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Recursive Nearest Neighbor Search in a Sparse and Multiscale Domain for Comparing Audio Signals

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Abstract

We investigate recursive nearest neighbor search in a sparse domain at the scale of audio signals. Essentially, to approximate the cosine distance between the signals we make pairwise comparisons between the elements of localized sparse models built from large and redundant multiscale dictionaries of time-frequency atoms. Theoretically, error bounds on these approximations provide efficient means for quickly reducing the search space to the nearest neighborhood of a given data; but we demonstrate here that the best bound defined thus far involving a probabilistic assumption does not provide a practical approach for comparing audio signals with respect to this distance measure. Our experiments show, however, that regardless of these non-discriminative bounds, we only need to make a few atom pair comparisons to reveal, e.g., the origin of an excerpted signal, or melodies with similar time-frequency structures.

Keywords: Multiscale decomposition; sparse approximation; time-frequency dictionary; audio similarity

1. Introduction

- Sparse approximation is essentially the modeling of data with few terms
- ³ from a large and typically overcomplete set of atoms, called a "dictionary" [24].

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Consider an $\mathbf{x} \in \mathbb{R}^K$, and a dictionary \mathcal{D} composed of N unit-norm atoms in the same space, expressed in matrix form as $\mathbf{D} \in \mathbb{R}^{K \times N}$, where $N \gg K$. A pursuit is an algorithm that decomposes \mathbf{x} in terms of \mathbf{D} such that $||\mathbf{x} - \mathbf{D}\mathbf{s}||^2 \le \epsilon$ for some error $\epsilon \ge 0$. (In this paper, we work in a Hilbert space unless otherwise noted.) When \mathbf{D} is overcomplete, \mathbf{D} has full row rank and there exists an infinite number of solutions to choose from, even for $\epsilon = 0$. Sparse approximation aims to find a solution \mathbf{s} that is mostly zeros for ϵ small. In that case, we say that \mathbf{x} is sparse in \mathcal{D} .

Matching Pursuit (MP) is an iterative descent sparse approximation method

Matching Pursuit (MP) is an iterative descent sparse approximation method based on greedy atom selection [17, 24]. We express the *n*th-order model of the signal $\mathbf{x} = \mathbf{H}(n)\mathbf{a}(n) + \mathbf{r}(n)$, where $\mathbf{a}(n)$ is a length-*n* vector of weights, $\mathbf{H}(n)$ are the *n* corresponding columns of \mathbf{D} , and $\mathbf{r}(n)$ is the residual. MP augments the *n*th-order representation, $\mathcal{X}_n = \{\mathbf{H}(n), \mathbf{a}(n), \mathbf{r}(n)\}$, according to:

$$\mathcal{X}_{n+1} = \begin{cases}
\mathbf{H}(n+1) = [\mathbf{H}(n)|\mathbf{h}_n], \\
\mathbf{a}(n+1) = [\mathbf{a}^T(n), \langle \mathbf{r}(n), \mathbf{h}_n \rangle]^T, \\
\mathbf{r}(n+1) = \mathbf{x} - \mathbf{H}(n+1)\mathbf{a}(n+1)
\end{cases}$$
(1)

using the atom selection criterion

$$\mathbf{h}_n = \arg\min_{\mathbf{d} \in \mathcal{D}} ||\mathbf{r}(n) - \langle \mathbf{r}(n), \mathbf{d} \rangle \mathbf{d}||^2 = \arg\max_{\mathbf{d} \in \mathcal{D}} |\langle \mathbf{r}(n), \mathbf{d} \rangle|$$
(2)

where $||\mathbf{d}|| = 1$ is implicit. The inner product here is defined $\langle \mathbf{x}, \mathbf{y} \rangle \stackrel{\Delta}{=} \mathbf{y}^T \mathbf{x}$. This criterion guarantees $||\mathbf{r}(n+1)||^2 \le ||\mathbf{r}(n)||^2$ [24]. Other sparse approximation methods include orthogonal MP [28], orthogonal least squares (OLS) [41, 33], molecular methods [9, 38, 19], cyclic MP and OLS [36], and minimizing a relaxed 21 sparsity measure [6]. These approaches have higher computational complexities 22 than MP, but can produce data models that are more sparse. 23 Sparse approximation is data-adaptive and can produce parametric and multiscale models having features that function more like mid-level "objects" than low-level projections onto sets of vectors [9, 19, 38, 8, 22, 27, 32, 34, 43]. These 26 aspects make sparse approximation a compelling complement to state-of-the-art 27 approaches for and applications of comparing audio signals based upon, e.g., monoresolution cepstral and redundant time-frequency representations, such as fingerprinting [42], cover song identification [26, 10, 3, 35], content segmentation, indexing, search and retrieval [16, 4], artist or genre classification [39].

In the literature we find some existing approaches to working with audio sig-32 nals in a sparse domain. Features built from sparse approximations can provide 33 competitive descriptors for music information retrieval tasks, such as beat track-34 ing, chord recognition, and genre classification [40, 32]. Sparse representation classifiers have been applied to music genre recognition [27, 5], and robust speech recognition [12]. Parameters of sparse models can be compared using histograms 37 to find similar sounds in acoustic environment recordings [7, 8], or atoms can be learned to compare and group percussion sounds [34]. Biologically-inspired 30 sparse codings of correlograms of sounds can be used to learn associations between descriptive high-level keywords and audio features such that new sounds can be automatically categorized, and large collections of sounds can be queried 42 in meaningful ways [22]. Outside the realm of audio signals, sparse approxi-43 mation has been applied to face recognition [44], object recognition [29], and 44 landmine detection [25]. 45

In this paper, we discuss the comparison of audio signals in sparse domains, but not specifically for fingerprinting or efficient audio indexing and search — 47 two tasks that have been convincingly solved [42, 13, 16, 18]. We explore the possibilities and effectiveness of comparing, atom-by-atom, audio signals modeled 49 using sparse approximation and large overcomplete time-frequency dictionaries. Our contributions are three-fold: 1) we generalize an iterative nearest-neighbor search algorithm to comparing subsequences [14, 15]; 2) we show that though 52 sparse models of audio signals can be compared by considering pairs of atoms, 53 the best bound so far derived [14, 15] does not make a practical procedure; 54 and 3) we show experimentally that the hierarchic comparison of audio signals in a sparse domain still provides intriguing and informative results. Overall, our work here shows that a sparse domain can facilitate comparisons of audio 57 signals in "hierarchical" ways through comparing individual elements of each sparse data model organized roughly in order of importance.

In the next two sections, we discuss and elaborate upon a recursive method of nearest neighbor search in a sparse domain [14, 15]. We extend this method to comparing subsequences, and examine the practicality of probabilistic bounds on the distances between neighbors. In the fourth section, we describe several experiments in which we compare a variety of audio signals through comparisons of their sparse models. We conclude with a discussion about the results and several future directions.

₆₇ 2. Nearest Neighbor Search by Recursion in a Sparse Domain

8 Consider a set of signals

$$\mathcal{Y} \stackrel{\Delta}{=} \{ \mathbf{y}_i \in \mathbb{R}^K : ||\mathbf{y}_i|| = 1 \}_{i \in \mathcal{I}}$$
 (3)

where $\mathcal{I} = \{1, 2, ...\}$ indexes this set, and a query signal $\mathbf{x}_q \in \mathbb{R}^K$, $||\mathbf{x}_q|| = 1$.

Assume that we have generated sparse approximations for all of these signals $\widehat{\mathcal{Y}} \stackrel{\Delta}{=} \{\{\mathbf{H}_i(n_i), \mathbf{a}_i(n_i), \mathbf{r}_i(n_i)\} : \mathbf{y}_i = \mathbf{H}_i(n_i)\mathbf{a}_i(n_i) + \mathbf{r}_i(n_i)\}_{i \in \mathcal{I}}$ using a dictionary \mathbf{D} that spans the space \mathbb{R}^K , and giving the n_q -order representation $\{\mathbf{H}_q(n_q), \mathbf{a}_q(n_q), \mathbf{r}_q(n_q)\}$ for \mathbf{x}_q . Since \mathcal{D} spans \mathbb{R}^K , \mathcal{D} is "complete," and any signal in \mathbb{R}^K is "compressible" in \mathcal{D} , meaning that we can order the representation weights in $\mathbf{a}_i(n_i)$ or $\mathbf{a}_q(n_q)$ in terms of decreasing magnitude, i.e.,

$$0 < |[\mathbf{a}_i(n_i)]_{m+1}| \le |[\mathbf{a}_i(n_i)]_m| \le Cm^{-\gamma}, m = 1, 2, \dots, n_i - 1$$
 (4)

for n_i arbitrarily large, with C>0, and where $[\mathbf{a}]_m$ is the mth element of the column vector \mathbf{a} . This can be seen in the magnitude representation weights in Fig. 1, which are weights of sparse representations of piano notes, described in Section 4.1. With MP and a complete dictionary, we are guaranteed $\gamma>0$ because $||\mathbf{r}(n+1)||^2<||\mathbf{r}(n)||^2$ for all n [24].

