e. we just create a mapping between x and t using a linear interpolation matrix and take the maximal scalar product.

Mikolov et al do not observe improved loss when using more advanced neural networks..

Enforcing an orthogonality constraint, this boils down to the procrustes problem as given by

$$W^{\star} = \operatorname{argmin}_{W \in O_d(\mathbf{R})} \|WX - Y\|_F = UV^T$$
(3.6)

where U and  $V^T$  respectively the left and right Eigenvectors of  $YX^T$ .

We can create an adversarial unsupervised approach by taking random samples from

$$X = \{x_1, \dots, x_n\}$$
 and  $Y = \{y_1, \dots, y_m\}$ 

where the discriminator has to detect which of the two datasets the sample stemmed from. The "generator" (not really a generator) W is trained to prevent the discriminator from making accurate decisions. Specifically, this is modeled as a two-player game, where WX and Y should be as similar as possible.

As such, the discriminator objective is:

$$\mathcal{L}_D(\theta_D|W) = \tag{3.7}$$

$$= -\frac{1}{n} \sum_{i=1}^{n} \log P_{\theta_D}(\text{source} = 1|Wx)$$
 (3.8)

$$-\frac{1}{m}\sum_{i=1}^{n}\log P_{\theta_D}(\text{source} = 0|Wx)$$
(3.9)

and the generator objective is:

$$\mathcal{L}_W(W|\theta_D) = \tag{3.10}$$

$$= -\frac{1}{n} \sum_{i=1}^{n} \log P_{\theta_D} \text{ (source } = 0 | Wx_i \text{)}$$
 (3.11)

$$-\frac{1}{m}\sum_{i=1}^{m}\log P_{\theta_D} \text{ (source } = 1|y_i)$$
 (3.12)

Here, the discriminator is trained on the labels whether the word came from the first embedding, or the second embedding.

(-¿ Is there a mathematical way to simply this..?)

The resulting rotation matrix is not on-par with supervised approaches. As such, the authors propose to make use of the frequencies as additional information. As such, we use the most common words as anchors.

We then generate a dictionary from the two embeddings. Then we solve the procrustres to arrive at the final orthogonal matrix.

Use of the CSLS

### 3.2 Methods using Normalising Flows

# 3.2.1 Density Matching for Bilingual Word Embeddings (Zhou 2019)

Zhou et. al consider the problem by creating two smooth embedding spaces.

For simplicity, I will refer to embedding space  $\mathcal{X} \in \mathbf{R}^d$  as the source embeddings, and the embedding space  $\mathcal{Y} \in \mathbf{R}^d$  as the target embeddings.

We denote  $x_i$  and  $y_j$  as the individual words with respective embeddings  $x \in \mathcal{X}$  and  $y \in \mathcal{Y}$ .

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The general approach of the authors is as follows:

We learn two mapping functions  $f_{xy}: \mathcal{X} \to \mathcal{Y}, f_{yx}: \mathcal{Y} \to \mathcal{X}$ 

We maximize the density of data points in the source space. We use a loglikelihood term in the loss-function.

For monolingual embeddings, we define tractable density functions p(x), p(y).

We use gaussian embeddings, where each single gaussian module represents a single word.

$$p(x) \approx p(y) \approx GMM$$
 (3.13)

This implies that for both  $\forall x \in \mathcal{X}$  and  $\forall y \in \mathcal{Y}$  we have

$$p(x) = \sum_{i \in 1, \dots, N_x} \pi(x_i) \tilde{p}(x|x_i)$$
(3.14)

where the final probability term

$$\tilde{p}(x|x_i) = \mathcal{N}(x|x_i, \sigma_x^2 I)$$

describes the probability density of the single word-embedding, represented through a Gaussian in this case, and  $N_x$  defines the total number of components / words in dictionary C.

To initiate the embeddings, the authors use the word2vec (WAS IT WORD2VEC) point-vector as the mean of the Gaussian distribution, and use a fixed variance term  $\sigma^2$  for all Gaussians.

