Radar Pulse Deinterleaving with Transformer Based Deep Metric Learning

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Abstract—When receiving radar pulses it is common for a recorded pulse train to contain pulses from many different emitters. The radar pulse deinterleaving problem is the task of separating out these pulses by the emitter from which they originated. Notably, the number of emitters in any particular recorded pulse train is considered unknown. In this paper, we define the problem and present metrics that can be used to measure model performance. We propose a metric learning approach to this problem using a transformer trained with the triplet loss on synthetic data. This model achieves strong results in comparison with other deep learning models with an adjusted mutual information score of 0.882.

I. Introduction

Radar pulse deinterleaving aims to separate out a train of radar pulses by the emitters from which they originated. We want to transform a single interleaved pulse train into many smaller deinterleaved pulse trains where each train contains all the pulses from a single emitter and only pulses from that emitter. A major challenge in this problem is that the number of emitters and hence the number of pulse trains we need to separate out is unknown.

Deinterleaving is a key part of the electronic warfare (EW) [1] pipeline, falling under the signal analysis stage. It aims to extract useful information from the detected signals, allowing for further analysis. It is much easier to analyse many smaller deinterleaved signals than a single large interleaved signal for tasks such as determining the radar's operating mode [2], specific emitter identification (SEI) [3], and operator intention identification [4].

The main challenges faced in deinterleaving arise from congested and contested environments. In congested environments there is a large number of emitters often with significant overlap in parameter space, meaning it is difficult to differentiate between them. In contested environments we are faced with radars designed to be hard to detect. This can lead to dropped pulses due to low signal to noise ratio as well as high variance between pulses from the same emitter. Some examples of these are emitters that hop between frequencies and those with a variable pulse repetition interval (PRI), meaning there is varying time between the start of emitted pulses. This could result in a system reporting that there are more radars than are actually present.

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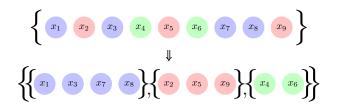


Fig. 1. The radar pulse deinterleaving problem requires partitioning a set of radar pulses by the emitter from which the pulses originated. Here the colours represent the emitters and the features vectors x_i denote PDWs. This can be approached using metric learning where we generate an embedding for each pulse in the context of the whole pulse train and cluster these embeddings using an off the shelf clusterer.

Classically, deinterleaving was achieved through identifying the PRI for each emitter. Techniques for this include the cumulative difference histogram (CDIF) [5] and the sequential difference histogram (SDIF) [6]. These techniques work well when detecting pulses from emitters with a constant PRI. However, as the electromagnetic environment becomes more crowded we will require techniques that are capable of deinterleaving trains containing emitters which are more agile, often in an adversarial way.

Recently the rise of deep learning has yielded new methods which can leverage large amounts of data to identify complex patterns and provide robustness to noise. The authors believe these approaches will be vital to successful deinterleaving in the presence of ever more agile digital radars. However, current applications of these methods have strict limitations. They can be largely divided into two categories. The first [7]-[14] assumes that the number of emitters is fixed thus reducing the deinterleaving problem to a classification problem. This assumption is unlikely to hold in practice and the methods that make this assumption cannot easily adapt to different numbers of emitters. In this category a large number of works use variants of recurrent neural networks (RNNs) [7]-[12] motivating our choice of baseline. Similar to classical deinterleavers, the second category of deep learning approaches to deinterleaving [15]–[18] attempt to identify pulse repetition intervals (PRIs) from times of arrival (ToAs) to deinterleave pulse trains. These methods often ignore additional pulse descriptor word (PDW) features extracted from raw IQ data, thereby restricting performance.

This paper seeks to address these limitations by proposing a pipeline that can leverage the full set of features while making no assumptions about the number of emitters.

Our key contributions are:

- We define the deinterleaving problem where the number of emitters is unknown and show how to measure performance with identified metrics (Section II).
- We combine metric learning with sequence-to-sequence models to propose a class of solutions to the problem as formulated. (Section III-A)
- We analyse the types of model that would be most appropriate for solving this problem, concluding that the transformer is a strong candidate (Section III-B).
- We propose the use of the triplet loss with the batch all triplet mining strategy during training of the model (Section III-C).
- We train a sequence-to-sequence transformer with the triplet loss on a synthetic dataset and show that it has strong performance when compared to benchmark techniques (Section IV).

