# Spatially Adaptive Inference with Stochastic Feature Sampling and Interpolation

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Abstract. In the feature maps of CNNs, there commonly exists considerable spatial redundancy that leads to much repetitive processing. Towards reducing this superfluous computation, we propose to compute features only at sparsely sampled locations, which are probabilistically chosen according to activation responses, and then densely reconstruct the feature map with an efficient interpolation procedure. With this sampling-interpolation scheme, our network avoids expending computation on spatial locations that can be effectively interpolated, while being robust to activation prediction errors through broadly distributed sampling. A technical challenge of this sampling-based approach is that the binary decision variables for representing discrete sampling locations are non-differentiable, making them incompatible with backpropagation. To circumvent this issue, we make use of a reparameterization trick based on the Gumbel-Softmax distribution, with which backpropagation can iterate these variables towards binary values. The presented network is experimentally shown to save substantial computation while maintaining accuracy over a variety of computer vision tasks.

**Keywords:** Sparse convolution, sparse sampling, feature interpolation

#### 1 Introduction

On many computer vision tasks, significant improvements in accuracy have been achieved through increasing model capacity in convolutional neural networks (CNNs) [11, 32]. These gains, however, come at a cost of greater processing that can hinder deployment on resource-limited devices. Towards greater practicality of deep models, much attention has been focused on reducing CNN computation.

A common approach to this problem is to prune weights and neurons that are not needed to maintain the network's performance [19, 10, 9, 34, 13, 20, 28, 37].

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Fig. 1. (a) Input image. (b) Deterministic sampling for efficient inference. (c) Content-aware stochastic sampling by our method (without *Grid Prior* for better viewing), which yields the same detection accuracy as (b) but with less overall computation. (d) Content-aware stochastic sampling with the same overall computation as (b) but with better detection accuracy.

Orthogonal to these architectural changes are methods that eliminate computation at inference time conditioned on the input. These techniques are typically based on feature map sparsity, where the locations of zero-valued activations are predicted so that the computation at those positions can be skipped [6, 30, 1]. As illustrated in Fig. 1(b), this approach deterministically samples predicted foreground areas while avoiding computational expenditure on the background.

In this paper, we seek a more efficient allocation of computation over a feature map that takes advantage of its spatial redundancy<sup>5</sup>. Rather than exhausting all the computation on areas estimated to have the most significant activation, our approach is to treat the predicted activation map as a probability field, stochastically sample a sparse set of locations based on their probabilities, and then interpolate the features at these samples to reconstruct the rest of the feature map. This strategy avoids wasting computation at locations where the features can simply be interpolated, and it allows feature computation to expand into presumed low-activation regions, which can compensate for errors in activation prediction. With the generated sampling distributions shown in Fig. 1(c) and (d), this sampling-interpolation approach can reduce the computation needed to match the accuracy of feature map sparsity methods, or alternatively increase accuracy through more comprehensive sampling of the feature map while maintaining the same level of overall computation.

To identify sparse points for interpolation, our network trains a content-aware stochastic sampling module that produces a binary sampling mask over the activations. Due to the inability to backpropagate through networks containing binary variables, we employ a reparameterization trick where the non-differentiable mask values are replaced by differentiable samples from a Gumbel-Softmax distribution, which can be smoothly annealed into binary values during training [17]. The module learns to spatially adjust the sampling density according to predicted activations, and the interpolation is performed with a kernel whose parameters are jointly learned with the sampling module. To aid in interpolation of areas far from the content-aware samples, very sparse uniform samples over the feature map are added before interpolation, which we refer to as a grid prior.

<sup>&</sup>lt;sup>5</sup> We note that CNNs commonly capitalize on spatial redundancy by downsampling input image resolutions or employing strides in convolution.

With this content-based sampling-interpolation approach, our network obtains appreciable reductions in computation without much loss in accuracy on COCO object detection, Cityscapes semantic segmentation, and ImageNet classification. An extensive ablation study validates the sampling and interpolation components of our algorithm, and comparisons to related techniques show that the proposed method provides a superior FLOPs-accuracy tradeoff for object detection and semantic segmentation, and comparable performance for image classification.

# 2 Related work

In this section, we briefly review related approaches for reducing computation in convolutional neural networks.