Furthermore, instead of using a uniform prior over the mixture weights  $\pi(x_i)$ , the authors use the observation that frequencies may give bettern information, as as such initiate the priors as the relative frequencies of the words, i.e.

$$\pi(x_i) = \frac{\operatorname{freq}(x_i)}{\sum_j \operatorname{freq}(x_j)}$$

.

This implies that the experiment is not executed in a manner where the Gaussian Mixutres are carefully chose, but rather by augmenting and generalizing the word2vec point vectors to Gaussian distributions. This has the effect that the resulting probability space becomes smooth (rather than a discrete space of point-vectors) [CITE??? SMOOTHNESS] Another assumption made is that a gaussian distribution is best-able to capture this mixtures.

Every training step obtains samples form the Gaussian mixture space.

The density calculation can be defined by using volume-preserving invertible transformations. Perhaps it could make sense to also apply convolutional operators?

To stabilize the unsupervised training, the author employs the following methods.

- 1. back-translation-loss to train both directions (this is conceptually similar to the cyclic dependency in GANs.
- 2. the author uses a weakly loss. identical text-strings in both languages are mapped to similar places in the space.
- 3. frequency of the words is used as a prior to the GMM prior weights.

The authors indicate robust training which is not strongly dependent on initialization

The training functions in such as a way as to increase the log-probability density. The general approach is:

1. A continuous vector x is sampled from the GMM. Specifically,

$$x_i \sim \pi(x_i), x \sim \tilde{p}(x|x_i) \tag{3.15}$$

2. We then apply  $f_{xy}$  to get the inferred target embedding  $y = f_{xy}(x)$ . Specifically, we define the mapping operation as:

$$f_{xy}(\cdot) = W_{xy} \cdot \tag{3.16}$$

We choose W as an invertible matrix, thus, the inverse function is defined as

$$x = f(y) = W_{xy}^{-1}y (3.17)$$

setting the volume of this operation to

$$J(f^{-1}(x)) = W_{xy}$$

which is used for the log-density calculations. Specifically

$$\log p(x; W_{xy}) = \log p(y) + \log |\det(W_{xy})| \tag{3.18}$$

3. We then maximize the log-likelihood, which equivalent to minimizing the expectation of the KL-divergence between the prior and modal distributions.

$$KL(p(x)||p(x;W_{xy}) \tag{3.19}$$

We arrive at the conditional unsupervised mapping-loss function of (incorporating the observation that each mixture represents a single word token)

$$\mathcal{L}_{xy} = \mathbf{E}x \sim p(x) \left[ \log p(y) + \log \left| \det(W_{xy}) \right| \right]$$
 (3.20)

Where the first term is the log-density of the data, and the second term describes a regualizer (prior) over the sample model which we use for the mapping. This equation corresponds to the normalising flow which is volume-preserving.

The maximum log-likelihood is proportional to the minimum expectation of the KL-divergence between the prior and modal distribution. (DAFUQ IS

# 3.3 Methods viewing this problem as an Optimal Transport (OT) Problem

# 3.3.1 Gromov-Wasserstein Alignment of Word Embedding Spaces

The problem of aligning word embedding spaces can also be seen as an Optimal Transport  $(\mathbf{OT})$  problem.

Some work in using the Gromov-Wasserstein distance function has been proposed (CITE Jaakkola).

The authors argue that the supervised case is

```
Algorithm 1: How to write algorithms

Result: Write here the result
initialization;
while While condition do

instructions;
if condition then

instructions1;
instructions2;
else

instructions3;
end
end
```

Supervised Case: Procrustes

Assuming we have

such that

$$T(x^{(i)}) \approx y^{j}$$

where  $w_i^y$  is a translation of  $w_i^x$ .

Let X and Y be the matrices with the column-vectors describing the word-embeddings.

We can then find T by solving the equation

$$\min_{T \in \mathcal{F}} \|X - T(Y)\|_F^2 \tag{3.21}$$

The quality of the resulting alignment depends on how well we can find T, as well as on the choice of the space  $\mathcal{F}$ .

