

# EBLearn: Open-Source Energy-Based Learning in C++

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## Abstract

*Energy-based learning (EBL) is a general framework to describe supervised and unsupervised training methods for probabilistic and non-probabilistic factor graphs. An energy-based model associates a scalar energy to configurations of inputs, outputs, and latent variables. Learning machines can be constructed by assembling modules and loss functions. Gradient-based learning procedures are easily implemented through semi-automatic differentiation of complex models constructed by assembling predefined modules. We introduce an open-source and cross-platform C++ library called EBLearn<sup>1</sup> to enable the construction of energy-based learning models. EBLearn is composed of two major components, **libidx**: an efficient and flexible multi-dimensional tensor library, and **libelearn**: an object-oriented library of trainable modules and learning algorithms. The latter has facilities for such models as convolutional networks, as well as for image processing. It also provides graphical display functions.*

## 1. Introduction

Energy-based learning [11] (EBL) provides a unified framework for probabilistic and non-probabilistic machine learning methods. Energy based learning models have been successfully used in a number of applications such as object recognition [6, 16, 7], outdoor unstructured robotics vision [5], signal processing [13], time series modeling [12], manifold learning [1, 4], financial prediction [2], document recognition [10], natural language processing [3], unsupervised learning of feature hierarchies [18, 15, 8] and text classification [17]. **Inference** in energy based models for a given input  $X$  is performed by finding a configuration of output  $Y$  and latent variables  $Z$  that minimize an **energy function**  $E(X, Y)$ . For **Learning**, the energy function is parameterized by a parameter vector  $W$ . Learning is performed by minimizing a suitable **loss functional** with respect to  $W$ . For a given input  $X$ , the purpose of learning is to shape the energy surface so that corresponding

desired output configuration(s) have lower energy than all other configurations.

A simplified EBL model where  $X$  represents the inputs to the system (a sample from the data),  $Y$  represents the output variable of the model, the inference process is formulated as:

$$Y^* = \arg \min_{Y \in \mathcal{Y}} E(X, Y) \quad (1)$$

where  $\mathcal{Y}$  is a set of possible outputs. When  $\mathcal{Y}$  is a discrete set with few elements, exhaustive search can be used, but when  $\mathcal{Y}$  has high cardinality or is a continuous set, suitable minimization algorithms must be employed. In unsupervised scenarios, the energy function has no observation  $X$ , and the model  $E(Y)$  simply indicates whether a particular  $Y$  is similar to training samples (low energies) or dissimilar (higher energies). Different types of problems can be formulated under this model, such as classification, detection, regression, ranking, density estimation, clustering, and others.

Probabilistic models are a special case of EBL in which the energy is integrable with respect to  $Y$ . The distribution can be obtained using the Gibbs formula:

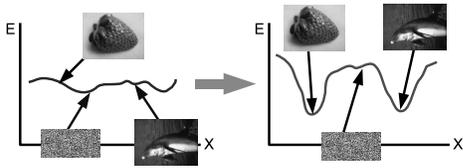
$$P(Y|X) = \frac{e^{-\beta E(X, Y)}}{\int_y e^{-\beta E(X, y)}} \quad (2)$$

where  $\beta$  is a positive constant, and the denominator is called the *partition function* (variable  $X$  is simply dropped for unsupervised scenarios).

When latent variables  $Z$  are present, they can be minimized over or marginalized over. With minimization, the energy function is simply redefined as  $E(X, Y) = \min_{Z \in \mathcal{Z}} E(X, Y, Z)$ , and with marginalization as  $E(X, Y) = -\frac{1}{\beta} \log \int_z e^{-\frac{1}{\beta} E(X, Y, Z)}$ . If necessary, this integral can be approximated through sampling or using variational methods.

Given a training set  $\{(X^1, Y^1), \dots, (X^P, Y^P)\}$ , training a model consists in shaping the energy surface  $E(W, X, \cdot)$  (expressed as a function of  $Y$ ) parametrized by  $W \in \mathcal{W}$  by minimizing a suitable loss functional with respect to  $W$ ,

<sup>1</sup><http://ebllearn.sourceforge.net>



**Figure 1. Energy Based Training:** Before training, the energy surface produced by an *EBL* model is not distinctive around training data. After training, the energy surface is shaped lower around training data.

averaged over the training set:

$$W^* = \min_{W \in \mathcal{W}} \frac{1}{P} \sum_i \mathcal{L}(E(W, X^i, \cdot), Y^i) \quad (3)$$

The loss can more simply be expressed as a function of  $W$  :  $\frac{1}{P} \sum_i \mathcal{L}(W, X^i, Y^i)$ . The purpose of loss is to measure whether the “correct” output  $Y^i$  for a given  $X^i$  has lower energy than all other outputs. As a result, the system produces lower energy values for regions around observed  $Y$  values, as shown in Figure 1. Building an *EBL* model can therefore be achieved by designing:

1. the *architecture* of the  $E(W, X, Y)$ ,
2. the *inference algorithm* that will be used to infer outputs  $Y$  that minimize  $E(W, X, Y)$  for a given  $X$  and fixed  $W$ ,
3. the *loss function*  $\mathcal{L}(E(W, X, \cdot), Y)$  and
4. the *learning algorithm* that will be used to find the best  $W$  that minimizes the loss averaged over a training set.