Model pruning A widely investigated approach for improving network efficiency is to remove connections or filters that are unimportant for achieving high performance. The importance of these network elements has been approximated in various ways, including by connection weight magnitudes [10,9], filter norms [20, 34, 12], and filter redundancy within a layer [13]. To reflect network sensitivity to these elements, importance has also been measured based on their effects on the loss [19, 28] or the reconstruction error of the final response layer [37] when removing them. Alternatively, sparsity learning techniques identify what to prune in conjunction with network training, through constraints that zero out some filters [34], cause some filters to become redundant and removable [5], scale some filter or block outputs to zero [22], or sparsify batch normalization scaling factors [25, 35]. Model pruning techniques as well as other architecture-based acceleration schemes, such as low-rank factorizations of convolutional filters [16] and knowledge distillation of networks [15], are orthogonal to our approach and could potentially be employed in a complementary manner. Early stopping Rather than prune network elements, early stopping techniques reduce computation by skipping the processing at later stages whenever it is deemed to be unnecessary. In [7], an adaptive number of ResNet layers are skipped within a residual block for unimportant regions in object classification. The skipping mechanism is controlled by halting scores predicted at branches to the output of each residual unit. In [21], a deep model for semantic segmentation is turned into a cascade of sub-models where earlier sub-models handle easy regions and harder cases are progressively fed forward to the next sub-model for further processing. In [18], various predefined downsampling configurations are randomly used during training, and the appropriate configuration is applied according to the computation budget during inference. Like our method, these techniques spatially adapt the processing to the input content. However, they process all spatial positions at least to some degree, which limits the achievable computational savings.

Activation sparsity The activations of rectified linear units (ReLUs) are commonly sparse. This property has been exploited for network acceleration by excluding the zero values from subsequent convolutions [31, 29]. This approach has been extended by estimating the activation sparsity and skipping the computation for predicted insignificant activations. The sparsity has been predicted from

prior knowledge of road and sidewalk locations in autonomous driving applications [30], from model-predicted foreground masks at low resolution [30], from a small auxiliary layer that supplements each convolutional layer [6], and from a highly quantized version of the convolutional layer [1]. Our work instead reconstructs activation maps by interpolation from a sparse set of samples selected in a content-aware fashion, thus avoiding computation at locations where features can be easily reconstructed. Moreover, our probabilistic sampling distributes computation among feature map locations with varying levels of predicted activation, providing greater robustness to activation prediction errors.

Sparse sampling To reduce processing cost, PerforatedCNNs compute only sparse samples of a convolutional layer's outputs and interpolate the remaining values [8]. The sampling follows a predefined pattern, and the interpolation is done by nearest neighbors. Our method also takes a sparse sampling and interpolation approach, but in contrast to the input-independent sampling and generic interpolation of PerforatedCNNs, the sampling in our network is adaptively determined from the input such that the sampling density reflects predicted activation values, and the interpolation parameters are learned. As shown later in the experiments, this approach allows for much greater sparsity in the sampling. In [27], high-resolution predictions are generated from low-resolution results. Instead, our method is used for features rather than final outputs.

Gumbel-based selection Random selection based on the Gumbel distribution has been used in making discrete decisions for network acceleration. The Gumbel-Softmax trick was utilized in adaptively choosing network layers to apply on an input image [33] and in selecting channels or layers to skip [14]. In contrast to these techniques which determine computation based on image-level semantics for image classification, our sampling is driven by the spatial organization of features and is geared towards accurately reconstructing positional content. As a result, our method is well-suited to spatial understanding tasks such as object detection and semantic segmentation.

## 3 Methodology

In this section, we first present a general introduction of the stochastic sampling-interpolation network and then describe the stochastic sampling module and interpolation module in detail. Next, we introduce the grid prior which is found to be helpful for interpolation. Finally, we illustrate how to integrate the sampling and interpolation network modules into residual blocks.

## 3.1 Stochastic Sampling-Interpolation Network

Convolutions in neural networks typically generate an output feature map  $Y \in \mathbb{R}^{C_{out} \times H \times W}$  pixel by pixel from an input feature map  $X \in \mathbb{R}^{C_{in} \times H \times W}$ :

$$Y(p) = \sum_{p' \in R_k} W_c(p') X(p+p'), p \in \Omega, \tag{1}$$

where H and W represent the height and width of the feature map,  $C_{in}$  and  $C_{out}$  denote the input and output feature dimensions,  $\Omega = \{(i, j) | i \leq W, j \leq H, i, j \in \mathbb{Z}^+\}$  represents the spatial domain of the feature map,  $R_k$  indicates the

