CODE: Coherence Based Decision Boundaries for Feature Correspondence

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Abstract—A key challenge in feature correspondence is the difficulty in differentiating true and false matches at a local descriptor level. This forces adoption of strict similarity thresholds that discard many true matches. However, if analyzed at a global level, false matches are usually randomly scattered while true matches tend to be coherent (clustered around a few dominant motions), thus creating a coherence based separability constraint. This paper proposes a non-linear regression technique that can discover such a coherence based separability constraint from highly noisy matches and embed it into a correspondence likelihood model. Once computed, the model can filter the entire set of nearest neighbor matches (which typically contains over 90% false matches) for true matches. We integrate our technique into a full feature correspondence system which reliably generates large numbers of good quality correspondences over wide baselines where previous techniques provide few or no matches.

Index Terms—Feature matching, wide-baseline matching, visual correspondence, RANSAC

1 Introduction

Orrespondence between image pairs involve finding the projections of the same scene points in both images. By linking multiple images together, correspondence is a critical first input for many vision systems [1], [2], [3]. As applications vary wildly, correspondence algorithms must accommodate a wide range of baselines and scenarios, e.g. Internet images, noisy infrared images, high resolution images, low-resolution video frames, etc. In addition, the desired stability of downstream systems requires correspondence algorithms to find as many matches as possible while keeping false matches to a minimum. The simultaneous need for large numbers of matches, robustness and flexibility places a heavy strain on correspondence algorithms and has motivated intensive research along this direction.

To date, feature matching [4], [5], [6] is the correspondence solution of choice for many computer vision systems. While lacking the correspondence density offered by optical flow alternatives [7], [8], [9], [10], feature matchers provide an attractive blend of wide baselines, fast speed and fine localization. The goal of feature matching is to correspond sparsely scattered, distinctive key-points. Distinctiveness is enhanced by consolidating local patch information into transformation invariant descriptors and matching decisions are based on descriptor comparison. While generally effective, feature correspondence typically discards many true matches to suppress the number of false matches [11],

[12], [13]. This can cause a paucity of matches which negatively impacts downstream algorithms. The problem is especially severe at the extremes of feature correspondence's working range shown in Fig. 1A). Lowering match acceptance thresholds provides many more (sometimes by a few orders of magnitude [11]) true matches. However, false matches increase more rapidly in number, leading to the mess shown in Fig. 1B).

Despite appearances, matches in Fig. 1B) may actually be separable into true and false matches. This is because true matches tend to be coherent, with neighboring pixels sharing similar motions, while wrong matches are usually randomly scattered. This leads us to propose a novel approach termed COherence based DEcision boundaries or CODE, which computes a coherence based partition of the potential correspondence space into true and false regions. At first glance, this approach appears self-contradictory, since coherence estimation and thus the partition is intrinsically coupled with the unknown (or highly imperfectly estimated) correspondence. However, we observe that a coarse coherence estimate is sufficient as we can eventually rely on the feature's spatial localization for fine matching. This permits quasi-decoupling of the coherence and correspondence steps. In CODE, we develop a robust, nonlinear regression formulation by treating feature matches as noisy data points. This is used to model a coherence cost (or the likelihood) for every possible motion. As the regression model only needs to be coarsely estimated, it can be effectively approximated from very noisy point correspondences obtained from low acceptance thresholds. Once estimated, the model forms a coherence based decision boundary for verifying correspondence hypotheses. The result is efficient, coherence enforcement for feature correspondences.

The proposed approach has some distinctive theoretical and practical advantages. 1) Global motion modeling: By applying our optimization on the bilateral domain which contains both spatial and motion coordinates [x, y, u, v],

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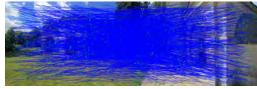
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A) Traditional A-SIFT feature matches



B) A-SIFT feature matches with lowered acceptance thresholds



C) A-SIFT feature correspondence with CODE regression

Fig. 1: Feature matching for an example image pair. A) A-SIFT [6], a highly regarded wide-baseline feature matcher. B) Relaxed acceptance thresholds result in many more true matches hidden in a pile of false matches. C) Our fusion of A-SIFT with the proposed technique termed *CODE*. *CODE* regression can model the motion adequately even under such noisy matches in B). Note that only the right side of the main arch structure is visible in the second image.