Consider the Euclidean distance between two signals of the same dimension, which is the cosine distance for unit-norm signals. Thus, with respect to this distance, the $\mathbf{y}_i\in\mathcal{Y}$ nearest to \mathbf{x}_q is given by solving

$$\min_{i \in \mathcal{I}} ||\mathbf{y}_i - \mathbf{x}_q|| = \max_{i \in \mathcal{I}} \langle \mathbf{x}_q, \mathbf{y}_i \rangle.$$
 (5)

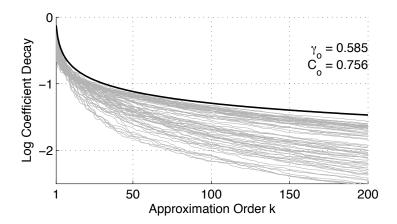


Figure 1: Gray lines show decays of representation weight magnitudes as a function of approximation order k for several decompositions of unit-norm signals (4-second recordings of single piano notes described in Section 4.1). Thick black line shows a global compressibility bound with its parameters.

We can express this inner product in terms of sparse representations

$$\langle \mathbf{x}_q, \mathbf{y}_i \rangle = \langle \mathbf{H}_q(n_q) \mathbf{a}_q(n_q) + \mathbf{r}_q(n_q), \mathbf{H}_i(n_i) \mathbf{a}_i(n_i) + \mathbf{r}_i(n_i) \rangle$$

$$= \mathbf{a}_i^T(n_i) \mathbf{H}_i^T(n_i) \mathbf{H}_q(n_q) \mathbf{a}_q(n_q) + O[\mathbf{r}_q, \mathbf{r}_i].$$
(6)

- With a complete dictionary we can make $O[\mathbf{r}_q,\mathbf{r}_i]$ negligible by choosing ϵ ar-
- bitrarily small, so we can express (5) as

$$\max_{i \in \mathcal{I}} \langle \mathbf{x}_q, \mathbf{y}_i \rangle \sim \max_{i \in \mathcal{I}} \mathbf{a}_i^T(n_i) \mathbf{G}_{iq} \mathbf{a}_q(n_q) = \max_{i \in \mathcal{I}} \sum_{m=1}^{n_i} \sum_{l=1}^{n_q} [\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}]_{ml}$$
(7)

- where $[\mathbf{B} \bullet \mathbf{C}]_{ml} = [\mathbf{B}]_{ml} [\mathbf{C}]_{ml}$ is the Hadamard, or entry wise, product, $[\mathbf{B}]_{ml}$
- is the element of **B** in the *m*th row of the *l*th column, $\mathbf{G}_{iq} \stackrel{\Delta}{=} \mathbf{H}_{i}^{T}(n_{i})\mathbf{H}_{q}(n_{q})$ is a
- 89 $n_i \times n_q$ matrix with elements from the Gramian of the dictionary, i.e., $\mathbf{G} \stackrel{\Delta}{=} \mathbf{D}^T \mathbf{D}$,
- and finally we define the outer product of the weights

$$\mathbf{A}_{iq} \stackrel{\Delta}{=} \mathbf{a}_i(n_i) \mathbf{a}_q^T(n_q). \tag{8}$$

- 91 2.1. Recursive Search Limited by Bounds
- Since we expect the decay of the magnitude of elements in $\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}$ to be
- fastest in diagonal directions by (4), we define a recursive sum along the M

anti-diagonals starting at the top left:

$$S_{iq}(M) \stackrel{\Delta}{=} S_{iq}(M-1) + \sum_{m=1}^{M} [\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}]_{m(M-m+1)}$$

$$\tag{9}$$

for $M=2,3,\ldots,\min(n_i,n_q)$, and setting $S_{iq}(1)=[\mathbf{A}_{iq}\bullet\mathbf{G}_{iq}]_{11}$. With this we can express the argument of (7) as

$$\langle \mathbf{x}_q, \mathbf{y}_i \rangle \approx \sum_{m=1}^{n_i} \sum_{l=1}^{n_q} [\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}]_{ml} = S_{iq}(M) + R(M)$$
 (10)

where at step M, we are comparing M additional pairs of atoms to those considered in the previous steps. R(M) is a remainder that we will bound. The total number of atom pairs contributing to $S_{iq}(M)$ (9) is

$$P(M) \stackrel{\Delta}{=} \sum_{m=1}^{M} m = M(M+1)/2.$$
 (11)

The approach taken by Jost et al. [14, 15] to find the nearest neighbors of \mathbf{x}_q in \mathcal{Y} bounds the remainder R(M) by compressibility (4). Assuming we have a positive upper bound on the remainder, i.e., $R(M) \leq \widetilde{R}(M)$. we know lower and upper bounds on the cosine distance $L_{iq}(M) \leq \langle \mathbf{x}_q, \mathbf{y}_i \rangle \leq U_{iq}(M)$, where

$$L_{iq}(M) \stackrel{\Delta}{=} S_{iq}(M) - \widetilde{R}(M) \tag{12}$$

$$U_{iq}(M) \stackrel{\Delta}{=} S_{iq}(M) + \widetilde{R}(M). \tag{13}$$

Finding elements of \mathcal{Y} close to \mathbf{x}_q with respect to (5) can be done recursively over the approximation order M. For a given M, we find $\{S_{iq}(M)\}_{i\in\mathcal{I}}$, compute the remainder $\tilde{R}(M)$, and eliminate signals that are not sufficiently close to \mathbf{x}_q with respect to their cosine distance by comparing the bounds. This 107 approach is similar to hierarchical ones, e.g., [21], where the features become 108 more discriminable as the search process runs. (Also note that compressibility 109 is similar to the argument made in justifying the truncation of Fourier series in 110 early work on similarity search [1, 11, 30], i.e., that power spectral densities of 111 many time-series decay like $\mathcal{O}(|f|^{-b})$ with b > 1.) 112 Starting with M = 1, we compute the sets $\{L_{iq}(1)\}_{i \in \mathcal{I}}$ and $\{U_{iq}(1)\}_{i \in \mathcal{I}}$, 113 that is, the first-order upper and lower bounds of the set of distances of \mathbf{x}_q

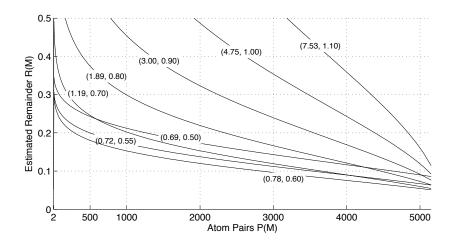


Figure 2: Estimated remainder (assuming unit-norm signals) using bound in (16) with p=0.2 (probability that remainder does not exceed bound) and n=100 (number of elements in each sparse model) as a function of the number of atom pairs already considered for several pairs of compressibility parameters (C, γ) estimated from the dataset used to produce Fig. 1.

from all signals in \mathcal{Y} . Then we find the index of the largest lower bound 115 $i_{\max} = \arg \max_{i \in \mathcal{I}} L_{iq}(1)$, and reduce the search space to $\mathcal{I}_1 \stackrel{\Delta}{=} \{i \in \mathcal{I} : U_{iq}(1) \geq$ 116 $L_{i_{\max}q}(1)$, since all other data have a least upper bound on their inner prod-117 uct with \mathbf{x}_q than the greatest lower bound in the set. For the next step, we 118 compute the sets $\{L_{iq}(2)\}_{i\in\mathcal{I}_1}$ and $\{U_{iq}(2)\}_{i\in\mathcal{I}_1}$, find the index of the maximum $i_{\max} = \arg \max_{i \in \mathcal{I}_1} L_{iq}(2)$, and construct the reduced set $\mathcal{I}_2 \stackrel{\Delta}{=} \{i \in \mathcal{I}_1 :$ 120 $U_{iq}(2) \geq L_{i_{\max}q}(2)$. Continuing in this way, we find the elements of \mathcal{Y} closest 121 to \mathbf{x}_q at each M with respect to the cosine distance by recursing into the sparse 122 approximations of the signals. 123

2.2. Bounding the Remainder

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To reduce the search space quickly we desire that (12) and (13) converge quickly to the neighborhood of $\langle \mathbf{x}_q, \mathbf{y}_i \rangle$, or in other words, that the bounds on the remainder quickly become discriminative. Jost et al. [14, 15] derive three different bounds on R(M). From the weakest to the strongest, these are:

1. $[\mathbf{G}_{iq}]_{ml} = 1$ (worst case scenario, and impossible for n > 1)

$$R(M) \le C^2 (||\mathbf{c}_M^{\gamma}||_1 + ||\mathbf{d}^{\gamma}||_1)$$
 (14)