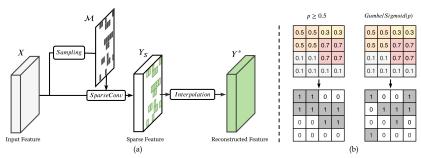


Fig. 2. (a) Stochastic sampling-interpolation network. The stochastic sampling module generates a sparse sampling mask  $\mathcal{M}$  based on the input feature map X, and then calculates features only at the sampling points, forming a sparse feature map  $Y_s$ . The features of unsampled points are interpolated by the interpolation module to form the output feature map  $Y^*$ . (b)(Left) In deterministic sampling, points with the same confidence are either sampled altogether or not sampled at all. (Right) In stochastic sampling, a random subset of the points with the same confidence will be sampled, with a sampling density determined by their confidence.

support region of kernel offsets with kernel size k (e.g., for a  $3 \times 3$  convolution,  $R_k = \{(-1, -1), (-1, 0), ..., (1, 1)\}$  and k = 3), and  $W_c \in \mathbb{R}^{C_{in} \times C_{out} \times k \times k}$  denotes convolution weights.

Spatial redundancy commonly exists in feature maps, such that features at certain points can be approximated by interpolating the features from surrounding positions. Therefore, exhaustive computation across the entire space is not required. Our method takes advantage of this using a content-aware stochastic sampling module and a trainable interpolation module.

A basic illustration of our method is shown in Fig. 2(a). The sampling module generates a mask  $\mathcal{M} \in \mathbb{R}^{H \times W}$ . In the inference phase,  $\mathcal{M}$  is binary and it is calculated first before computing Y. Points masked as 1 in  $\mathcal{M}$  are sampled, and convolution is applied only on these points, resulting in a sparse feature map  $Y_s$ , which is calculated as:

$$Y_s(p) = \begin{cases} 0 & \mathcal{M}(p) = 0 \\ Y(p) & \mathcal{M}(p) = 1. \end{cases}$$
 (2)

Then, the features of unsampled points are constructed by the interpolation module C. Together with the features of sampled points, they constitute the reconstructed output feature  $Y^*$ :

$$Y^*(p) = \begin{cases} C(Y_s)(p) & \mathcal{M}(p) = 0\\ Y_s(p) & \mathcal{M}(p) = 1. \end{cases}$$
(3)

Since the cost of calculating  $\mathcal{M}$  is less than that of Y and C, if  $\mathcal{M}$  is sparse, then the computation cost can be reduced. In our experiments, the sparsity of  $\mathcal{M}$  can be greater than 70% on average.

A technical challenge of this approach is that the binary sampling mask  $\mathcal{M}$  is non-differentiable, making this sampling module incompatible with backpropagation in the training stage. To circumvent this issue, a mask  $\mathcal{M}$  that gradually changes from soft to hard during training is used. Therefore, the mask can undergo optimization from the beginning of training and then becomes roughly

consistent to a hard mask used in the inference stage by the end of training. During training, with this soft mask, the output  $Y_s$  at each point p is calculated as follows:

$$Y_s(p) = \mathcal{M}(p) \odot Y(p), \tag{4}$$

and the full output feature  $Y^*$  is calculated by:

$$Y^*(p) = (1 - \mathcal{M}(p)) \odot C(Y_s)(p) + \mathcal{M}(p) \odot Y_s(p), \tag{5}$$

where  $\odot$  denotes broadcast multiplication.

# 3.2 Stochastic Sampling Module

In previous works [30,6,1], sampling is usually done in a deterministic manner, where points with confidence greater than a certain threshold are sampled, as shown in Fig. 2(b)(Left). But in our stochastic sampling, a higher confidence only indicates a higher probability of the point being sampled, as shown in Fig. 2(b)(Right). Due to the spatial redundancy over the feature map, adjacent points may have similar features and confidences, so deterministic sampling typically samples or not samples adjacent points together. However, stochastic sampling can sample a portion of the points, and the features of the other unsampled points can be obtained from the interpolation module. Therefore, sparser sampling can be achieved while maintaining relatively accurate feature maps.

Our sampling is based on the two-class Gumbel-Softmax distribution, which was first introduced in reinforcement learning [17] to simulate stochastic discrete sampling. Thus, the mask  $\mathcal{M}$  is defined as:

$$\mathcal{M}(p) = \frac{exp((-log(\pi_1(p)) + g_1(p))/\tau)}{\sum_{j \in \{0,1\}} exp((-log(\pi_j(p)) + g_j(p))/\tau)},$$
(6)

where  $\pi$  denotes confidence map generated from a two-class Softmax activation. g represents noise sampled from a standard Gumbel distribution, and it provides the randomness of Gumbel-Softmax. If the noise g is eliminated, Gumbel-Softmax degenerates into a deterministic function that is approximately equal to the Softmax function with a temperature term.  $\tau$  is a temperature parameter. When  $\tau$  approaches 0,  $\mathcal{M}(p)$  becomes approximately binary.