II. PROBLEM FORMULATION

Given a *pulse train* of n radar pulses $X = \{x_1, \ldots, x_n\}$ from an unknown number of emitters N, we want to *partition* the pulse train by the emitters from which the pulses originated. For example if the pulses x_1, x_2, x_3, x_4, x_5 came from emitters 1, 1, 2, 2, 1 respectively the ground truth partition would be $\{\{x_1, x_2, x_5\}, \{x_3, x_4\}\}$. The goal of deinterleaving is to create a model where for each pulse train in the test dataset the predicted partition $U = \{U_1, U_2, \ldots\}$ is as similar as possible to the ground truth partition $V = \{V_1, V_2, \ldots\}$, where U_i and V_j are the subsets of the partitions known as *blocks*. To evaluate this we use standard *extrinsic clustering* metrics and take their mean over every pulse train in the dataset. The metrics are:

Adjusted Mutual Information (AMI):

The mutual information between two partitions of a set is

$$MI(U,V) = \sum_{i=1}^{|U|} \sum_{j=1}^{|V|} \frac{|U_i \cap V_j|}{n} \log \frac{n|U_i \cap V_j|}{|U_i||V_j|}$$
(1)

This does not adjust for chance so a random prediction will give a non-zero score. To account for this we use *adjusted* mutual information [19] which is given by

$$AMI(U,V) = \frac{MI(U,V) - \mathbf{E}[MI(U,V)]}{\frac{1}{2}(H(U) + H(V)) - \mathbf{E}[MI(U,V)]}$$
(2)

where n is the number of pulses in the train and $H(\cdot)$ is the entropy of a partition. AMI gives a value of 1 when the partitions are identical and 0 when the mutual information is equal to the expected mutual information.

Adjusted Rand Index (ARI):

Rand index [20] considers pairs of points. A correct prediction is when a pair is predicted to be in the same block when they are truly in the same block (true positive) or if they

are predicted to be in different blocks when they are truly in different blocks (true negative). We then define rand index as

$$RI(U, V) = \frac{\text{\#correct predictions}}{\text{\#pairs}}$$
 (3)

This can be seen as the accuracy of a pairwise binary classification problem asking 'are these two points from the same block'. Much like with AMI we adjust for chance to get the adjusted rand index with a value of 1 when the partitions are identical and 0 when the rand index is equal to the expected rand index.

V-measure:

V-measure [21] is defined as the harmonic mean of homogeneity and completeness. Homogeneity quantifies to what degree each cluster contains only members of a single class. Completeness quantifies to what degree all members of a given class are assigned to the same cluster. It is often useful to examine homogeneity and completeness individually to indicate whether a model has more of a tendency to split up points from the same ground truth cluster (low completeness) or to combine points from different ground truth clusters (low homogeneity). V-measure has a minimum value of 0 and a maximum value of 1 where higher is better. It is not adjusted for chance.

III. METHOD

A. A metric learning approach to deinterleaving

Since the number of emitters is unknown, the number of blocks in the ground truth partition for any pulse train is also unknown. This means that familiar deep learning classification pipelines are not applicable.

Instead we take a *metric learning* [22] approach. We use a *sequence-to-sequence* model to generate an *embedding* z_i for each pulse x_i in a pulse train X. We train this model so that embeddings for pulses from the same emitter are close together and those from different emitters are far apart (as measured by their Euclidean distance). During inference we can cluster these embeddings using a non-parametric clustering algorithm such as HDBSCAN [23] to form a predicted partition for the original pulse train. A visual representation of the training and inference pipelines can be seen in Figure 2.

For the deinterleaving problem individual pulses have little meaning in isolation. It is therefore important to use a sequence-to-sequence model to ensure that their embeddings are generated in the *context* of the rest of the pulse train. In other words, the model must consider the *relationships* between the pulses.