2. $[\mathbf{G}_{iq}]_{ml} \sim \text{iid Bernoulli}(0.5), \ \Omega = \{-1, 1\} \text{ (impossible for } n > 1)$

$$R(M) \le C^2 \sqrt{\ln 4} \left(||\mathbf{c}_M^{\gamma}||_2^2 + ||\mathbf{d}^{\gamma}||_2^2 \right)^{1/2}$$
 (15)

3. $[\mathbf{G}_{iq}]_{ml} \sim \text{iid Uniform, } \Omega = [-1, 1],$

$$R(M) \le C^2 \sqrt{2/3} \operatorname{Erf}^{-1}(p) \left(||\mathbf{c}_M^{\gamma}||_2^2 + ||\mathbf{d}^{\gamma}||_2^2 \right)^{1/2}$$
 (16)

with probability $0 \le p \le 1$

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where we define the following vectors for $n \stackrel{\Delta}{=} \min(n_i, n_q)$ and $M = 2, \dots, n_q$

$$\mathbf{c}_{M}^{\gamma} \stackrel{\Delta}{=} \{ [l(m-l+1)]^{-\gamma} : m = M+1, \dots, n; l = 1, \dots, m \}$$
 (17)

$$\mathbf{d}^{\gamma} \stackrel{\Delta}{=} \{ [l(n-m+1)]^{-\gamma} : m = 1, \dots, n-1; l = m+1, \dots, n \}.$$
 (18)

Appendix A gives derivations of these bounds, as well as the efficient computation of (16) for the special case of $\gamma=0.5$. The parameters (C,γ) describe the compressibility of the signals in the dictionary (4). The bounds of (15) and (16) are much more discriminative than (14) because they involve an ℓ_2 -norm at the price of uncertainty in the bound. The bound in (16) is attractive because we can tune it with the parameter p, which is the probability that the remainder will not exceed the bound. Figure 2 shows bounds based on (16) for several pairs of compressibility parameters for the dataset used to produce Fig. 1.

2.3. Estimating the Compressibility Parameters

The bounds (14)–(16), and consequently the number of atom pairs we must consider before the bounds become discriminable, depend on the compressibility parameters, (C, γ) — which themselves depend in a complex way on the signal, the dictionary, and the method of sparse approximation. Figure 3 shows the error surface, feasible set, and the optimal parameters for the dataset used to produce Fig. 1. We describe our parameter estimation procedure in Appendix B. The resulting bound is shown in black in Fig. 1. These compressibility parameters also agree with those seen in Fig. 2.

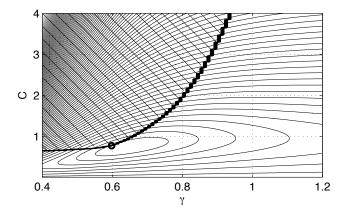


Figure 3: Error surface as a function of the compressibility parameters for the dataset used to produce Fig. 1, with the feasible set shaded at top left, and optimal parameters marked by a circle.

3. Subsequence Nearest Neighbor Search in a Localized Sparse Domain

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The recursive nearest neighbor search so far described has the obvious limitation that it cannot be applied to comparing subsequences of large data vectors, as is natural for comparing audio signals. Thus, we must adapt its structure to work for comparing subsequences in a set of data

$$\mathcal{Y} \stackrel{\Delta}{=} \{ \mathbf{y}_i \in \mathbb{R}^{N_i} : N_i \ge K \}_{i \in \mathcal{I}}$$
(19)

(note that now we do not restrict the norms of these signals). We can create from the elements of \mathcal{Y} a new set of all subsequences having the same length as a K-dimensional query \mathbf{x}_q ($K < N_i$):

$$\mathcal{Y}_K \stackrel{\Delta}{=} \left\{ \mathbf{P}_t \mathbf{y}_i / ||\mathbf{P}_t \mathbf{y}_i|| : t \in \mathcal{T}_i = \{1, 2, \dots N_i - K + 1\}, \mathbf{y}_i \in \mathcal{Y} \right\}$$
(20)

where \mathbf{P}_t extracts a K-length subsequence in \mathbf{y}_i starting a time-index t (it is an identity matrix of size K starting a column t in a $K \times N_i$ matrix of zeros). The set \mathcal{T}_i are times at which we create length-K subsequences from \mathbf{y}_i . If we decompose each of these by sparse approximation, then we can use the framework in the previous section. However, sparse approximation is an expensive operation that

we want to do only once for the entire signal, and independent of the length of \mathbf{x}_q .

To address this problem, we instead approximate each element in \mathcal{Y}_K by building local sparse representations from the global sparse approximations of each \mathbf{y}_i , and then calculating their distance to \mathbf{x}_q using the framework in the previous section. From here on we consider only the K-length subsequences of a single element $\mathbf{y}_i \in \mathcal{Y}$ without loss of generality (i.e., all other elements of \mathcal{Y}_i can be included as subsequences). Toward this end, consider that we have decomposed the N_i -length signal \mathbf{y}_i using a complete dictionary to produce the representation $\{\mathbf{H}_i(n_i), \mathbf{a}_i(n_i), \mathbf{r}_i(n_i)\}$. From this we construct the local sparse representations of \mathbf{y}_i :

$$\widehat{\mathcal{Y}}_K \stackrel{\Delta}{=} \left\{ \left\{ \mathbf{P}_t \mathbf{H}_i(n_i), \xi_t \mathbf{a}_i(n_i), \mathbf{P}_t \mathbf{r}_i(n_i) \right\} : t \in \mathcal{T}_i \right\}$$
(21)

where the time partition \mathcal{T}_i is the set of all times at which we extract a Klength subsequence from \mathbf{y}_i , and ξ_t is set such that $||\xi_t \mathbf{P}_t \mathbf{y}_i|| = 1$, i.e., each
length-K subsequence is unit-norm. For each K-dimensional subsequence, (7)
now becomes

$$\max_{t \in \mathcal{T}_{i}} \langle \mathbf{x}_{q}, \mathbf{P}_{t} \mathbf{y}_{i} \rangle = \max_{t \in \mathcal{T}_{i}} \left[\langle \mathbf{H}_{q}(n_{q}) \mathbf{a}_{q}(n_{q}), \xi_{t} \mathbf{P}_{t} \mathbf{H}_{i}(n_{i}) \mathbf{a}_{i}(n_{i}) \rangle + O[\mathbf{r}_{q}, \mathbf{r}_{i}] \right]$$

$$\approx \max_{t \in \mathcal{T}_{i}} \xi_{t} \mathbf{a}_{i}^{T}(n_{i}) \mathbf{H}_{i}^{T}(n_{i}) \mathbf{P}_{t}^{T} \mathbf{H}_{q}(n_{q}) \mathbf{a}_{q}(n_{q})$$

$$= \max_{t \in \mathcal{T}_{i}} \xi_{t} \sum_{m=1}^{n_{i}} \sum_{l=1}^{n_{q}} [\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}(t)]_{ml} \tag{22}$$

where \mathbf{A}_{iq} is defined in (8), we define the time-localized Gramian

$$\mathbf{G}_{iq}(t) \stackrel{\Delta}{=} \mathbf{H}_{i}^{T}(n_{i}) \mathbf{P}_{t}^{T} \mathbf{H}_{q}(n_{q})$$
(23)

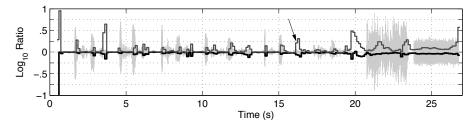
and we have excluded the terms involving the residuals because we can make them arbitrarily small.

3.1. Estimating the Localized Energy

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The only thing left to do is find an expression for ξ_t so that each subsequence is comparable with the others with respect to the cosine distance. We

(a) Six Speech Signals (0-20 s), Music Excerpt (21-23 s), Realization of GWN (24-27 s)



(b) Music: Orchestra

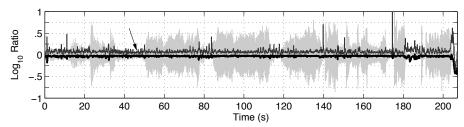


Figure 4: Short-term energy ratios, $\log_{10}(\sum_{j=1}^{n_t} w_j a_j^2/||\mathbf{P}_t \mathbf{y}_i||^2)$, over 1 s windows (hopped 125 ms) for MP decompositions using 8xMDCT [31] to a global residual energy 30 dB below the initial signal energy. Arrow points to line (top, gray) using weighting $w_j = 1$. The other line (bottom, black) uses (25). Data in (a) are described in Section 4.2; data in (b) are described in Section 4.3.

assume that the localized energy can be approximated from the local sparse representation in the following way assuming $||\mathbf{P}_t \mathbf{y}_i|| > 0$

$$\xi_t = ||\mathbf{P}_t \mathbf{y}_i|| \approx \sqrt{\mathbf{a}_i^T(n_i)\mathbf{H}_i^T(n_i)\mathbf{P}_t^T \mathbf{P}_t \mathbf{H}_i(n_i)\mathbf{a}_i(n_i)} \approx \sqrt{\sum_{j=1}^{n_t} w_j a_j^2}$$
(24)

where the n_t weights $a_j \in \{[\mathbf{a}_i(n_i)]_m : [\mathbf{H}_i^T(n_i)\mathbf{P}_t^T\mathbf{P}_t\mathbf{H}_i(n_i)]_{ml} \neq 0, 1 \leq m, l \leq n_i\}$ are those associated with atoms having support in [t, t + K), and w_j we define to weigh the contribution of a_j^2 to the localized energy estimate. We set $\xi_t = 0$ if $\sum_{j=1}^{n_t} a_j^2 = 0$.