### 1.1 Loss Functions for Energy-Based Learning

In energy based learning designing proper loss functions for different types of architectures is required to avoid trivial solutions where the energy surface becomes flat. Many different loss functions have been proposed in machine learning literature and in this section we formulate several popular loss functions in energy based learning framework.

**Energy Loss:** simply means that the loss function per each sample is equal to the energy function:  $\mathcal{L}_{en}(W, X^i, Y^i) = E(W, X^i, Y^i)$ . Energy loss is the simplest possible loss function that can be used in wide variety of cases. The energy loss will ensure that the energy surface is *pulled down* around desired data, but it does not guarantee that the energy surface is *pulled up* at all other locations. This might lead to situations where the energy for all samples becomes constant. Linear regression using mean squared error (MSE) can be formulated using energy loss:  $\mathcal{L}(W, X^i, Y^i) = E(W, X^i, Y^i) = \|Y^i - WX^i\|^2$  where  $W$  is the matrix of parameters with respect to which  $\mathcal{L}$  is minimized.

**Perceptron Loss:** defined per each sample and over all possible configurations of outputs takes the following form

$$\mathcal{L}_{perceptron}(W, X^i, \mathcal{Y}) = E(W, X^i, Y^i) - \min_{y \in \mathcal{Y}} E(W, X^i, y) \quad (4)$$

Perceptron loss pulls down the energy of the correct configuration (first term) and it pulls up the energy of current prediction. When the machine prediction is correct, the loss

is equal to zero and positive otherwise. One can see that, this loss function does not enforce a margin between correct and incorrect configurations, thus might lead to almost flat surfaces.

**Negative Log-Likelihood Loss:** is suitable to train a model to produce probability estimates for  $P(Y|X)$ :

$$\mathcal{L}_{nll} = E(W, X^i, Y^i) + \frac{1}{\beta} \log \int_y e^{-\beta E(W, X^i, y)} \quad (5)$$

where  $\beta \in \mathcal{R}^+$ . As in maximum likelihood solutions for probabilistic models, the integral in the second term of equation 5 might be intractable to compute or might not have an analytical solution in most cases. Approximate solutions to this integral can be obtained by approximate analytical solutions, sampling methods and variational methods.

**Contrastive Loss:** enforces a gap between the energy of correct answer  $Y^i$  and the energy of the *most offending incorrect answer*  $\bar{Y}^i$ , defined as the *wrong* answer with the smallest energy. The most commonly used contrastive loss is the hinge loss:

$$\mathcal{L}_{hinge} = \max(0, m + E(W, X^i, Y^i) - E(W, X^i, \bar{Y}^i)) \quad (6)$$

where  $m$  is the positive margin. The model is updated whenever the energy of incorrect answer is less than  $m$  larger than energy of correct answer.

### 1.2 Modules in Energy-Based Learning

Until now, we have stated that the energy function is parametrized by  $W$ . In most supervised scenarios, this is conveniently achieved by using a functional module  $G_W(X)$  which maps samples from input space to output space, and by capping it with a “distance” module that measure the discrepancy between  $G_W(X)$  and the output  $Y$ . A very simple example of a functional module is a single matrix multiplication that projects the input along the space defined by its columns. In more complicated cases it can be combination of linear and nonlinear functions. Common modules include linear modules as explained, simple nonlinear modules that are applied on each element of the input state independently, like sigmoid functions, and convolutional modules that are very similar to linear modules, but are applied as convolution operations on input image maps. The loss function  $\mathcal{L}$  is minimized with respect to the parameters  $W$  of the functional modules.

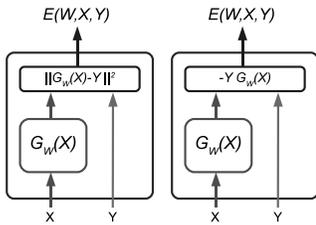
### 1.3 Architectures for Energy-Based Learning

In this section we provide *EBL* models for some widely used learning algorithms.