a smooth curve can model spatially discontinuous motion (conceptually similar to edge preserving bilateral filters [14]). This dispenses with the need to handle discontinuities by weakening the smoothing function or demanding explicit discontinuity detection. Instead, a global as-smoothas-possible function accommodates a wide variety of motions, making modeling both flexible and robust. 2) Computational tractability: We show that an as-smooth-as-possible regression function can be formulated as a convex cost with a guaranteed global minimum. This avoids initialization problems and facilitates mathematically elegant solutions. 3) Soft modeling: Given our input feature matches' noisy nature, modeling motion as a hard one-to-one mapping assignment between images is unlikely to be accurate. The proposed likelihood model expresses this ambiguity, and it adds a layer of robustness since (unlike in one-to-one modeling). As a result, a mistakenly high score for an incorrect motion need not affect the score of the true motion, and this allows graceful degradation on difficult scenes. 4) Efficiency: Naive correspondence likelihood modeling is potentially very expensive. If the likelihood of every possible motion for every pixel is stored as a matrix, even a small 640×480 image, with potential horizontal and vertical motion ranges of [-640, 640] and [-480, 480], would require a matrix of size $640 \times 480 \times (2 \times 640) \times (2 \times 480)$. If expressed as a likelihood based regression model, we can estimate every matrix entry with a few hundred variables. This makes correspondence likelihood modeling computationally tractable.

In practice, CODE consists of a number of smooth likeli-

hood decision boundaries for validating feature correspondences. These can be rapidly computed from noisy correspondences hypotheses (50% wrong matches are acceptable at this stage). Once computed, the decision boundaries can filter the entire nearest-neighbor matching set (which often has over 90% outliers) for inliers, with the one-to-one correspondence forming a natural complement to the likelihood based CODE boundaries. This process retains many more correspondences than traditional descriptor based thresholding methods while eliminating nearly all wrong matches. Our regression can be computed in a few seconds and since verification is cheap, the proposed technique scales efficiently to large numbers of features. In practice, CODE filtering time is O(N) where N is the pre-defined number of feature matches used in the regression.

To make a complete feature correspondence package, we fuse *CODE* with our re-implementation of GPU A-SIFT [15] and Muja and Lowe's [16] fast approximate nearest neighbor matching. As *CODE*'s overhead is low, our full system is faster than the original A-SIFT implementation [6], while procuring many more matches. An example is shown in Fig. 1C).

This current work extends our previous conference papers [11], [17]. In this paper, we provide a thorough exposition of the proposed technique, and improve its scalability to handle large numbers of matches through a theoretical approximation. In addition, we perform a number of new and expanded experiments to validate the proposed technique against a range of state-of-the-art methods. In summary, the main contributions are listed as follows:

- We propose *CODE*, a principled approach expressing the motion smoothness constraints as a matching likelihood estimate of every possible motion of every image pixel. *CODE* can be efficiently computed from highly noisy feature matches and forms an effective mechanics of discerning the difference between true and false matches.
- We integrate *CODE* with A-SIFT to create a feature matching system which yields high quality matches at wide baselines where previous techniques provide few or no matches. Despite the aggressive matching, it avoids matching images of different scenes/objects. We share our optimized implementation, in the form of executable files, to the research community at: http://www.kind-of-works.com/home/code.
- We demonstrate our high numbers of reliable correspondences bring compelling improvements when integrated into the existing structure-from-motion systems.

1.1 Related Work

This paper owes a significant debt to an array of pioneering research on affine invariant, wide-baseline features [6], [18], [19]. However, the original works could not fully exploit their features' invariance due to difficulty of identifying false matches. This fact forced these methods to adopt very strict similarity criteria which discarded many true match hypotheses. By modeling the true underlying motion from very noisy matches, our *CODE* technique can reliably differentiate true and false matches. This results in a flood of good matches shown in Fig. 1.