In our implementation,  $\pi$  is generated by a  $3 \times 3$  convolutional layer with a two-class Softmax activation, and  $\tau$  is exponentially decreased over iterations according to  $\tau = \alpha^{iter}\tau_0$ , where  $\alpha$  is the decay factor, *iter* is the number of iterations, and  $\tau_0$  is the initial temperature. In our experiments, we set  $\tau_0 = 1$ . Therefore, at the beginning of training, the mask is soft, allowing the sampling module to be trained. At the end of training,  $\tau$  becomes close to 0, so the mask generated by the sampling module is nearly binary, as desired for our discrete inference. To encourage the network to produce sparse sampling masks, the sparse loss is introduced on the confidence map  $\pi_1$  for all layers during training:

$$L_{sparse} = \sum_{l} \|\pi_1^l\|_1 \tag{7}$$

where  $\pi_1^l$  indicates l-th layer's confidence map, and  $\|\cdot\|_1$  indicates L1-norm. Different levels of sparsity are achieved by adjusting the weight of the sparse loss. Therefore, the training objective is:

$$L = L_{task} + \gamma L_{sparse}. \tag{8}$$

where  $L_{task}$  is the task specific objective and  $\gamma$  is the sparse loss weight.

# 3.3 Interpolation Module

In previous works [6, 7], the features of unsampled points are obtained by reusing the previous features at the corresponding points [7]:

$$Y^*(p) = \mathcal{M}(p) \odot Y_s(p) + (1 - \mathcal{M}(p)) \odot X(p) \tag{9}$$

or just by setting them to zero [6]:

$$Y^*(p) = \mathcal{M}(p) \odot Y_s(p) + (1 - \mathcal{M}(p)) \odot \mathbf{0} = \mathcal{M}(p) \odot Y_s(p). \tag{10}$$

However, these approaches ignore the spatial redundancy of feature maps, which could be used to obtain the features of unsampled points, leading to a more accurate feature map.

Our method capitalizes on this spatial redundancy to generate relatively accurate feature maps. This is done by interpolating the features of unsampled points from those of sampled points as indicated in Eq. (3). The interpolation is formulated as:

$$C(Y_s)(p) = \frac{\sum_{s_i} W_I(p, s_i) Y_s(s_i)}{\sum_{s_i} W_I(p, s_i) + \epsilon}, s_i \in \Omega,$$
(11)

where  $W_I(p,s_i) \geq 0$  represents interpolation weights,  $\Omega$  represents the spatial domain like in Eq. (1), and  $\epsilon$  is a small constant which is set to  $10^{-5}$ . In this formula, the features of the unsampled points are represented by a weighted average of the features at sampled points. However, interpolation by considering all sampled points is costly, even if  $Y_s$  is sparse in the inference phase. Fortunately, since neighboring points commonly have stronger feature correlation, we can alleviate this problem by computing the interpolation only using samples that lie within a window centered on the given unsampled point. Specifically, we restrict s to a window of radius r centered on the unsampled point p, defined as  $\mathcal{R}_s^r(p) = \{s_i | s_i \in \Omega, ||s_i - p||_{\infty} \leq r\}$ . Thus, the windowed interpolation module is formulated as:

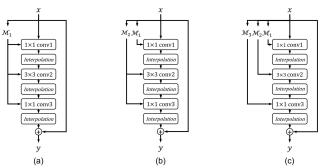
$$C(Y_s)(p) = \frac{\sum_{s_i} W_I(p, s_i) Y_s(s_i)}{\sum_{s_i} W_I(p, s_i)}, s_i \in \mathcal{R}_s^r(p).$$
 (12)

The best formulation of  $W_I$  in irregular spatial sampling is an open problem. We explored different design choices:

**RBF Kernel** The radial basis function (RBF) kernel is a commonly used interpolation function, defined as:

$$W_I(p_1, p_2) = exp(-\lambda^2 ||p_1 - p_2||^2), \tag{13}$$

where  $p_1$ ,  $p_2$  are two points and  $\lambda$  is a learnable parameter. With this kernel, an interpolated point is more greatly affected by closer points. The learnable RBF kernel is adopted by default in our approach.



**Fig. 3.** Integration of our sampling and interpolation modules into a standard residual block. (a) Applying a single mask to all convolutions; (b) Applying a separate mask on conv1, and sharing a mask for conv2 and conv3. This approach is adopted as the default setting in our experiments; (c) Applying three separate masks for each of the convolutions.