If all atoms contributing to the subsequence have their entire support in [t, t + K), and are orthonormal, then we can set each $w_j = 1$. This does not hold for subsequences of a signal decomposed using an overcomplete dictionary, as shown by Fig. 4. For much of the time we see $\sum_{j=1}^{n_t} a_j^2 \ge ||\mathbf{P}_t \mathbf{y}_i||^2$, which

means our localized estimate of the segment energy is greater than its real value.

This will make ξ_t and consequently (22) smaller.

Instead, we make a more reasonable estimate of $||\mathbf{P}_t \mathbf{y}_i||$ by accounting for

Instead, we make a more reasonable estimate of $||\mathbf{P}_t \mathbf{y}_i||$ by accounting for the fact that atoms can have support outside [t, t+K). For instance, if an atom has some fraction of support in the subsequence we multiply its weight by that fraction. We thus weigh the contribution of the jth atom to the subsequence norm using

$$w_{j} = \begin{cases} 1, & u_{j} \geq t, u_{j} + s_{j} \leq t + K \\ (K/s_{j})^{2}, & u_{j} < t, u_{j} + s_{j} \geq t + K \\ (u_{j} + s_{j} - t)^{2}/s_{j}^{2}, & u_{j} < t, t < u_{j} + s_{j} \leq t + K \\ (t + K - u_{j})^{2}/s_{j}^{2}, & t \leq u_{j} < t + K, u_{j} + s_{j} > t + K \end{cases}$$

$$(25)$$

where u_j and s_j are the position and scale, respectively, of the atom associated with the weight a_j . In other words, if an atom is completely in [t, t + K), it contributes all of its energy to the approximation; otherwise, it contributes only a fraction based on how its support intersects [t, t + K). With this we are now slightly underestimating the localized energies, as seen in Fig. 4. In both of these cases for $\{w_j\}$, however, we can assume by the energy conservation of MP [24] that as the subsequence length becomes larger our error in estimating the subsequence energy goes to zero, i.e.,

$$\lim_{K \to N_i} ||\mathbf{P}_t \mathbf{y}_i||^2 - \sum_{j=1}^{n_t} w_j a_j^2 = ||\mathbf{P}_t \mathbf{r}_i(n_i)||^2.$$
 (26)

With a complete dictionary, the right hand side can be made zero. Significant departures from the energy estimate of subsequences can be due to the interactions between atoms [37].

3.2. Recursive Subsequence Search Limited by Bounds

Now, similar to (9) and (10), we can say,

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$$\langle \mathbf{x}_q, \mathbf{P}_t \mathbf{y}_i \rangle \approx \xi_t \sum_{m=1}^{n_t} \sum_{l=1}^{n_q} [\mathbf{A}_{iq}(t) \bullet \mathbf{G}_{iq}(t)]_{ml} = S_{iq}(t, M) + R(t, M)$$
 (27)

where for $M=2,3,\ldots,\min(n_i,n_q)$, and with $S_{iq}(t,1)=\xi_t[\mathbf{A}_{iq}(t)\bullet\mathbf{G}_{iq}(t)]_{11}$,

$$S_{iq}(t,M) \stackrel{\Delta}{=} S_{iq}(t,M-1) + \xi_t \sum_{m=1}^{M} [\mathbf{A}_{iq}(t) \bullet \mathbf{G}_{iq}(t)]_{m(M-m+1)}. \tag{28}$$

The problem of finding the subsequence closest to \mathbf{x}_q with respect to the cosine distance can now be done iteratively over M by bounding each remainder R(t,M) using (14), (15), or (16), and the method presented in in Section 2.1.
Furthermore, we can compare only a subset of all possible subsequences using
a coarser time partition \mathcal{T}_i .

222 3.3. Practicality of the Bounds for Audio Signals

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The experiments by Jost et al. [14, 15] use small images (128 square) and 223 orthogonal wavelet decompositions, which do not translate to audio signals de-224 composed over redundant time-frequency dictionaries. Jost et al. [14, 15] do 225 not state the compressibility parameters they use, but for the high-dimensional 226 audio signals with which we work in this paper it is not unusual to have $\gamma \approx 0.5$ 227 when using MP and highly overcomplete dictionaries. We find that decom-228 posing 4 s segments of music signals (single channel, 44.1 kHz sample rate, 229 representation weights shown in Fig. 1) using the dictionary in Table 1 requires on average 2,375 atoms to reduce the residual energy 20 dB below the initial 23 signal energy. Thus, for the bound (16) using n = 2,375 atoms, and with the 232 parameters $(C, \gamma) = (0.4785, 0.5)$ (in the feasible set), Fig. 5 clearly shows that 233 in order to have any discriminable bound (say ± 0.2 for unit-norm signals) we 234 must either select a low value for p — in which case we are assuming the first 235 atom comparison is approximately the cosine distance — or we must make over a million pairwise comparisons. 237

This is not practical for signals of large dimension, and dictionaries containing billions of time-frequency atoms. There is no possibility of tabulating the dictionary Gramian for quick lookup of atom pair correlations; and the cost of looking up atoms in million-atom decompositions is expensive as well. It is clear then that the tightest bound given in (16) is not practical for efficiently discrim-

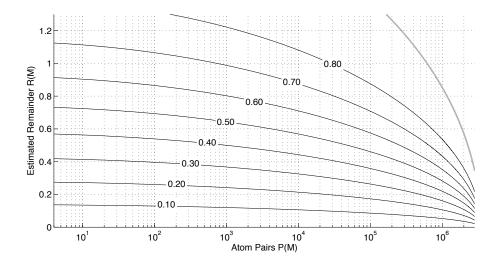


Figure 5: Estimated remainder (assuming unit-norm signals) as a function of the number of atom pairs already considered for dataset used to produce Fig. 1. Gray: bound in (15). Black, numbered: bound in (16) for several labeled p (probability that remainder does not exceed bound) with n = 2,375 (number of elements in each sparse model), and $(C, \gamma) = (0.4785, 0.5)$.

inating distances between audio signals with respect to their cosine distance (5)
decomposed by MP and time-frequency dictionaries.

4. Experiments in Comparing Audio Signals in a Sparse Domain

Though approximate nearest neighbor subsequence search of sparsely approximated audio signals with the bound (16) is impractical, we have found that approximating the cosine distance in a sparse domain has some intriguing behaviors. We now present several experiments where we compare different types of audio data in a sparse domain under a variety of conditions. All signals are single channel, and have a sampling rate of 44.1 kHz. We decompose each by MP [17] using either the dictionary in Table 1, or the 8xMDCT dictionary [31].

4.1. Experiment 1: Comparing Piano Notes

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In this experiment, we look at how well low-order sparse approximations of sampled piano notes embody their harmonic characteristics by comparing

s (samples/ms)	$\Delta_u \text{ (samples/ms)}$	Δ_f (Hz)
128/3	32/0.7	43.1
256/6	64/2	43.1
512/12	128/3	43.1
1024/23	256/6	43.1
2048/46	512/12	21.5
4096/93	1024/23	10.8
8192/186	2048/46	5.4
16384/372	4096/93	2.7
32768/743	8192/186	1.3

Table 1: Time-frequency dictionary parameters (44.1 kHz sampling rate): atom scale s, time resolution Δ_u , and frequency resolution Δ_f . Finer frequency resolution for small-scale atoms is achieved with interpolation by zero-padding.

them using the methods presented in Section 2. The data in set 'A' are 68 notes 257 (chromatically spanning A0 to G#6) on a real and somewhat in-tune piano; and 258 in set the data 'B' are 39 notes (roughly a C major scale C0 to D6) on a real and very out-of-tune piano with very poor recording conditions. We truncate all signals to have a dimension of 176,400 (4 seconds), and decompose each 261 by MP [17] over a redundant dictionary of time-frequency Gabor atoms, the 262 parameters of which are summarized in Table 1. We stop each decomposition 263 once its residual energy drops 40 dB below the initial energy. We normalize the weights of each model by the square root energy of the respective signal. We do 265 not align the time-domain signals such that the note onsets occur at the same 266 time. Figure 1 shows the ordered decays of the weights in the sparse models of 267 data set 'A'. 268