B. Transformer embedding model

The transformer [24] is a popular architecture that has demonstrated state of the art results across natural language processing and computer vision. One of the key features that makes it powerful is the way it explicitly models the

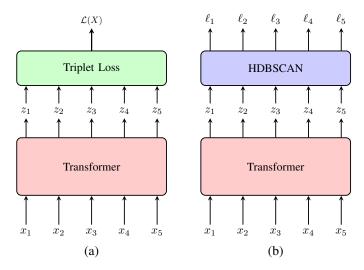


Fig. 2. The architecture for the training (a) and inference (b) pipelines share the common component of a sequence-to-sequence transformer used as a pulse embedding model. We use vanilla dot product attention and no positional encodings. (a) During training we use the triplet loss to push apart embeddings of pulses from different emitters and bring together those from the same emitter. (b) During inference we use a clustering algorithm (HDBSCAN) to generate an integer label l_i for each embedding from which we can trivially determine the predicted partition for the set of pulses.

relationship between every pair of tokens through its attention mechanism. Every token can influence every other token and the strength and form of this influence is learned during training. Given the importance of relationships between pulses in the deinterleaving problem, as discussed in the previous section, the transformer's explicit modelling of these relationships makes it a compelling choice for our model.

Specifically, the model we consider is a vanilla sequence-to-sequence transformer with dot product self-attention. The order of detection of the pulses may be inferred by their ToA meaning positional encodings are not necessary and thus were not used. Preliminary experiments showed they had a small negative effect on performance, likely due to redundant information.

C. Clustering embeddings with the triplet loss

As stated in Section III-A we want to train the model so that, within a pulse train, embeddings for pulses from the same emitter are close together and those from different emitters are far apart. The triplet loss [25] is a common loss used for deep metric learning. The loss considers triplets of embeddings, an anchor z, a positive embedding z^+ which is from the same emitter as the anchor, and a negative embedding z^- which is from a different emitter to the anchor. It compares the distance between the anchor and the positive to the distance between the anchor and the negative. It aims to satisfy the triplet inequality stated in equation 4.

$$d(z, z^+) + \alpha < d(z, z^-) \tag{4}$$

This says that we want the distance between the anchor and the positive to be less than the distance between the anchor and the negative by some margin α so that the embedding of

pulses from the same emitter are closer together than those from different emitters. When this inequality is satisfied we do not want to update the model, so the loss must be 0. When it is not satisfied we want to penalise the model by the degree to which it is violated. We therefore use the hinge loss in equation 5.

$$\mathcal{L}(z, z^+, z^-) = \max\{d(z, z^+) - d(z, z^-) + \alpha, 0\}$$
 (5)

For each pulse train there is on the order of $\mathcal{O}(n^3)$ possible triplets where n is the number of pulses in the pulse train. Often most of these triplets are relatively uninformative to the model. For example, easy triplets which satisfy equation 4 have a loss of 0. Therefore, training using all possible triplets is very inefficient.

To increase the efficiency of training with the triplet loss we need a method to choose triplets that will be informative for training. This is called a triplet mining strategy. Our experiments found that the triplet mining strategy with the best performance was the batch all [26] strategy where we take the mean of the loss across all non-easy triplets.

The batch all triplet loss is defined as follows. Given the embeddings $Z = \{z_1, \ldots, z_n\}$, let the set of all non-easy triplets be

$$\mathcal{T}^{+}(Z) = \left\{ (z_i, z_j, z_k) \in Z^3 \middle| \begin{array}{c} i \sim j \not\sim k, \\ i \neq j \neq k, \\ d(z_i, z_j) + \alpha \ge d(z_i, z_k) \end{array} \right\}$$

where \sim is the equivalence relation where $i\sim j$ indicates that the pulses at index i and j come from the same emitter. The batch all triplet loss is

$$\mathcal{L}(Z) = \frac{1}{|\mathcal{T}^{+}(Z)|} \sum_{\substack{(z_i, z_j, z_k) \in \mathcal{T}^{+}(Z)}} \max \{ d(z_i, z_j) - d(z_i, z_k) + \alpha, 0 \}$$

In this work we use the Euclidean distance metric.

IV. EXPERIMENTAL WORK

A. Datasets

Using synthetic data is overwhelmingly the norm in deinter-leaving. The authors are only aware of one paper [18] that uses real data. This is for two main reasons. The first is that real data is very difficult to ground truth. Usually labels are assigned to pulses by an existing deinterleaver. Training on this data would result in the performance of our models converging to that of the existing deinterleavers. Since the goal of this work is to develop models that outperform existing deinterleavers this approach would be pathologically constrained. The second reason is that as far as the authors are aware there are no real publicly available datasets suitable for deinterleaving.