T orthogonal matrix is a popular choice, results iin the orthogonal Procrustes problem

$$\min_{P \in O(n)} \|X - PY\|_F^2 \tag{3.22}$$

with  $O(n) = P \in \mathbf{R}^{nxn} | P^T P = I$ .

In this case, we can find a closed-form solution in terms of the SVD of PX and Y. Given an SVD of  $U\Sigma V^T$  of  $XY^T$ , the solution is given by  $P^* = UV^T$ .

Because the loss must be able to be measured, it only solves the supervised problem.

#### Unsupervised Maps: Optimal Transport

Optimal transport formalizes the problem of finding a minimum cost mapping between two point sets, viewed as discrete distributions.

We assume two empirical delta-distributions over embeddings (which are as such equivalent to point-vectors).

#### WHAT WAS THIS SYMBOL AGAIN WHICH LOOKS LIKE A V

$$\mu = \sum_{i=1}^{n} p_i \delta_{x(i)}, v = \sum_{j=1}^{m} q_j \delta_{y(i)}$$
(3.23)

where  $p_i$  and  $q_j$  can be considered as the prior distributions, usually set to the uniform distribution or word-frequency.

The transport problem is then realized as

$$\inf_{T} \{ c(x, T(x)) d\mu(x) | T\mu = v \}$$
(3.24)

where the cost is usually simply defined as ||x - T(x)|| and  $T\mu = v$  implies that the source points must exactly map to the target points.

We use Kantorovich's relaxed formulation, where the set of transportation plans is a polytope:

$$\prod(p,q) = \{ \Gamma \in \mathbf{R}_{+}^{n \times m} | \Gamma \mathbb{1}_{n} = p, \Gamma^{T} \mathbb{1}_{n} = q \}$$
 (3.25)

In the end, this is just a matrix-factorization problem of the type, including a regularization on the total entropy H of the loss:

$$\min_{\Gamma \in \prod(p,q)} \langle \Gamma, C \rangle - \lambda H(\gamma) \tag{3.26}$$

The authors propose "Transport across unaligned spaces", addressing the need for vectors across two spaces. Instead of samples between two spaces, we now focus on the distance of the two spaces as a whole.

Upsides:

- 1. Discrete Vectors
- 2. Some formulation of "minimal" rotation

Shortcoming:

1. An overall single rotation will most likely not solve this issue

How do we now proceed with using the distances between two functions as the main information to map the two words.

### 3.3.2 Graph translation using normalizing flows

This follows from the observation that any document x token matrix can be modelled as a graph.

# 3.3.3 Using monolingual word-embeddings to map one to another

#### Cross-lingual word-mappings (Artetxe 2018)

Extension of work for more distant language pairs for looking at tokens.

Artetxe does the following series of oprations

- 1. Embedding normalization: Normalizes the length of the word-embeddings and then mean-center each dimension.
- 2. Fully unsupervised initialization: Uses distances to measure similarity  $(M_X = X^T X \text{ and } M_Y = Y^T Y)$ . Common alternatives include finding common letters and numerals across two words. Distances should be isometric up to some permutation. Main assumption: Distance between the two spaces must be equivalent, else unsupervised is not possible! Compute sorted $(\sqrt{M_X}, \sqrt{M_Y})$ . This didnt quite work, so we randomly keep some elements with probability p. This allows for easier matching because there are many ways to generate this.. [This is basically like batch-sampling for training, no? This noise (if batchsize is chosen well), should be able to well-solve this]. We only focus on the top 20000 words in each language for simplicity. Instead of kNN, we use the CLSL retrieval. Use bidirectional learning to avoid local optima where one word is not available in another language.

3. Re-weighting [What??] by whitening the cross-correlation matrix. Ultimately, this again also just performs on distances.

$$USV^T = X^TDZ$$

looking at  $W_X = US^{\frac{1}{2}}$ , then we whiten by

$$U^T(X^TX)^{\frac{1}{2}}U$$

.

CSLS often used because of the "hubness" problem [Perhaps we can copy all these oprations in a probabilistic context with Wassterstein?]