Similar to optical flow [7], [8], [9], [10], [20], we use motion smoothness to facilitate correspondence matching. However, we retain feature correspondence's design philosophy, trading matching density and fineness in favor for wide-baselines, higher speeds and robustness. As such our technique may ignore subtle motion details which might be retained by optical flow algorithms. If desired, density and fine motions may be recovered through a subsequent recomputation, but this is beyond the current paper's scope.

Our formulation builds on the motion coherence framework [21], [22]. Unlike the smoothness prior enforced in most optical flow algorithms which directly penalizes motion differences, motion coherence seeks the smoothest continuous motion field consistent with observed data. The global data integration has repeatedly demonstrated impressive levels of stability [22], [23], [24]. However, the original formulation does not accommodate motion discontinuities and is vulnerable to local minima. By adapting the motion coherence into a motion discontinuity-preserving, global regression technique, we avoid these problems while maintaining its robustness.

Edge-preserving bilateral filters [14] were a major inspiration for this work. However, unlike bilateral filtering which is a local operation, our regression on the bilateral domain computes a global model. This allows us to connect information from across the image for robustness to outliers. In addition, while bilateral filters return a single value at each pixel, *CODE* will return a continuous function encoding a distribution of likelihoods for all possible motion values a pixel can take.

We also draw inspiration from previous attempts to fuse smoothness constraints with a sparse feature correspondence. These include graph matching [13], [25], [26], [27], [28], bounded distortion [29] and mesh-based correspondence reasoning [30], with a number of approaches [29], [30] explicitly focused on removing false matches. However, the graph and distortion based techniques [29], [30] are vulnerable to local minima. Further, their computation cost increases with the number of features, making the techniques less scalable to high resolution imagery. While scalability and non-convexity are problems with most correspondence formulations, assuming pre-computed, albeit noisy correspondence creates a sub-problem which we show is amenable to convex regression modeling, thus avoiding both these issues.

In practical terms, a family of RANSAC techniques [31], [32], [33], including the recent branch-and-bound formulation [34], share our goal of removing false matches. However, the basic RANSAC formulation is designed for small (e.g. 10-20 variables) linearizable models, making them application specific. Our general smoothness based constraints cannot be enforced with RANSAC, thus motivating our investigation of regression techniques. In practice, it is best to use our method to boost the number of true matches first, then followed by RANSAC to estimate application specific parameters when appropriate.

Also worth mentioning are region growing correspondence algorithms [35], [36] which provide quasi-dense correspondence while handling large baselines and occlusions. These techniques provide many more correspondences than ours. However our solution retains the feature based techniques are region growing correspondences.

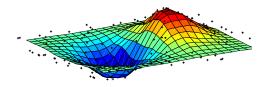


Fig. 2: Regression can be understood as finding a continuous surface that explains scattered data points (denoted by "+").

niques' innate advantage in speed, correspondence accuracy and handling large scale changes.

Finally we emphasize that different from the objectives of SIFT flow [37] or object matching [38], we do not desire correspondence between different objects. Given two images of physically different objects, our desired result is no correspondence.

2 OUR APPROACH

This section outlines our general formulation for coherence based data regression. We begin by designing a multifunction regression in Sect. 2.1 based on motion coherence [21], [22]. Sect. 2.2 explains the bottlenecks of this approach and proposes an accelerated approximation. We provide a succinct summary of the regression process in Sect. 2.3. Based on the theoretical foundation laid in this section, Sect. 3 applies the regression techniques to the correspondence problem.

2.1 Generalized Coherence for Data Regression

The problem is formulated as fitting a smooth function to observed data points. The fitting function is $f: \mathbf{p} \mapsto q$, where $\mathbf{p} \in \mathbb{R}^D$ and q are its domain and co-domain. We assume $f(\mathbf{p})$ is a linear combination of K smooth functions $\{f_k(\mathbf{p})\}_{k=1,2,\ldots,K}$, such that

$$q = f(\mathbf{p}) = \sum_{k=1}^{K} a_k(\mathbf{p}) f_k(\mathbf{p}) , \qquad (1)$$

where $a_k(\mathbf{p})$ are known weighting functions over the D-dimensional domain of \mathbf{p} , which provide formulation flexibility we exploit later.¹

The *observed data* consists of N noisy scalar values $\{\hat{q}_j\}$ at corresponding locations $\{\mathbf{p}_j\}$, which are assumed to be noisy observations of $f(\mathbf{p})$, such that

$$\hat{q}_j = f(\mathbf{p}_j) + n_j = \sum_{k=1}^K a_k(\mathbf{p}_j) f_k(\mathbf{p}_j) + n_j ,$$
 (2)

with n_j representing noise. With respect to Fig. 2, \mathbf{p} is a spatial point on the two-dimensional X-Y domain. The regression function $f(\mathbf{p})$ is represented by a continuous valued surface, while \hat{q}_j is the observed value for a spatial point \mathbf{p}_j .