We generated synthetic training, validation, and test sets using a simulator. These contain 100,000 pulse trains in the train set, 10,000 in the validation set, and another 10,000 pulse trains in the test set. Each pulse train consists of 1000 pulses, and contains between 2 and 20 emitters. Although

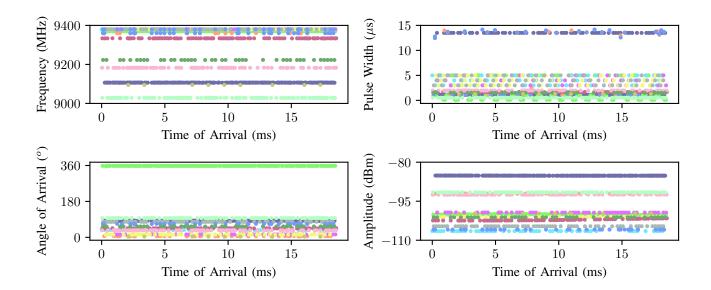


Fig. 3. A plot of the different PDW features against time of arrival for an example pulse train in the test dataset. The pulse train contains 13 emitters with significant overlap in parameters. The colour of each data point indicates which of the emitters the pulse originated from.

processing pulse trains with only one emitter in them is of interest, when this is the case no valid triplets can be formed. Therefore, training on them results in zero loss and the model is not updated.

For feature vectors to represent pulses we use pulse descriptor words (PDWs) with 5 dimensions: time of arrival (ToA), centre frequency, pulse width (PW), angle of arrival (AoA), and amplitude. Figure 3 shows an example of the PDWs from one of the pulse trains in our test set.

B. Model description

We use 8 transformer layers with 8 attention heads, a feed-forward hidden size of 2048, and a residual size of 256. The model is trained with Adam [27] with a learning rate of 0.0001, and a batch size of 8 for 8 epochs. The triplet loss margin α is 1.9, and dropout is 0.05. Embeddings z_i are projected to 8 dimensions. Hyperparameters were selected via a parameter sweep on the validation set.

C. Data Normalization

We normalise the data internally to each pulse train only. To do this we treat each element of the PDWs separately. The ToAs are linearly re-scaled so that the lowest ToA is 0 and the highest is 1. The centre frequency, pulse width, and amplitude are all statistically normalised by taking away the mean and dividing by the standard deviation in the pulse train. Finally AoA is divided by 360° to place it in the range [0, 1].

D. Baselines

As far as we are aware there are no readily available deep learning based benchmarks for this problem as formulated. In addition to this, the datasets on which other results were tested are not publicly available making comparison between solutions challenging. Instead, to gain a performance baseline we trained a sequence-to-sequence gated recurrent unit (GRU) [28] model using the same triplet loss discussed in Section III-C. Similar architectures to this are widely used in deinterleaving related literature as discussed in Section I. The key point to note about the GRU is that the relationships between pulses are expressed less explicitly than in the transformer. An internal state vector is propagated through the network meaning tokens only attend to those before it in the sequence. We used 8 GRU layers with a hidden size of 512 resulting in the number of parameters being roughly the same as that of the transformer.

As a sanity check we also include results obtained using an identity model where the PDWs themselves are clustered using HDBSCAN. We found the minimum cluster size parameter of HDBSCAN had a large effect on performance. For our experiments we set it to 20 for all models.

E. Results and discussion

As shown in Table I, the transformer outperforms our chosen baseline models in all performance metrics from Section II

Figure 4 displays the results of the transformer when predicting the number of clusters in each pulse train. These values are extracted directly from the deinterleaved pulse trains by counting the number of clusters present. This, along

TABLE I RESULTS ON THE TEST SET

Deinterleaver	Transformer	GRU	Identity
AMI	0.882	0.850	0.761
ARI	0.817	0.811	0.610
V-measure	0.884	0.852	0.765
Homogeneity	0.938	0.856	0.899
Completeness	0.853	0.872	0.697

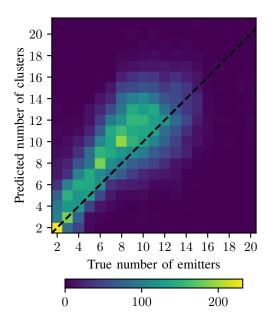


Fig. 4. A confusion matrix of the true number of emitters in the synthetic test set against the number of clusters in the partition predicted by our transformer. The black line represents an ideal scenario where the predicted number of clusters is the same as the true number of clusters. It can be seen that the transformer generally predicts there to be more clusters than there are emitters.