Figure 6(a) shows the magnitude correlations between all pairs of signals in set 'A' evaluated in the time-domain. The overtone series is clear as diagonal lines offset at 12, 19, 24, and 28, semitones from the main diagonal. Figure 6(b) shows the approximated magnitude correlations (9) using only M = 10 atoms from each signal approximation (thus P(10) = 55 atom pairs). Even though the

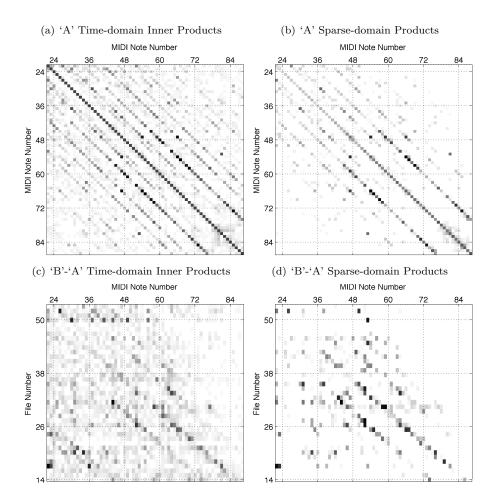


Figure 6: $|S_{iq}(10)|$ (9) for two sets of recorded piano notes. (a) and (b): Set 'A' compared with itself in time and sparse domains (M = 10). (c) and (d): Set 'B' (rows) compared with set 'A' (columns) in time and sparse domains (M = 10). Elements on main diagonals in (a) and (b) are scaled by 0.25 to increase contrast of other elements.

mean number of atoms in this set of models is about 7000 we still see portions of the overtone series. The diagonal in Fig. 6(b) does not have a uniform color because low notes have longer sustain times than high notes, and the sparse models thus have more time-frequency atoms with greater energies spread over the support of the signal. Figure 6(c) show the magnitude correlations between sets 'B' and 'A' evaluated in the time-domain; and Fig. 6(d) shows the

magnitude correlations (9) using only M=10 atoms from each model. In a sparse domain, we can more clearly see the relationships between the two sets because the first 10 terms of each model are most likely related to the stable harmonics of the notes, and not to the noise. We can see a diatonic scale starting around MIDI number 36, as well as the fact that the pitch scale in data set 'B' lies somewhere in-between the semitones in data set 'A'.

Figure 7(a) shows the approximate magnitude correlations $|S_{iq}(M)|$ (9), as 286 well as the upper and lower bounds on the remainder using the tightest bound (16) with p = 0.2, and n = 100, for the signal A3 from set 'A' and the rest of the 288 set. Here we can see that the lower bound for the largest magnitude correlation 289 exceeds the upper bound of all the rest after comparing only M=19 atoms 290 from each decomposition. All but five of the signals can be excluded from the 291 search after M = 6. The four other signals having the largest approximate magnitude correlation are labeled, and are harmonically related to the signal 293 through its overtone series. With a signal selected from set 'B' and compared to 294 set 'A', Fig. 7(b) shows that we must compare many more atoms between the 295 models until the bounds have any discriminability. After P(M) = 1500 atom 296 comparisons we can see that the largest magnitudes $|S_{iq}(M)|$ (9) are roughly 297 harmonically related to the detuned D5 from set 'B'. 298

As a final part of this experiment, we look at the effects of comparing atoms 299 with parameters within some subset. As done in Fig. 6(d), we compare the 300 sparse approximations of two different sets of piano notes, but here we only 301 consider those atoms that have scales greater than 186 ms. This in effect means that we look for signals that share the same "long-term" time-frequency be-303 haviors. The resulting $|S_{iq}(10)|$ (9) is shown in Fig. 8. We see correlations 304 between the notes much more clearly compared with Fig. 6(d). Removing the 305 short-term phenomena improves "tonal"-level comparisons between the signals because non-overlapping yet energetic short atoms are replaced by atoms representative of the note harmonics.

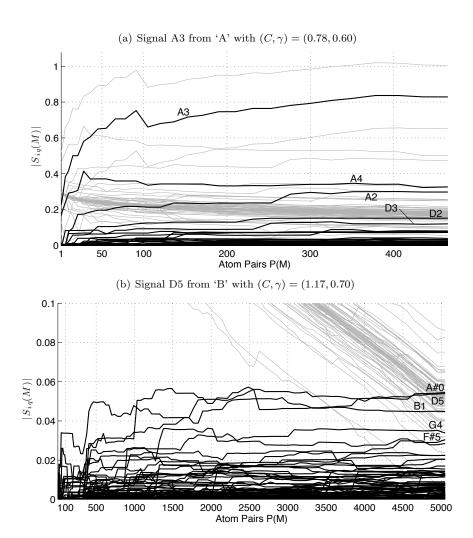


Figure 7: Black: $|S_{iq}(M)|$ (9) as a function of the number of atom pairs considered for the set of piano notes in 'A' with a signal from either (a) 'A' (note A3) or (b) 'B' (note D5 approximately). Gray: for each $S_{iq}(M)$, magnitudes of $L_{iq}(M)$ (12) and $U_{iq}(M)$ (13) using bound in (16) with p = 0.2 (probability that remainder does not exceed bound), and n = 100 (number of elements in each sparse model). Largest magnitude correlations are labeled. Note differences in axes.

4.2. Experiment 2: Comparing Speech Signals

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In this experiment, we test how efficiently using (28) we can find in a speech signal the time from which we extract some \mathbf{x}_q . We also test how distortion in the query affects these results. We make a signal by combining six segments of

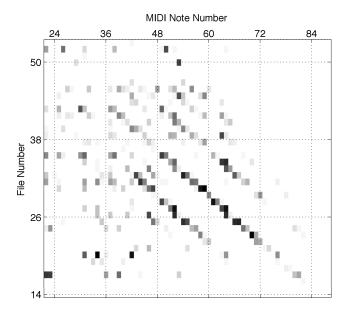


Figure 8: $|S_{iq}(10)|$ (9) for two sets of recorded piano notes in a sparse domain using only the atoms with duration at least 186 ms. Compare with Fig. 6(d).

speech, a short music segment, and white noise, shown in Fig. 4(a). The six 313 speech segments are the same phrase spoken by three females and three males: 314 "Cottage cheese with chives is delicious." We extract from one of these speech 315 signals the word "cheese," to create \mathbf{x}_q with duration of 603 ms, shown at top in 316 Fig. 9. We decompose this signal using MP and the 8xMDCT dictionary [31]. 317 We distort the query in two ways: with additive WGN (AWGN), and with 318 an interfering sound having a high correlation with the dictionary. In the first 319 case, shown in the middle in Fig. 9, the signal $\mathbf{x}'_q = (\alpha \mathbf{x}_q + \mathbf{n})/||\alpha \mathbf{x}_q + \mathbf{n}||$ is 320 the original \mathbf{x}_q distorted by a unit-norm AWGN signal \mathbf{n} . We set $\alpha=0.3162$ 321 such that $10\log_{10}(||\alpha \mathbf{x}_q||^2/||\mathbf{n}||^2) = 20\log_{10}(|\alpha|) = -10$ dB. For this signal, 322 we find the following statistics from 10,000 realizations of the AWGN signal: 323 $\mathrm{E}[|\langle \mathbf{x}_q, \mathbf{n} \rangle|] \approx 1 \times 10^{-5}, \ \mathrm{Var}[|\langle \mathbf{x}_q, \mathbf{n} \rangle|] \approx 4 \times 10^{-6}.$ We also find the following 324 lowing statistics for the 8xMDCT dictionary: $E[\max_{\mathbf{d} \in \mathcal{D}} |\langle \mathbf{n}, \mathbf{d} \rangle|] \approx 5 \times 10^{-4}$, $\operatorname{Var}\left[\max_{\mathbf{d}\in\mathcal{D}}|\langle\mathbf{n},\mathbf{d}\rangle|\right]\approx 2\times 10^{-5}$. Thus, the noise signal is not well-correlated 326 either with the original signal or the 8xMDCT dictionary. In the second case,

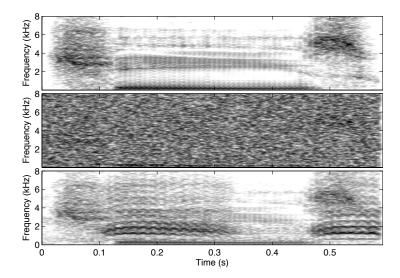


Figure 9: Log spectrograms of the query signals with which we search. Top: query of male saying "cheese." Middle: query distorted with additive white Gaussian noise (AWGN) with SNR = -10 dB. Bottom: query distorted with interfering crow sound with SNR = -5 dB.