Plain Convolution Plain convolution can also be used as a learnable interpolation kernel. However, negative convolution weights may result in a zero denominator for Eq. (12). Therefore, we use the absolute value of the convolution weights in the denominator instead.

**Average Pooling** Average pooling is the simplest interpolation kernel. It assigns the same weight to all pixels in the interpolation window. We use this method as a baseline.

## 3.4 Grid Prior

With our stochastic sampling, the windowed sample set  $\mathcal{R}_s^r$  at some unsampled points may be empty. To handle this situation, one way is to fill the features of these points as zeros. To allow interpolation at these points, we instead build an equal-interval sampling mask  $\mathcal{M}_{grid}$  of stride s and combine it with the mask  $\mathcal{M}_{sample}$  generated by the sampling module:  $\mathcal{M} = \max(\mathcal{M}_{sample}, \mathcal{M}_{grid})$ . The combined mask is used in the interpolation module, and we find experimentally that it does not affect the performance but better stabilizes the training process in comparison to zero filling, as shown in Sec. 4.2.

#### 3.5 Integration with Residual Block

Our sampling and interpolation modules can be easily integrated with existing network architectures. Here, we use the residual block [11] as an example to show how these modules can be used. A natural way to insert them is by applying a separate sampling mask for each convolution (Fig. 3 (c)). However, since conv3 is a pointwise convolution (kernel size is  $1 \times 1$ ), its mask can be shared with conv2 without changing the output (Fig. 3 (b)), meanwhile avoiding the computation for generating an extra mask. Another approach is to share a single mask among all convolutions within the residual block (Fig. 3 (a)). However, since conv2 is not a pointwise convolution (kernel size is  $3 \times 3$ ), sampling a point in conv2 requires the corresponding  $3 \times 3$  points from conv1 to be sampled, which in turn places strong conditions on the sampling of conv2 if conv1 and conv2 were to share the same mask. Thus, sharing one mask for all convolutions within a block

is not an effective solution. We found the two sampling mask approach (Fig. 3 (b)) to be slightly better than others in experiments, so it is used by default.

## 4 Experiments

In this section, we validate our approach on three tasks: object detection, semantic segmentation, and image classification. Comparisons between our approach and other baseline models are conducted in terms of speed-accuracy tradeoff. Average floating point operations (FLOPs) in the backbone network over the whole validation set is used to evaluate inference speed.

## 4.1 Experimental settings

Object Detection Our models are trained on the 118k images of the COCO 2017 [24] train set, and evaluated on the 5k images of the COCO 2017 validation set. The standard mean average precision (mAP) is used to measure accuracy. The baseline model is based on Faster R-CNN with Feature Pyramid Network (FPN) [23] and deformable convolution [4, 38]. The other components and hyperparameters are based on mmdetection [2]. More details are given in Appendix.

For our modules, the sparse sampling module and interpolation module are integrated into all the residual blocks as shown in Fig. 3 (b). The window size r for interpolation is set to 7, and the stride s of the grid prior is set to 11. During training, the parameter  $\lambda$  used in the learnable RBF kernel is initialized to 3, and the decay factor  $\alpha$  of the stochastic sampling module is set to make the Gumbel-Softmax temperature parameter  $\tau=0.01$  at the end of training. During inference, we use the same  $\tau$  to produce masks, and points with mask values below than 0.5 are marked as unsampled.

Semantic Segmentation Our models are trained on the 2975 finely annotated images of the Cityscapes [3] train set and evaluated on the 500 images of the validation set. Accuracy is measured by the standard mean IoU. The baseline model is ResNet-50 based dilated FCN [26] with deformable convolution layers [4]. The implementation and hyper-parameters are based on the open-source implementation of TorchCV [36]. More details are given in Appendix. For our modules, the same hyper-parameters are used as in object detection.

**Image Classification** Our models are trained on the ImageNet-1K training set, and follow both the training and inference settings of [6]. We choose ResNet-34 as our baseline for fair comparison. Following [6], our sampling and interpolation modules are incorporated in all residual blocks, except for the first block in each stage. Since ResNet-34 is composed of basic blocks (two  $3 \times 3$  convolutions), we apply separate masks for each convolution.

Different from the previous two tasks, experiments on ImageNet use a much lower image resolution (i.e.  $224\times224$ ). Thus, for our modules, we reduce the window size r of the interpolation module to 5, and the stride s of the grid prior to 2, while keeping other hyper-parameters the same as in object detection.