Each individual $f_k(.)$ function is composed of two terms:

$$f_k(\mathbf{p}) = H_k + \phi_k(\mathbf{p}) . (3)$$

1. The formulation is unchanged if q is a vector. However for computational speed, q is a scalar throughout this paper.

 H_k is an (optional) unknown scalar offset and $\phi_k(\mathbf{p})$ is a smooth function evaluated with a motion coherence smoothness penalty [21], [22]

$$\Psi_k = \int_{\mathbb{R}^D} \frac{|\overline{\phi}_k(\omega)|^2}{\overline{g}(\omega)} d\omega , \qquad (4)$$

where $\overline{\phi}_k(.)$ denotes the Fourier transform of a function $\phi_k(.)$, while $\overline{g}(\omega)$ is the Fourier transform of a Gaussian function with a spatial standard deviation γ . Hence, Eqn. (4) achieves smoothness by penalizing high frequency terms.

Our goal is to find the smoothest possible set of $f_k(.)$ functions consistent with the given data points $\{\mathbf{p}_j, \hat{q}_j\}$. This is expressed as the following energy minimization:

$$E = \sum_{j=1}^{N} C \left(\hat{q}_j - f(\mathbf{p}_j) \right) + \lambda \sum_{k=1}^{K} \Psi_k$$

$$= \sum_{j=1}^{N} C \left(\hat{q}_j - \sum_{k=1}^{K} a_k(\mathbf{p}_j) f_k(\mathbf{p}_j) \right) + \lambda \sum_{k=1}^{K} \Psi_k .$$
(5)

Here, C(.) represents the Huber function as used in [17]:

$$C(z) = Huber(z) = \begin{cases} z^2 & \text{if } ||z|| \le \epsilon \\ 2\epsilon ||z||_1 - \epsilon^2 & \text{if } ||z|| > \epsilon \end{cases}$$
 (6)

and λ is the weight given to the smoothness constraint Ψ_k .

Directly minimizing E with respect to functions $f_k(.)$ appears intractable as $\{f_k(.)\}$ are continuous functions. However, we can reduce the problem to an optimization over a finite number of variables as shown below.

Note the Fourier transform relation, $\phi_k(\mathbf{p}) = \int_{\mathbb{R}^D} \overline{\phi}_k(\omega) e^{2\pi\iota\langle \mathbf{p},\omega\rangle} d\omega$. We know that at the minimum of the energy E of Eqn. (5), its derivative is zero. Hence,

$$\frac{\delta E}{\delta \overline{\phi}_{k}(\mathbf{z})} = 0, \forall \mathbf{z} \in \mathbb{R}^{D}, k \in \{1, 2, \dots, K\}
\Rightarrow \sum_{j=1}^{N} \mathbf{w}_{k}(j) \int_{\mathbb{R}^{D}} \frac{\delta \overline{\phi}_{k}(\omega)}{\delta \overline{\phi}_{k}(\mathbf{z})} e^{2\pi\iota \langle \mathbf{p}_{j}, \omega \rangle} d\omega
+ \lambda \int_{\mathbb{R}^{D}} \frac{\delta}{\delta \overline{\phi}_{k}(\mathbf{z})} \frac{|\overline{\phi}_{k}(\omega)|^{2}}{\overline{g}(\omega)} d\omega = 0
\Rightarrow \sum_{j=1}^{N} \mathbf{w}_{k}(j) e^{2\pi\iota \langle \mathbf{p}_{j}, \mathbf{z} \rangle} + 2\lambda \frac{\overline{\phi}_{k}(-\mathbf{z})}{\overline{g}(\mathbf{z})} = 0$$
(7)

where \mathbf{w}_k is a $N \times 1$ vector that serves as a place-holder for more complicated terms.