with the homogeneity and completeness scores, indicates that the transformer and identity models have a general trend to overestimate the number of clusters, with a RMS error of 3.07, and 3.48 respectively, whereas the GRU tends to underestimate the number of clusters, with an RMS error of 2.57. The transformer achieves a completeness score of 0.938 meaning clusters overwhelmingly contain only pulses from a single emitter. From our perspective, correctly forming clusters that contain pulses from only one emitter, even if we incorrectly split that emitter's pulses into separate clusters, is better than mixing pulses from different emitters together. This is because it suggests the deinterleaver is very effective at distinguishing between different emitters. Furthermore, we believe this incorrect splitting of pulses from a single emitter into multiple clusters could be corrected during a post processing step which combines related clusters - thus increasing completeness and V-measure.

Figure 5 shows the mean adjusted mutual information for models across an increasing number of emitters. The performance is strong for 5 or more emitters. However, when there are few ground truth clusters performance is poor, regardless of the model used. This suggests the issue is not related to the training scheme, as the untrained identity model exhibits a similar trend. The problem must therefore be due to the dataset or the clustering algorithm itself. All the models tend to over-split clusters when the number of emitters is low. They maintain high homogeneity, but completeness drops rapidly for fewer than 5 emitters. This could be due to three reasons, each identifiable in Figure 6:

1) The ground truth clusters are generally much larger for

- low numbers of emitters.
- 2) The variability between ground truth cluster sizes is much higher for low numbers of emitters.
- 3) Imbalance between cluster sizes is more prominent for small numbers of emitters, with one big cluster and many smaller ones.

All these effects are likely amplified by the fact that pulse trains have been limited to length 1000. By addressing these potential issues in the dataset or clustering approach, the performance for low numbers of emitters could likely be improved.

V. FUTURE WORK

Future work will explore addressing the issues concerning performance at low numbers of emitters as discussed in Section IV-E. Equally, scaling this model to longer and variable length sequences would also prove beneficial as the dynamics of the previously discussed clipping issues could be reduced. In practice, real pulse trains are longer and variable in length so this scenario would be more realistic. Finally, we will explore applying this model to real data collects. The main challenge involved in this is obtaining reliable ground truth labels as discussed in section IV-A.

VI. CONCLUSION

In this paper we defined the deinterleaving problem and proposed a metric learning approach as a solution using a transformer trained with the triplet loss on a synthetic dataset. We showed that it achieves strong results in comparison with an identity baseline indicating that this method is effective for solving the deinterleaving problem. Additionally we showed that, in this pipeline, the transformer architecture outperforms a GRU architecture commonly used in the literature. Future work on transformers with this pipeline can be exploited to further improve deinterleaving performance.

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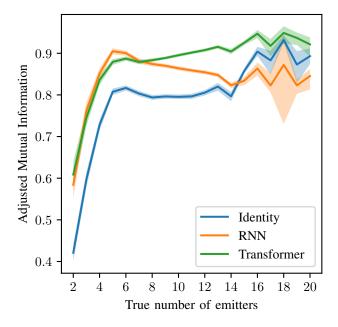


Fig. 5. The true number of emitters against the mean AMI for that number of clusters. The shaded area represents the bootstrapped [29] 0.1-0.9 confidence interval. It is notable that the width of the confidence interval is much larger for high numbers of clusters where we have few data points (The distribution of data points can be seen in Figure 4).

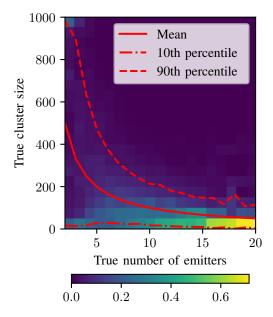


Fig. 6. The true number of emitters against the true cluster size (the number of pulses from each emitter in a pulse train). The lines represent the mean, 10th percentile and 90th percentile. They show that the size and variability in size of clusters decreases as the number of emitters increases. The heat map colours represent the proportion of clusters in pulse trains with a particular number of emitters that are within a given size bin. They show that for low numbers of emitters there is significant imbalance between cluster sizes, as the number of emitters increases the cluster sizes become more balanced.

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