**Regression:** is one of the most common algorithms used in machine learning. A regressor model (Fig. 2) can be obtained by using a squared error energy function

$$E(W, X^i, Y^i) = \frac{1}{2} \|G_W(X^i) - Y^i\|^2 \quad (7)$$

together with energy loss. When  $G_W$  is a linear operator, this model becomes equivalent to solving the least squares problem.



**Figure 2. Architectures for Energy Based Model:** Left: Regression can be formulated by using a squared distance energy function combined with energy loss and a module  $G_w(X)$ . Right: Two class classification can be formulated similarly using perceptron loss.

**Two-Class Classification:** can be formulated using a simple energy function as shown in Figure 2.

$$E(W, X^i, Y^i) = -Y^i G_W(X^i) \quad (8)$$

Any of the perceptron loss, hinge loss or negative log likelihood loss can be used with this energy function to solve two class classification problems.

**Multi-Class Classification:** can be done by replacing the energy function in Figure 2 with

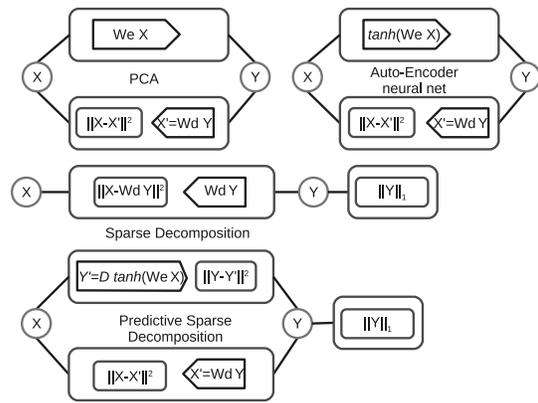
$$E(W, X^i, Y^i) = \sum_{k=1}^c \delta(Y^i - k) g_k \quad (9)$$

where  $\delta(u)$  is Kronecker delta function and  $G_W = [g_1 \ g_2 \ \dots \ g_c]$ . As with the two-class classification problem, perceptron, hinge and negative log likelihood loss functions can be used.

One can imagine that complicated architectures can be built by combining several functional modules and energy functions as long as the combined energy function can be minimized with respect to the desired outputs  $Y$  and the final loss function can be minimized with respect to the parameters  $W$ .

It has to be noted that, any factor graph can also be modeled using energy based learning. In a simple factor graph an observed state is connected to a latent state through a factor node which models the constraint in between two states. The combination of two states are assigned high likelihood under the compatibility constraint defined by the factor node. One can also separate the dependency constraints between two states into two directional factor nodes representing the compatibility between one node and transformation of the other.

Most unsupervised learning algorithms can be modeled in this framework. In Figure 3, we show several common unsupervised learning algorithms. *PCA* is a linear model, where the transformation from observed input  $X$  to latent representation  $Y$  is a linear projection. Accordingly, transformation from  $Y$  to input space  $X'$  is also a linear projection. The model has to be trained under the reconstruction compatibility constraint such that transformed reconstruction  $X'$  has to minimize the squared reconstruction error between original input  $X$  and projected input  $X'$ . *Auto-encoder neural networks* are very similar to *PCA*, except the projection from input  $X$  to latent variable  $Y$  is



**Figure 3. Modeling Unsupervised Learning Algorithms:** Many unsupervised learning algorithms can be represented in the factor graph model and trained with energy based learning.

non-linear. *Sparse Decomposition* [14] is a uni-directional model where, there is no direct projection from input  $X$  to latent variable  $Y$ . Instead, for each input  $X$ , the system has to carry out an optimization process to infer latent representation  $Y$ . In addition to reconstruction compatibility constraint, the latent representation has to minimize the  $L_1$  norm constraint. *Predictive Sparse Decomposition* [9] is an extension to sparse coding models, where a nonlinear predictor function is also trained to infer latent variable  $Y$  from input  $X$  without requiring any optimization process.

## 2. libidx: Tensor Descriptors and Operators

The *idx* library (or *libidx*) provides convenient and efficient tensor (multi-dimensional arrays) manipulations, used as a basis for the *eblearn* library. There are three main components to the library: tensor descriptors and iterators, content operators and image-specific operators.