#### 4.2 Ablation Study

We validate several design choices on the COCO2017 object detection benchmark. All the models are trained and evaluated on images with a shorter side of 1000 pixels and with the sparse loss weight set to 0.1.

Table 1. Comparison of different interpolation kernels on COCO2017 validation

	Avg Pool	Plain Conv	RBF Kernel
mAP	41.8	41.9	42.0
GFlops	110.0	109.8	96.6

Table 2. Validation of the interpolation module on COCO2017 validation

	Fill Zeros	Reuse Features	Ours
mAP	42.0	42.1	42.0
GFlops	164.9	226.1	96.6

Table 3. Comparison of different grid prior settings on COCO2017 validation

	s = 9	s = 11	s = 13	w/o Grid Prior
mAP	41.9	42.0	41.8	41.8
GFlops	95.4	96.6	95.3	95.0

**Different interpolation kernels** We first compare the different interpolation kernels mentioned in Sec. 3.3: *learnable RBF kernel*, *plain convolution* and *average pooling*. Results are shown in Table 1, with the sparse loss weight in the training phase set so that the three kernels yield similar accuracy. It can be seen that the *learnable RBF kernel* consumes fewer FLops.

Effect of removing interpolation We further study the effect of removing the interpolation module by replacing it with reusing features of the previous layer [7] (see Eq. (9)) and directly filling zeros [6] (see Eq. (10)). Results are shown in Table 2. Reusing features achieves 42.1 mAP with 226.1 GFLOPs, while Filling zeros obtains 39.7 mAP with 121.8 GFLOPs. Both of these methods consume noticeably more computation than our interpolation module with similar or worse accuracy, which indicates that they are inferior to our method in terms of FLOPs-accuracy tradeoff.

Effect of Grid Prior We compare performance under different grid stride s or by removing the grid prior. Results are shown in Table 3. We found the performance in different settings to be similar, but the training process is sometimes not stable at s=13 and without the grid prior. Therefore, we choose s=11 as our default setting.

## 4.3 Object Detection

We compared our method to other baselines on the COCO2017 object detection benchmark. To better illustrate the FLOPs-accuracy tradeoff under different parameters and settings, we display charts rather than tables and present the numerical results in Appendix.

Uniform sampling The direct way to sample uniformly is by using a mask with sampling points at equal intervals. However, since the interval must be integer, the minimal interval would be 2, which limits its feasibility for handling arbitrary sampling ratios. Thus, instead of using an equal-interval sampling mask, we choose to downsample the input images and not include our modules, which can also be seen as a uniform sampling method. We compared these two approaches with a downsampling rate of 2, and experimental results show the performance of these two approaches to be very similar<sup>6</sup>.

<sup>&</sup>lt;sup>6</sup> 38.5 mAP for downsampled images vs. 38.7 mAP for equal-interval sampling masks.

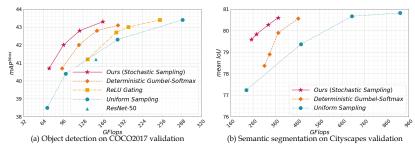


Fig. 4. (a) Tradeoff curves for different sampling methods on object detection (COCO2017 validation). Curve of "Uniform Sampling" is drawn from various resolutions, and others are drawn from various sparse loss weights (with shorter side of 1000 pixels). (b) Experiments on the Cityscapes semantic segmentation benchmark. Except for "Uniform Sampling", all other models are trained and evaluated on images with a shorter side of 1024 pixels.

For uniform sampling, we resize the shorter sides of input images to {1000, 800, 600, 500} to draw the FLOPs-accuracy curve. For our method, the curve is drawn according to different sparse loss weights, i.e. {0.2, 0.1, 0.05, 0.02}. Results are shown in Fig. 4(a). For our baseline without any sampling or interpolation modules, which is trained and evaluated on images with a shorter side of 1000 pixels, it obtains 43.4 mAP with 289.5 GFLOPs. In comparison, our stochastic sampling method performs at 43.3 mAP with only 160.4 GFLOPs, saving nearly half of the computation cost with no accuracy drop.

**Deterministic sampling** We next compare our stochastic sampling strategy to deterministic sampling methods. For fair comparison, we only replace the sampling module, while leaving the other parts unchanged. Two deterministic sampling methods are examined:

- ReLU Gating: Similar to LCCL [6], a ReLU function is applied on sampling confidence to generate a sparse mask, by trimming values smaller than 0 during training and inference. The sampling confidence is generated by a  $3 \times 3$  convolution with a 1-dimensional output. Since the output of ReLU is not binary, Eq. (5) is used to calculate the full output feature.
- Deterministic Gumbel-Softmax: Eliminating the noise g in Gumbel-Softmax naturally results in deterministic sampling, as described in Sec. 3.2, and is approximately equal to the Softmax function with a temperature term.