shown at the bottom of Fig. 9, we distort the signal by adding the sound of 328 a crow **c** so that $\mathbf{x}'_q = (\alpha \mathbf{x}_q + \mathbf{c})/||\alpha \mathbf{x}_q + \mathbf{c}||_2$ with $||\mathbf{c}|| = 1$. Here, we set 329 $\alpha = 0.5623$ given by $20 \log_{10}(|\alpha|) = -5$ dB. For this interfering signal, we find that $|\langle \mathbf{x}_q, \mathbf{c} \rangle| \approx 2 \times 10^{-3}$, but $\max_{\mathbf{d} \in \mathcal{D}} |\langle \mathbf{c}, \mathbf{d} \rangle| \approx 0.21$, which is higher than $\max_{\mathbf{d}\in\mathcal{D}}|\langle \mathbf{x}_q,\mathbf{d}\rangle|\approx 0.17$. In this case, unlike for the AWGN interference, it is 332 likely that the sparse approximation of the signal with the crow interference will 333 have atoms in its low-order model due to the crow and not the speech. We do 334 not expect the AWGN interference to be a part of the signal model created by 335 MP until much later iterations. 336

Fig. 10 shows $|S_{iq}(t, M)|$ (28) aligned with the original signal for four values of M using the sparse approximations of the clean and distorted signals. We plot at the rear of these figures the localized magnitude time-domain correlation of the windowed and normalized signal with the query \mathbf{x}_q . In Fig. 10(a), using the clean \mathbf{x}_q , we clearly see its position even when using a single atom pair for each 100 ms partition of the time-domain. We see the same behavior in Fig. 10(b)—

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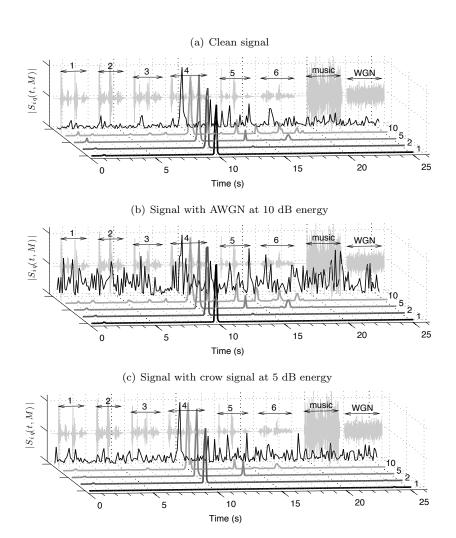


Figure 10: $|S_{iq}(M,t)|$ (28) as a function of time and the number of atoms M (labeled at right) considered from each representation for each localized sparse approximation. Localized magnitude correlation of each signal with query is shown by the thin black line in front of the gray time-domain signal at rear.

(c) for the two distorted signals, but in the case where the crow sound interferes we find the query for $M \geq 2$, or with at least three atom pair comparisons. The first atom of the decomposed query with the crow is modeling the crow and not the content of interest, and so we must increase the order of the model to find the location of \mathbf{x}_q . As we increase the number of pairs considered we also find

	Clean Signal			Signal + WGN		Signal + Crow			
#	t (s)	$ S_{iq} $	content	t (s)	$ S_{iq} $	content	t (s)	$ S_{iq} $	content
1	10.0	0.798	"cheese"	10.0	0.236	"cheese"	10.0	0.409	"cheese"
2	13.6	0.199	"cheese"	13.6	0.080	"cheese"	13.6	0.060	"cheese"
3	11.3	0.153	"-ives is-"	15.1	0.051	"delicious"	16.9	0.030	"cheese"
4	16.9	0.149	"cheese"	11.3	0.045	"-ives is-"	6.9	0.012	"cheese"
5	15.1	0.141	"delicious"	16.9	0.042	"cheese"	1.3	0.011	"cheese"
6	18.3	0.076	"delicious"	18.3	0.028	"delicious"	18.3	0.010	"delicious"
7	1.3	0.057	"cheese"	8.1	0.014	"delicious"	13.2	0.010	"cottage"
8	8.1	0.035	"delicious"	12.0	0.012	"-licious"	15.1	0.009	"delicious"
9	2.4	0.026	"delicious"	5.2	0.011	"delicious"	16.0	0.004	"cott-"
10	6.9	0.024	"cheese"	6.8	0.010	"cheese"	22.8	0.003	WGN

Table 2: Times, values and signal content for first 10 peaks in $|S_{iq}(t, 10)|$ (P(10) = 55) in Figs. 10(a)–(c). Highest-rated distances in each (bold) points to the origin of signal.

other segments that point in the same direction as \mathbf{x}_q . Table 2 gives the times and content of the ten largest values in $|S_{iq}(t,10)|$. For the clean and AWGN distorted \mathbf{x}_q , "cheese" appears five of the six times it exists in the original signal. Curiously, these same five instances are the five smallest distances when \mathbf{x}_q has the crow interference.

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We perform the same test as above but using a much longer speech signal (about 21 minutes in length) excerpted from a book-on-CD, "The Old Man and the Sea" by Ernest Hemingway, read aloud by a single person. From this signal we create several queries \mathbf{x}_q , from words to sentences to an entire paragraph of duration 35 s. We decompose each signal over the dictionary in Table 1 until the global residual energy is 20 dB below the initial energy. The approximation of the entire 21 m signal has 1,004,001 atoms selected from a dictionary containing 2,194,730,297 atoms.

One \mathbf{x}_q we extract from the signal is the spoken phrase, "the old man said" (861 ms in length). This phrase appears 26 times in the long excerpt. We evaluate $|S_{iq}(t, M)|$ (28) every 116 ms, and find the time at which \mathbf{x}_q originally appears using only M = 1 atom pair comparisons for each time partition. The

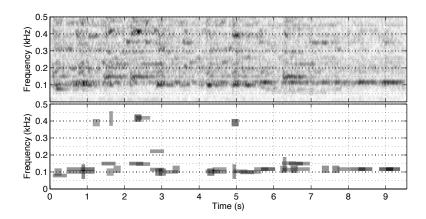


Figure 11: Polyphonic orchestral query: sonogram (top) and time-frequency tiles (bottom) of 50-order sparse approximation.

next highest ranked positions have values of 75% and 67% that of the largest $|S_{iq}(t,1)|$. When M=50, the values of the second and third largest values $|S_{iq}(t,50)|$ drop to 62% and 61% that of the largest value. In the top 30 ranked 367 subsequences for M=5 we find only one of the other 25 appearances of "the old 368 man said" (rank 26); but we also find "the old man agreed" (rank 11), and "the 369 old man carried" (rank 16). All other results have minimal content similarity to 370 the signal, but have time-frequency overlap in parts of the atoms of each model. 371 We perform the same test with a sentence extracted from the signal, "They 372 were as old as erosions in a fishless desert" (2.87 s), which only appears once. No 373 matter the M = [1, 50] we use, the origin of the excerpt remains at a rank of 6 374 with a value $|S_{iq}(t,50)|$ at 67.5% that of the highest rank subsequence. We find 375 that if we shift the time partition forward by 11.6 ms its ranking jumps to first, 376 with the second ranked subsequence at 73%. We observe a similar effect for a 377 query consisting of an entire paragraph (35 s). We find its origin by comparing 378 M=2 or more atoms from each model using a time partition of 116 ms. This result, however, disappears when we evaluate $|S_{iq}(t, M)|$ using a coarser time 380 partition of 250 ms.

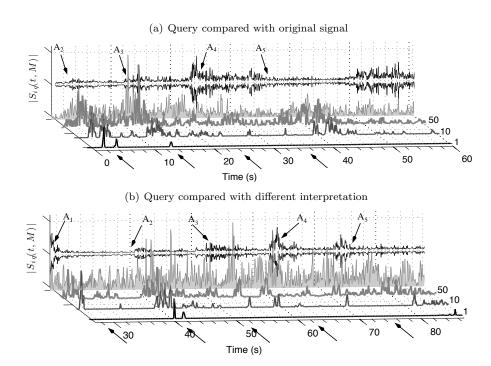


Figure 12: $|S_{iq}(t, M)|$ (28) for three values of M for the query and two different signals. Arrows mark the appearances of the 'A' theme, and their appearance number. Magnitude correlation of query with localized and normalized signal is shown by the solid gray area in front of the black time-domain signal at rear.

4.3. Experiment 3: Comparing Music Signals

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While the previous experiment deals with single-channel speech signals, in this experiment we make comparisons between polyphonic musical signals excerpted from a commercial recording of the fourth movement of *Symphonie Fantastique* by H. Berlioz. For the query, we use a 10.31 s segment of the third appearance of the 'A' theme of the movement (bars 33 - 39, located around 13 - 22 s in Fig. 4(b)). Figure 11 shows the sonogram and time-frequency tiles of the model of \mathbf{x}_q using the 50 atoms with the largest magnitude weights selected from the 8xMDCT dictionary [31]. We add no interfering signals as we do in the previous experiment.