### 2.1 idx: Tensor Descriptors

The *idx* class can be thought as a tensor pointer to a chunk of memory (or an *srg* class, *srg* standing for storage), and multiple *idx* can point to different subsets of that memory. A *idx* describes a tensor by its storage (*srg*), its offset on that storage, its number of dimensions (or order) and size and stride of each dimension. Declaring a new *idx* will allocate and initialize to zero a new storage of size and type specified to the constructor and class template. Here for example, we create a 3-dimensional tensor with double precision of size 32x32x3, which could also be interpreted as a 32x32 RGB image:

```
idx<double> t(32, 32, 3);
```

Being relatively cheap memory and computationally wise, an *idx* can be manipulated efficiently like a tensor pointer without affecting the actual tensor memory. For example, the user can select at virtually no cost the entire (2D) slice at position  $p$  of the  $d^{th}$  dimension of a tensor with *select* or narrow dimension  $d$  to size  $s$  starting at position  $p$  to create a 3D subset of the 3D tensor  $t$ :

```
idx<double> slice = t.select(d, p);
idx<double> subset3d = t.narrow(d, s, p);
```

Memory management is facilitated by the self garbage collection of *srg* (reference counters), i.e. a storage is automatically destroyed once no more *idx* point to it.

## 2.2 idx loops: Tensor Iterators

While tensor elements can be accessed individually via set and get methods, one will often need to loop over entire dimensions or entire tensors. Looping macros are provided for that effect, e.g. the *idx\_aloop2* macro loops over all elements of 2 tensors while the *idx\_bloop3* macro loops over the first dimensions of 3 tensors. For each tensor to be iterated, one must specify a temporary name for the new lower-order tensor, the original tensor and its type. For example, to compute the sum of multiple tensors one could write:

```
idx<double> td3d(32, 32, 3);
idx<int>    ti2d(32, 32);
int total = 0;
idx_aloop1(td0d, td3d, double) total += td0d.get();
idx_aloop1(ti0d, ti2d,    int) total += ti0d.get();
```

## 2.3 Tensor Operators: Content Manipulations

While *idx* descriptors are inexpensive pointers, *idx* content operators work with the tensor data yielding more expensive operations. We now describe a few important operators among others.

**Copy operator**, copy the content of *d1* to *f2* (same dimensions) with automatic type casting:

```
idx<double> d1(32, 32, 3);
idx<float>  f1(32, 32, 3);
idx_copy(d1, f1);
```

**I/O operators**, saving and loading tensors :

```
save_matrix(f1, "im.mat");
idx<float> f2 = load_matrix<float>("im.mat");
```

**Product operators**, the dot of two tensors (seen as vectors) or the matrix-vector multiplication:

```
float dot = idx_dot(f1, f2);
idx<float> f3(32, 16), f4(16), f5(32);
idx_m2dotm1(f3, f4, f5);
```

## 2.4 Tensor-based Image Operators

Images can be seen and manipulated as tensors. We present here some key image-specific operators.

**Image I/O operators**, load, save or resize images:

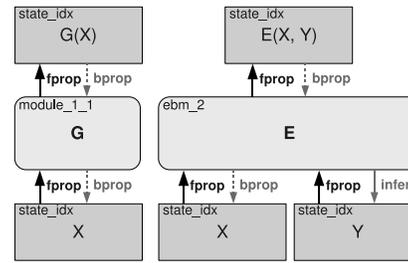
```
idx<float> im = load_image<float>("im.jpg");
save_image(im, "im.png");
im = image_resize(im, 16, 16);
```

**Image filtering**, local and global normalization:

```
idx<float> im2 = idx_copy<float>(im);
image_global_normalization(im);
image_local_normalization(im, im2, 9);
```

## 3. libblearn: Energy-Based Learning

The *libblearn* library is mainly constituted of modules of two types: *module\_1\_1* which takes 1 input and produces 1 output and *module\_2\_1* with 2 inputs and 1 output (Fig. 4). In particular for an *EBL* model, we derive the *ebm\_2* module from *module\_2\_1* to output an energy from its 2 inputs. Those two models are the basis for all modules in the library which can be combined into complex models. Each module implements the *fprop* (forward



**Figure 4.** The two basic types of modules. *module\_1\_1* (left) has 1 input and 1 output and *ebm\_2* has 2 inputs and 1 energy output. *state\_idx* store temporary results of calls to *fprop* (training and inference), *bprop* and *bbprop* (training only), and *infer2* (inference only).

propagation), *bprop* (backward propagation) and *bbprop* (usually back propagation of second derivatives) methods. *module\_2\_1* also implements the *infer2* method (infer second input). While *bprop* and *bbprop* methods are only used during training and *infer2* during inference, the *fprop* method is used during both phases. Intermediate results of *fprop*, *bprop* and *bbprop* calls are held in between modules in *state\_idx* objects.