For our stochastic sampling and the two deterministic sampling methods, the FLOPs-accuracy curve is drawn according to different sparse loss weights, i.e.  $\{0.2, 0.1, 0.05, 0.02\}$ . Compared to ReLU Gating, our stochastic sampling achieves clearly better performance than the ReLU based approach, especially at lower levels of computation. Deterministic Gumbel-Softmax performs better than ReLU Gating, but is still worse than stochastic sampling by a large margin.

Smaller backbones Replacing a large backbone with a smaller backbone is a common method for reducing computation. Thus, we also compare our method (based on ResNet-101) with a baseline using ResNet-50 as the smaller backbone.

Table 4. Performance comparison on the ImageNet validation set. All the methods are based on ResNet-34. Our models are trained with a loss weight of 0.01 and 0.015 to achieve accuracy or FLOPs similar to other methods for fair comparison. "w/o Interp" indicates removing the interpolation module and filling the features of unsampled positions with 0

Method	Type	Top-1/Top-5 Acc Drop(%)	FLOPs	Speedup
LCCL [6]	spatial	0.43/0.17	$2.7 \times 10^{9}$	24.8%
PFEC [20]	pruning	1.06/-	$2.7 \times 10^{9}$	24.2%
SFP [12]	pruning	2.09/1.29	$2.2 \times 10^{9}$	41.1%
FPGM [13]	pruning	1.29/0.54	$2.2 \times 10^{9}$	41.1%
Ours(0.01)	spatial	0.45/0.20	$2.53 \times 10^{9}$	
Ours(0.015)	spatial	1.19/0.47	$2.16 \times 10^{9}$	42.4%
w/o Interp(0.015)	spatial	1.07/0.46	$2.21 \times 10^{9}$	41.0%

ResNet-50 achieves 41.2 mAP with 149.2 GFLOPs, which indicates accuracy much worse than that of our method with similar computation costs.

#### 4.4 Semantic Segmentation

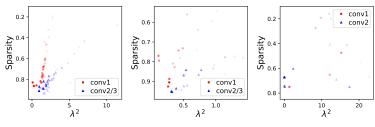
We also conduct experiments on the Cityscapes benchmark for semantic segmentation. Unless otherwise specified, all the models are trained and evaluated on images with a shorter side of 1024 pixels. We first compare our content-aware stochastic sampling to *uniform sampling* and deterministic sampling. For deterministic sampling, we choose *deterministic Gumbel-Softmax* for comparison because of its better performance exhibited on the object detection task.

Results are shown in Fig. 4(b) and present the numerical results in Appendix. For our method and deterministic Gumbel-Softmax, we draw curves according to different sparse loss weights {0.3, 0.2, 0.1, 0.05}. For uniform sampling, we resize the shorter sides of input images to {1024, 896, 736, 512}. The original baseline with a shorter side of 1024 pixels obtains 80.8 mean IoU with 920.6 GFLOPs. In comparison, our method obtains 80.6 mean IoU with only 373.2 GFLOPs, saving nearly 60% of the computation cost with almost no accuracy drop. Compared with other sampling methods, i.e. uniform sampling and deterministic Gumbel-Softmax, our method clearly achieves a better FLOPs-accuracy tradeoff.

## 4.5 Image Classification

We also compare our method to other state-of-the-art methods [6, 20, 12, 13] for reducing computation on the ImageNet-1K image classification benchmark. Similar to our method, LCCL [6] explores sparsity in the spatial domain, while PFEC [20], SFP [12] and FPGM [13] reduce computation by model pruning. Results are presented in Table 4. Our models are trained under different sparse loss weights, 0.01 and 0.015, to reach accuracy or FLOPs similar to other methods for fair comparison. Compared with LCCL [6] and PFEC [20], our approach achieves comparable accuracy with less FLOPs. Compared with SFP [12] and FPGM [13], our approach obtains a smaller accuracy drop with similar FLOPs.

We further remove the interpolation module from our method and fill the features of unsampled points with 0. Results show that removing interpolation



(a) Object Detection (b) Semantic Segmentation (c) Image Classification

**Fig. 5.** Relationship between interpolation parameter  $\lambda$  and sparsity. The transparency of each point represents its location in the network, where a darker color indicates greater depth. A larger  $\lambda$  yields a sharper interpolation kernel; when  $\lambda > 3$ , it approximates an identity kernel.

does not affect performance on the ImageNet validation set. This is inconsistent with object detection and semantic segmentation. We believe that this is because the classification network is focused on extracting global feature representations. Therefore, as long as the features of certain key points are calculated and preserved, the global features will not be affected and the performance will not be hurt. In other words, in the image classification task, it is not important to reconstruct the features of unsampled points by interpolation.