Fig. 12(a) shows $|S_{iq}(t, M)|$ (28) over the first minute of the original signal, for three values of M, including M = 50, the time-frequency representation of

which is shown at bottom of Fig. 11. For $|S_{iq}(t,50)|$ we can see a strong spike located around 13 s corresponding with the query, but we also see spikes at 395 about 2 s and around 43 s. The former set of spikes correspond with the second appearance of the 'A' theme, when only low bowed strings are playing the theme 397 in G-minor. This is quite similar to the instrumentation of the query: low bowed 398 strings and a legato bassoon in counterpoint in Eb-major. The latter set of spikes 399 is around the end of the fifth appearance of the theme, which is played in G-400 minor on low pizzicato strings with a staccato bassoon. For M=10, we see a 40 conspicuous spike at the time of the fifth appearance around 34 s, as well as of 402 the fourth appearance around 24 s, where the theme is played in Eb-major like 403 the query. Finally, we test how the sparse approximation of this query compares 404 with subsequences from a different recording of this movement, which is also in 405 a different tempo. Figure 12(b) shows $|S_{iq}(t,M)|$ (28) for three different values of M. We see high similarity with the first and second appearances of the main 407 theme, but not the third, which is what the query contains. 408

4.4. Discussion

There is no reason to believe that a robust and accurate speech or melody 410 recognition system can be created by comparing only the first few elements 411 of greedy decompositions in time-frequency dictionaries. What appears to be 412 occurring for the short signals, both the "cheese" and "the old man said," is 413 that the first few elements of their sparse and atomic decomposition create a prosodic representation that is comparable to others at the atomic level. For 415 the longer signals, such as sentences, paragraphs, and orchestral theme, a few 416 atoms cannot adequately embody the prosody, but we still see that by only 417 making a few comparisons we are able to locate the excerpted signal — as long 418 as the time partition is fine enough. This is due to the atoms of the models acting in some sense as a time-frequency fingerprint, an example of which is in 420 Fig. 11. Through the cosine distance, the relative time and frequency locations 421 of the atoms in the query and subsequence are being compared. Subsequences 422 that share atoms in similar configurations will be gauged closer to \mathbf{x}_q than those 424 that do not.

By using the cosine distance it is not unexpected that (28) will be extremely 425 sensitive to a partitioning of the time-domain. This comes directly from the 426 definition of the time-localized Gramian (23), as well as the use of a dictionary 427 that is not translation invariant. There is no need to partition the time axis 428 when using a parameterized dictionary if we assume that some of the atoms 429 in the model of \mathbf{x}_q will have parameters that are nearly the same as some of 430 those in the relevant localized sparse representations. In such a scenario, we can 43 search a sparse representation for the times at which atoms exist that are similar 432 in scale and modulation frequency to those modeling \mathbf{x}_q . Then we can limit our 433 search to those particular times without considering any uniform and arbitrary 434 partition of the time-domain. With non-linear greedy decomposition methods 435 such as MP and time-variant dictionaries, however, such an assumption cannot be guaranteed; but its limits are not yet well-known. 437

5. Conclusion

In this paper, we have extended and investigated the applicability of a 439 method of recursive nearest neighbor search [14, 15] for comparing audio signals using pairwise comparisons of model elements in a sparse domain. The multiscale descriptions offered by sparse approximation over time-frequency dictio-442 naries are especially attractive for such tasks because they provide flexibility in 443 making comparisons between data, not to mention a capacity to deal with noisy 444 signals. After extending this method to the task of comparing subsequences 445 of audio signals, we find that the strongest bound known for the remainder is too weak to quickly and efficiently reduce the search space. Our experiments 447 show, however, that by comparing elements of sparse models we can judge 448 with relatively few comparisons whether signals share the same time-frequency 449 structures, and to what degrees, although this can be quite sensitive to the time-domain partitioning. We also see that we can approach such comparisons hierarchically, starting from the most energetic content to the least, or starting 452

from the longest scale phenomenon to the shortest.

We are continuing this research in multiple directions. First, since we know 454 that the inner product matrix $G_{iq}(t)$ (23) will be very sparse for all t in time-455 frequency dictionaries, this motivates designing a tighter bound based on a 456 Laplacian distribution of elements in $G_{iq}(t)$ with a large probability mass ex-457 actly at zero. This bound would be much more realistic than that provided by 458 assuming the elements of the Gramian are distributed uniform (16). Another 459 part of the problem is of course that the sums in (9) and (28) are not such that 460 at step M the P(M) largest magnitude values of $\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}$ are actually being 46: summed. By our assumption in (4), we know that the decay of the magnitudes 462 of the elements in \mathbf{A}_{iq} will be quickest in diagonal directions, but dependent 463 upon the element position in the matrix. These diagonal directions are simply given by

$$\begin{bmatrix} \partial/\partial\gamma_i \\ \partial/\partial\gamma_q \end{bmatrix} m^{-\gamma_i} l^{-\gamma_q} = -m^{-\gamma_i} l^{-\gamma_q} \begin{bmatrix} \gamma_i/m \\ \gamma_q/l \end{bmatrix}$$
 (29)

where we now recognize that the weights of two different representations can decay at different rates. With this, we can make an ordered set of index pairs by

$$\Lambda = \{ (m, l)_{\lambda} : |[\mathbf{A}_{iq}]_{\lambda}| \ge |[\mathbf{A}_{iq}]_{\lambda+1}| \}_{\lambda=1,2,\dots,n_i n_q}$$
(30)

and define a recursive sum for $1 < m \le n_i n_q$

$$S_{iq}(m) \stackrel{\Delta}{=} S_{iq}(m-1) + [\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}]_{\Lambda_m}$$
(31)

setting $S_{iq}(1) = [\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}]_{11}$. We do not yet know the extent to which this approach can ameliorate the problems with the non-discriminating bound (16), as we have yet to design an efficient way to generate a satisfactory Λ , and estimate the bounds of the corresponding remainder — whether it is like that in (16), or another that uses the fact that $\mathbf{G}_{iq}(t)$ will be very sparse, even when $\mathbf{x}_q = \mathbf{y}_i$. We think that using a stronger bound and this indexing order will significantly reduce the number of pairwise comparisons that must be made

before determining a subsequence is not close enough with respect to the cosine distance. Furthermore, we can make the elements of \mathbf{A}_{iq} decay faster, and thus increase γ , by using other sparse approximation approaches, such as OMP [28, 23] or CMP [36]. And we cannot forget the implications of choosing a particular dictionary. In this work we have used two different parametric dictionaries, one of which is designed for audio signal coding [31]. Another interesting research direction is to use dictionaries better suited for content description than coding, such as content-adapted dictionaries [20, 2, 19].

Finally, and specifically with regards to the specific problem of similar-485 ity search in audio signals, the cosine distance between time-domain samples 486 makes little sense because it is too sensitive to signal waveforms whereas hu-487 man perception is not. Instead, many other possibilities exist for comparing 488 sparse approximation, such as comparing low-level histograms of atom parameters [7, 34]; comparing mid-level structures such as harmonics [9, 38, 8]; and 490 comparing high-level patterns of short atoms representing rhythm [32]. There 491 also exists the Matching Pursuit Dissimilarity Measure [25], where the atoms of 492 one sparse model are used to decompose another signal, and vice versa to see 493 how well they model each other. We are exploring these various possibilities 494 with regards to gauging more generally similarity in audio signals at multiple 495 levels of specificity within a sparse domain. 496

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Appendix A. Proof of Remainder Bounds

To show (14), we can bound R(M) loosely by assuming the worst case scenario of $[\mathbf{G}_{iq}]_{ml} = 1$ for all its elements. Knowing that R(M) is the sum of the elements of the matrix $\mathbf{A}_{iq} \bullet \mathbf{G}_{iq}$ except for the first P(M) values, and assuming (4), we can say

$$C^{-2}R(M) \le \sum_{m=M+1}^{n} \sum_{l=1}^{m} [l(m-l+1)]^{-\gamma} + \sum_{m=1}^{n-1} \sum_{l=m+1}^{n} [l(n-m+1)]^{-\gamma}$$
$$= ||\mathbf{c}_{M}^{\gamma}||_{1} + ||\mathbf{d}^{\gamma}||_{1}. \quad (A.1)$$

where \mathbf{c}_{M}^{γ} and \mathbf{d}^{γ} are defined in (17) and (18). This worst case scenario is not possible using MP because of its update rule (1).