## Bilingual Lexicon Induction through Unsupervised Machine Translation (Artetxe 2019)

Uses machine translation to build a lexicon induction model.

Input to algorithm:

- 1. Monolingual corpora fasttext through vecmap and the corpora used to train.
- 2. 400'000 most common bigrams and 400'000 most common tri-grams.
- 3. To find most similar pairs, instead of using kNN CLSL, proposed is

$$\phi(\bar{f}|\bar{e}) = \frac{\exp(\cos(\bar{e}, \bar{f})/\tau)}{\sum_{\bar{f}'} \exp(\cos(\bar{e}, \bar{f}')/\tau)}$$

to extract the top 100 sentences 5-gram language model.

- 4. The proposed method increases this by at least 10 percent points (P@1).
- 5. underlying language model uses 2000 sentences from each monolingual corpora, a language modeling loss, and combines this with a cyclic consistency loss.

- 3.4 Gaussian Embeddings and Token matching in other applications
- 3.4.1 Creating Gaussian Embeddings to represent Graphs (DEEP GAUSSIAN EMBEDDING OF GRAPHS: UNSUPERVISED INDUCTIVE LEARNING VIA RANKING)

# Analysis of the current state of the art

For the problem statement to be - generalizable to other datasets - unsupervised (perhaps some [noisy] inference on starting points possible..)

Some challenges that are faced are

- Hubness problem (MUSE, and CSLS)

(https://ruder.io/cross-lingual-embeddings/)

Functional modeling may be difficult to model. It may be able to model the relationship between words well, but it will not be able to distinguish the context or actual usage of the word within a sentence.

Word-order is completely ignored. [will probably ignore this aspect, simply because we dont want sequence-based approaches for generalizability] This may lead to different types of training..

Compositionality: word-representations can not be generalized further to sentence-generation and document-generation. Simply adding word-embeddings cannot result in a meaningful

Polysemy: a word may have multiple meanings. in a cross-lingual embedding

space, this feeling is amplified. there's some work in multi-sense embedding. this should enable to capture more fine-grained embeddings embeddings.

Feasability: Maybe not possible?

Upadhyay et al. argue that the choice of data is more important than the actual algorithm.

Include an analysis of all loss functions.

Include a list of assumptions made.

Include a list of observations made

Ablation study.

Unnormalized embeddings can modify training in that the loss function puts more weights to the high-occurring vectors.

Maybe Mikolov 2013b did not do proper hyperparameter tuning to check if deeper layers improve this performance..?

### Our Method

#### Ideas:

- 1. Generally, deviate more from first proposed paper.
- 2. (like EM) two-step training where dictionary is learned and then the underlying soft probability distribution (using parametrized models)
- 3. Instead of cross-correlation, look for some distance in the probability space which makes sense? Test multiple ones, and possibly arrive at a prior which is interesting
- 4. order of phrasing is different in korean / turkish to german
- 5. Enforce some sort of probability density on top?
- 6. Go deeper perhaps?
- 7. Generally collect a set of constraints and observations that seem to work, then combine them in a principled approach
- 8. Normalising flows seem invertible. Can we not just use  $f_{xy} = f_{yx}$ .
- 9. AutoML on normalising flows..?
- 10. Leave out individual variables
- 11. Jointly train both embedding and normalising flow

- 12. After two possible ways of embeddings are learned, can we use the weights of the neural networ (i.e. word2vec") to find a mapping that goes from one space into another? (... yes, we should be able to using least squares projection...)
- 13. Work on generalizability of this on other datasets (such as ESCO to australia, or Amazon Austria to Amazon Germany)
- 14. Regualizer that adds up to respect some corpus / frequency matrix properties...
- 15. Hubness problem by enforcing a prior to the dataset s.t. the embeddings are uniformly distributed in space?
- 16. Probability density matching using optimal transport not on Euclidean space, but on a Riemann manifold / hyperbolic space?

### Further Work

Can be used for more unstructured data, like graphs.

Evaluation

Conclusion