## Controlling Material Appearance by Examples — Supplemental Material

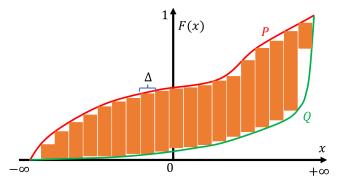


Figure 1: Practically, we should numerically evalute discrete CDF of P and Q from their samples and compute the integral of their difference (orange regions).

## 1. Sliced Cramér Loss

In our main paper, we introduce a resampling method to compute the Sliced Wasserstein loss between different number of samples. However, this loss can be accurately computed without resampling. We propose a sliced Cramér loss based on 1D Cramér distance [Cra28]. Given two distributions P and Q over R, their Cumulative Distribution Functions (CDFs) are respectively  $F_P$  and  $F_Q$ . The p-Cramér distance between P and Q is defined as:

$$l_p(P,Q) = \left( \int_{-\infty}^{+\infty} \left( F_P(x) - F_Q(x) \right)^p dx \right)^{\frac{1}{p}} \tag{1}$$

when p=1, the Cramér distance is equivalent to the first order Wasserstein distance. Practically, we have sample points drawn from P and Q (pixels on labeled regions in our case). We need to construct their empirical CDFs and numerically compute the integral shown in Fig. 1. In Listing 1, we show a Pytorch implementation, based on sorting with a time complexity of O((M+N)log(M+N)) where M and N are numbers of sample points for P and Q respectively. The implementation supports auto-differentiation with batch training.

Theoretically, a sliced Cramér loss should be a more precise comparison of two distributions than resampled Sliced Wasserstein. However, while evaluating the Cramér loss, we do not observe significant improvement over our resampling strategy. As the Cramér loss requires an additional sorting operation (line 10

in Listing 1), its evaluation is slightly slower. In Fig. 2, we show a few results generated using the sliced Wasserstein loss with resampling and the sliced Cramér loss respectively.

```
import torch as T
  def cramer_loss(u, v):
      \mbox{\tt\#} u is an array with shape [B, M] and v is an
       array with shape [B, N];
      # B is batch size, M and N are number of
      sample points
      # compute empirical CDFs of u and v
      u_{-} = T.sort(u, dim=1)
      v_{-} = T.sort(v, dim=1)
      w = T.cat((u, v), dim=1).sort(dim=1)
      u_cdf = T.searchsorted(u_, w[:, :-1], right=
      v_cdf = T.searchsorted(v_, w[:, :-1], right=
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      u_cdf = u_cdf / u.shape[1]
      v_{cdf} = v_{cdf} / v.shape[1]
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      # compute delta
16
17
      deltas = torch.diff(w, dim=1)
18
19
      # compute integral
      return ((u_cdf - v_cdf) * deltas).abs().sum(
      dim=1).mean()
```

Listing 1: Implementation of 1D Cramér loss

## References

[Cra28] CRAMÉR H.: On the composition of elementary errors: Statistical applications. Almqvist and Wiksell, 1928. 1

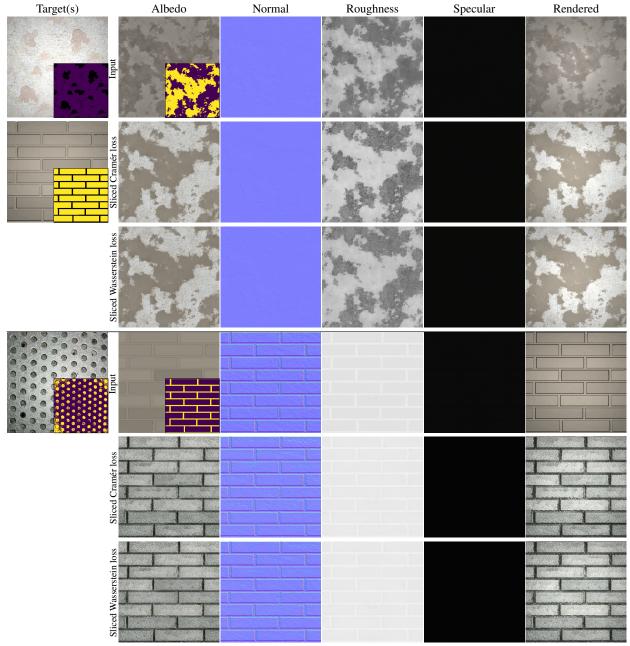


Figure 2: We compare material transfer results generated using the Sliced Cramér loss and resampled Sliced wasserstein loss. As mentioned, despite a theoretically better behaved formulation we do not observe significant benefit of the Sliced Cramér loss over the resampled Sliced Wasserstein loss.