## 4.6 Analysis of sampling and interpolation modules

In this section, we further study the relationship between the sampling and interpolation modules. Specifically, we analyze the relationship between sampling sparsity and the parameter  $\lambda$  in the RBF interpolation module. A larger  $\lambda$  indicates a sharper interpolation kernel, and when  $\lambda > 3$ , the RBF interpolation kernel approximates an identity kernel, for which there is no interpolation.

Results are shown in Fig. 5. For object detection (see Fig. 5 (a)) and semantic segmentation (see Fig. 5 (b)), sparsity and  $\lambda$  show a strong negative correlation. When sparsity is high, the  $\lambda$  is usually small, which means sampled points far away from the unsampled points also contribute greatly to the interpolation. However, for image classification (see Fig. 5 (c)), this correlation has not been observed.  $\lambda$  is quite large in most cases, which means the effect of the interpolation module is limited. This phenomenon is consistent with the experimental results of the "w/o Interp" entry in Table 4, that the results without interpolation are almost identical to our full model, further indicating that interpolation is not important for image classification.

Another observation is that the  $\lambda$  of conv1 are consistently smaller than  $\lambda$  of conv2 and conv3 in object detection and semantic segmentation. The reason is still unclear but we suspect that this phenomenon may be related to the receptive field of operators.

#### 4.7 Realistic run-time on CPU

Our method achieves good theoretical speed-accuracy trade-offs. In this section, we present a preliminary evaluation of the realistic run-time of the backbone

**Table 5.** Comparison of theoretical and realistic speedups on E5-2650 and I7-6650U. Baseline model is trained and evaluated on images with a shorter side of 1000 pixels. The CPU run-time is calculated on the COCO2017 validation set

Model	mAP	GFLOPs	Runtime(s/img)		Real Speedup		Theo Speedup
			E5-2650 v2	I7-6650U	E5-2650 v2	I7-6650U	Theo Speedup
Baseline	43.4	289.5	4.9	10.5	-	-	-
Our(0.02)	43.3	160.4	4.0	7.9	1.23	1.33	1.80
Our(0.05)	42.8	122.8	3.5	6.3	1.40	1.67	2.36
Our(0.1)	42.0	96.6	3.2	5.3	1.53	1.98	3.00
Our(0.2)	40.7	73.3	2.7	4.4	1.81	2.39	3.95

network on the CPU. According to Eq. (2), the mask  $\mathcal{M}$  is calculated before computing Y to decide which points in Y need to calculated. In order to show the realistic speedup of our method under different computing resources, we conducted experiments in two different hardware environments: a workstation (E5-2650 v2, 256G RAM and Ubuntu 16.04 OS) and a laptop (I7-6650U, 16G RAM and Ubuntu 16.04 OS). For fair comparison, we replace all convolutions by our implementation for all models, and enable multi-threading by default (32 threads for E5-2650 v2 and 4 threads for I7-6650U). Results on object detection are shown in Table 5. There is a gap between the theoretical speedup and the realistic speedup, but the gap for laptop (I7-6650U) is smaller than for workstation (E5-2650 v2). The main reason is that the workstation has better computing speed and more cores (E5-2650 v2 has 8 cores and I7-6650U has 2 cores), but has a memory access bottleneck. This suggests that our method is more suitable for devices with less computation speed but relatively faster IO speed, such as mobile or edge computing devices.

In addition, some works [30,29] have developed general techniques to accelerate sparse convolution based methods. For example, SCNN [29] designed a hardware accelerator for sparse convolution and demonstrate that ideal speedup of sparse convolution is achievable in such devices. SBNet [30] is a general method to accelerate sparse convolution in GPU. These general techniques of accelerating sparse convolution are compatible with our method, and can further close the gap between theoretical and actual speedup in real applications.

# 5 Conclusion

A method for reducing computation in convolutional networks was proposed that exploits the intrinsic sparsity and spatial redundancy in feature maps. We present a stochastic sampling and interpolation scheme to avoid expensive computation at spatial locations that can be effectively interpolated. To overcome the challenge of training binary decision variables for representing discrete sampling locations, Gumbel-Softmax is introduced to our sampling module. The effectiveness of this approach is verified on a variety of computer vision tasks.

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