## 3.1 Example: a Vision System

In this example, we build, train and execute in a few lines of code a convolutional neural network capable of object recognition as in [6]. The machine is a stack of convolution, subsampling and fully-connected modules (Fig. 5) taking input images to be classified as one of five categories. We now describe the construction of that system:

1. Build  $E(W, X, Y)$ , using a lenet7 neural network as  $G_W(X)$  and an euclidean energy module as energy function (see Fig. 5 for corresponding architecture):

```
parameter<double> W;
layers<double> 17(true);
17.add(new convolution_layer(W, 5, 5, 1, 1, ftbl(1, 8)));
17.add(new subsampling_layer(W, 4, 4, 4, 4, 8));
17.add(new convolution_layer(W, 6, 6, 1, 1, rtbl(8, 24, 4)));
17.add(new subsampling_layer(W, 3, 3, 3, 3, 24));
17.add(new convolution_layer(W, 6, 6, 1, 1, ftbl(24, 100)));
17.add(new full_layer(W, 100, 5));
euclidean_energy<double, int> eenergy;
machine<double, int> E(17, energy);
```

where the numbers are kernels sizes, strides and output sizes. *ftbl* and *rtbl* functions provide full or sparse random connections between layers. A shorter equivalent:

```
parameter<double> W;
lenet7<double> 17(W);
euclidean_machine<double, int> E(17);
```

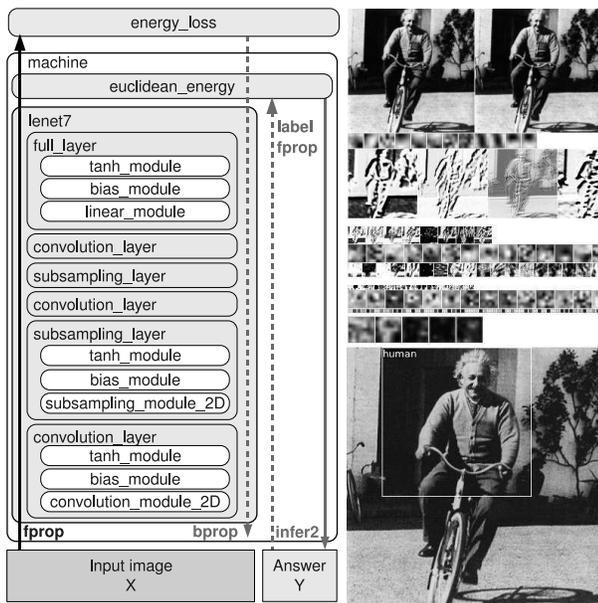
2. Build the loss and the trainer:

```
energy_loss eloss;
supervised_trainer<double, int> trainer(E, eloss);
```

3. Train the system with the NORB dataset [6] and a learning rate of 0.0001:

```
norb_datasource ds("/datasets/norb");
gd_param p(0.0001);
trainer.train(ds, p);
```

4. Execute the system:



**Figure 5.** A vision architecture (left) and a corresponding object detection example (right). This *machine* combines an euclidean energy with a 6-layer neural (*lenet7*). During both training and inference, the machine is first evaluated with *fprop* (bottom-up). Then for training only (dashed lines), an energy loss module comes on top of the machine and uses the training label to back-propagate (*bprop*) the gradient of the loss through the entire machine. During inference however, the loss module is not used and the answer is inferred via *infer2* following an *fprop*. The example (right) uses a machine trained with the 5-class NORB dataset (animal, car, human, plane, truck) and correctly classify Einstein as human. Internal states and kernels are represented top-down from input to output.

```
idx<double> image = load_image<double>("im.jpg");
state_idx<double> input(image);
int answer = E.infer2(input);
detector d(E, scales_number);
vector<bbox> answers = d.fprop(input);
```

## 4. Complementary Tools

In addition to core libraries *libelearn* and *libidx*, a set of tools are provided for display of tensors (*libidxgui*) and learning classes (*libelearnui*, see Fig. 5), for dataset generation and visualization (*libelearntools*) and for unit-testing (*tester*). Some demonstration projects are also available.

## 5. Conclusion

Energy based learning has been used in many different contexts of machine learning and provide a very efficient and flexible framework. We have showed that many supervised and unsupervised learning algorithms and factor graphs can be modeled using energy based learning framework. Inference and learning processes are formulated for many popular problems. More importantly, in this work we have presented an open source machine learning library (*EBLearn*) that can be used to built energy-based learning models. *EBLearn* is developed using C++ programming language for maximum portability and flexibility. We have also shown several code examples on how to use *EBLearn*

and several additional graphical display methods and image processing methods that are also included. With this work we introduce the availability of an open source machine learning library that can be used to train supervised, semi-supervised and unsupervised models. We believe this library contains one of the most extensive collection of machine learning algorithms.

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