We can find the tighter bound in (15) by assuming the distribution of signs of the elements of \mathbf{G}_{iq} is Bernoulli equiprobable, i.e., $P\{[\mathbf{G}_{iq}]_{ml}=1\}=P\{[\mathbf{G}_{iq}]_{ml}=-1\}=0.5$. Thus, defining a random variable $b_i:\mathbb{R}\mapsto\{-1,1\}$, and its probability mass function $f_B(b_i)=0.5\delta(b_i+1)+0.5\delta(b_i-1)$ using the Dirac function, $\delta(x)$, we create a random vector \mathbf{b} with $n^2-P(M)$ elements independently drawn from this distribution. Placing this into the double sums of (A.1) provides the bound

$$C^{-2}R(M) \le \left| \mathbf{b}^T \begin{bmatrix} \mathbf{c}_M^{\gamma} \\ \mathbf{d}^{\gamma} \end{bmatrix} \right| \le ||\mathbf{c}_M^{\gamma}||_1 + ||\mathbf{d}^{\gamma}||_1.$$
 (A.2)

This weighted Rademacher sequence has the property that [14]

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$$P\left\{\left|\mathbf{b}^{T}\mathbf{s}\right| > R\right\} \le 2\exp\left(-R^{2}/2||\mathbf{s}||_{2}^{2}\right), R > 0 \tag{A.3}$$

which becomes $P\{|\mathbf{b}^T\mathbf{s}| \leq R\} \geq \max\{0, 1 - 2\exp(-R^2/2||\mathbf{s}||_2^2)\}$ by the axioms of probability. With this we can find an R such that $P\{|\mathbf{b}^T\mathbf{s}| \leq R\}$ will be greater than or equal to some probability $0 \leq p \leq 1$, i.e.,

$$R(p) = \left(||\mathbf{c}_M^{\gamma}||_2^2 + ||\mathbf{d}^{\gamma}||_2^2 \right)^{1/2} \left[2 \ln \frac{2}{1-p} \right]^{1/2}. \tag{A.4}$$

This value can be minimized by choosing p=0, for which we arrive at the residual upper bound in (15). Note that even though we have set p=0, we still

have an unrealistically loose bound by the impossibility of MP of choosing two sets of atoms for which all entries of their Gramian \mathbf{G}_{iq} are in $\{-1,1\}$.

Finally, to show (16), we can model the elements of the Gramian as random variables, $u_i : \mathbb{R} \mapsto [-1, 1]$, independently and identically distributed uniformly

$$f_U(u_i) = \begin{cases} 0.5, & -1 \le u_i \le 1\\ 0, & \text{else.} \end{cases}$$
 (A.5)

Substituting this into (14) gives a weighted sum of random variables satisfying

$$C^{-2}R(M) \le \left| \mathbf{u}^T \begin{bmatrix} \mathbf{c}_M^{\gamma} \\ \mathbf{d}^{\gamma} \end{bmatrix} \right| \le ||\mathbf{c}_M^{\gamma}||_1 + ||\mathbf{d}^{\gamma}||_1.$$
 (A.6)

where **u** is the random vector. For large M, this sum has the asymptotic property [14, 15]:

$$P\{|\mathbf{u}^T\mathbf{s}| < R\} = \text{Erf}\sqrt{\frac{3R^2}{2||\mathbf{s}||_2^2}}.$$
 (A.7)

Setting this equal to $0 \le p \le 1$ and solving for R produces the upper bound (16). We can reach the upper bound (15) if we set p = 0.9586, but note that (16) can be made zero. This bound can still be extremely loose because the Gramian of two models in time-frequency dictionaries will be highly sparse.

Computing the ℓ_2 -norm in these expressions, however, leads to evaluating

Computing the ℓ_2 -norm in these expressions, however, leads to evaluating the double sums

$$||\mathbf{c}_{M}^{\gamma}||^{2} = \sum_{m=M+1}^{n} \sum_{l=1}^{m} \frac{1}{[l(m-l+1)]^{2\gamma}}$$
 (A.8)

$$||\mathbf{d}^{\gamma}||^{2} = \sum_{m=1}^{n-1} \sum_{l=m+1}^{n} \frac{1}{[l(n-m+1)]^{2\gamma}}$$
(A.9)

which can be prohibitive for large n. The dimensionality of \mathbf{c}_M^{γ} is n(n+1)/2 - P(M), and of \mathbf{d}^{γ} is n(n-1)/2. We approximate these values in the following way for $\gamma = 0.5$, using the partial sum of the harmonic series

$$\sum_{m=1}^{n} \frac{1}{m} = \ln n + \gamma_E + \frac{1}{2n} - \frac{1}{12n^2} + \frac{1}{120n^4} + \mathcal{O}(n^{-6})$$
 (A.10)

where $\gamma_E \approx 0.5772$ is the Euler-Mascheroni constant. To find $||\mathbf{d}^{0.5}||^2$

$$||\mathbf{d}^{0.5}||^2 = \sum_{m=1}^{n-1} \sum_{l=m+1}^n \frac{1}{l(n-m+1)}$$

$$= \sum_{m=1}^{n-1} \frac{1}{n-m+1} \left[\sum_{l=1}^n \frac{1}{l} - \sum_{l=1}^m \frac{1}{l} \right]$$

$$\approx \sum_{m=1}^{n-1} \frac{1}{n-m+1} \left[\ln n/m - \frac{n-m}{2nm} + \frac{n^2 - m^2}{12n^2m^2} \right]. \tag{A.11}$$

To find $||\mathbf{c}_{M}^{0.5}||^2$ we first use partial fractions and then the partial sum of the harmonic series:

$$||\mathbf{c}_{M}^{0.5}||^{2} = \sum_{m=M+1}^{n} \sum_{l=1}^{m} \frac{1}{l(m-l+1)}$$

$$= \sum_{m=M+1}^{n} \frac{1}{m+1} \sum_{l=1}^{m} \frac{1}{l} + \frac{1}{m-l+1}$$

$$\approx \sum_{m=M+1}^{n} \frac{1}{m+1} \left[\ln m + \gamma_{E} + \frac{1}{2m} - \frac{1}{12m^{2}} + \sum_{l=1}^{m} \frac{1}{l} \right]$$

$$\approx 2 \sum_{m=M+1}^{n} \frac{1}{m+1} \left(\ln m + \gamma_{E} + \frac{1}{2m} - \frac{1}{12m^{2}} \right). \tag{A.12}$$

With these expressions we can avoid double sums in calculating the bounds.

Appendix B. Estimating the Compressibility Parameters

The compressibility parameters (C, γ) must be estimated for the set of weights in $\widehat{\mathcal{Y}}$ (??), as well as those of \mathbf{x}_q . Since by (4) the parameters (C, γ) bound from above the decay of all the ordered weights, only the largest magnitude weights matter for their estimation. Thus, we define a vector, \mathbf{a} , of the largest n magnitude weights from each row in the set $\{\{\mathbf{a}_i(n_i)\}_{i\in\mathcal{I}}, \mathbf{a}_q(n_q)\}$, which is equivalent to taking the largest weights at each approximation order. Good compressibility parameters can be given by

$$\min_{C,\gamma} ||C\mathbf{z}^{\gamma} - \mathbf{a}||^2 + \lambda C \text{ subject to } C\mathbf{z}^{\gamma} \succeq \mathbf{a}$$
 (B.1)

where we define $\mathbf{z}^{\gamma} \stackrel{\Delta}{=} [1, 1/2^{\gamma}, \dots, 1/n^{\gamma}]^T$, and add a multiple of C in order to keep it from getting too large since the bounds (14)–(16) are all proportional to

it. The constraint is added to ensure all elements of the difference $C\mathbf{z}^{\gamma} - \mathbf{a}_i$ are positive such that (4) is true.

To remove the γ component from the exponent, and since all of the elements of **z** and **a** are positive and non-zero, we can instead solve the problem

$$\min_{C,\gamma} ||\ln C \mathbf{1} + \gamma \ln \mathbf{z} - \ln \mathbf{a}||^2 + \lambda \ln C$$

$$= \min_{C,\gamma} \left[(\ln C)^2 n + \gamma^2 ||\ln \mathbf{z}||^2 + ||\ln \mathbf{a}||^2 + \lambda \ln C + 2\gamma (\ln \mathbf{z})^T (\ln C \mathbf{1} - \ln \mathbf{a}) - 2 \ln C (\ln \mathbf{a})^T \mathbf{1} \right]$$
(B.2)

subject to the constraint $C\mathbf{z}^{\gamma} \succeq \mathbf{a}$. Taking the partial derivative of this with respect to γ and C, we find

$$\gamma_o = \frac{(\ln \mathbf{z})^T (\ln \mathbf{a} - \ln C\mathbf{1})}{||\ln \mathbf{z}||^2}$$
(B.3)

$$C_o = \exp\left[\lambda + \frac{1}{n} \sum_{i=1}^n [\ln \mathbf{a} - \gamma \ln \mathbf{z}]_i\right]. \tag{B.4}$$

Starting with some initial value of C then, we use the following iterative method

- 1. solve for γ given a C in (B.3);
- 2. find the new C in (B.4) using this γ ;
- 3. set $C' = \exp\left[\max(\ln \mathbf{a} \gamma_o \ln \mathbf{z})\right]$ and evaluate the error $||C'\mathbf{z}^{\gamma} \mathbf{a}||^2$;
- 4. repeat until the error begins to increase.
- The factor λ in effect controls the step size for convergence. A typical value we use is $\lambda = \pm 0.03$ based on experiments (the sign of which depends on if the objective function decreases with decreasing C).
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