Rearranging the terms gives

$$\overline{\phi}_k(\mathbf{z}) = \overline{g}(-\mathbf{z}) \sum_{j=1}^N \mathbf{w}_k(j) e^{-2\pi \iota \langle \mathbf{p}_j, \mathbf{z} \rangle} , \qquad (8)$$

Taking the inverse Fourier transform of Eqn. (8), we can write our continuous functions $\{\phi_k(\mathbf{p})\}$ in terms of a finite set of N-dimensional vectors $\{\mathbf{w}_k\}$,

$$\phi_k(\mathbf{p}) = \sum_{j=1}^N \mathbf{w}_k(j) g(\mathbf{p}, \mathbf{p}_j) = \sum_{j=1}^N \mathbf{w}_k(j) e^{-\frac{\|\mathbf{p} - \mathbf{p}_j\|}{\gamma^2}}, \qquad (9)$$

$$\forall k \in \{1, 2, ..., K\},$$

where $g(\mathbf{p}, \mathbf{p}_j)$ is a Gaussian radial basis function and $\{\mathbf{w}_k(j)\}$ are unknown variables. Thus Eqn. (9) states that the smoothest possible functions $\{\phi_k(.)\}$ which minimize

the energy in Eqn. (5) must lie in the space spanned by N radial basis functions $\{g(\mathbf{p}, \mathbf{p}_i)\}$.

For the smoothness constraint, substituting Eqn. (8) into Eqn. (4) allows the continuous regularization function Ψ_k to be expressed in terms of \mathbf{w}_k

$$\Psi_k = \mathbf{w}_k^T G \mathbf{w}_k , \qquad k \in \{1, 2, ..., K\} ,$$
 (10)

where G is a symmetric matrix with its elements given as

$$G(i,j) = g(\mathbf{p}_i, \mathbf{p}_j) = e^{-\|\mathbf{p}_i - \mathbf{p}_j\|^2 / \gamma^2}.$$
 (11)

Substituting Eqns. (9) and (10) into Eqn. (5) yields

$$\arg \min_{\{f_k(\mathbf{p})\}} \sum_{j=1}^{N} C\left(\hat{q}_j - f(\mathbf{p}_j)\right) + \lambda \sum_{k=1}^{K} \Psi_k$$

$$= \arg \min_{\{f_k(\mathbf{p})\}} \sum_{j=1}^{N} C\left(\hat{q}_j - \sum_{k=1}^{K} a_k(\mathbf{p}_j) f_k(\mathbf{p}_j)\right) + \lambda \sum_{k=1}^{K} \Psi_k$$

$$= \arg \min_{\{\mathbf{w}_k, H_k\}} \sum_{j=1}^{N} C\left(\hat{q}_j - \sum_{k=1}^{K} a_k(\mathbf{p}_j) \left(H_k + \sum_{i=1}^{N} \mathbf{w}_k(i) g(\mathbf{p}_j, \mathbf{p}_i)\right)\right)$$

$$+ \lambda \sum_{k=1}^{K} \mathbf{w}_k^T G \mathbf{w}_k,$$
(12)

where the energy is dependent only on a finite number of variables, i.e., $\{\mathbf{w}_k\}$ and H_k . Since G is a Gram matrix [22], this makes the coherence term Ψ_k in Eqn. (10) convex. As the Huber loss function C(.) is also convex and the sum of convex functions is convex, the overall energy minimization problem in Eqn. (12) is convex. Therefore, a gradient descent minimization of Eqn. (12) will lead to a guaranteed global minimum

Based on the estimated variables $\{\mathbf{w}_k\}$ and H_k , the global regression function $f(\mathbf{p})$ is constructed from $\{f_k(\mathbf{p})\}$:

$$f(\mathbf{p}) = \sum_{k=1}^{K} a_k(\mathbf{p}) f_k(\mathbf{p}) = \sum_{k=1}^{K} a_k(\mathbf{p}) (H_k + \phi_k(\mathbf{p}))$$
$$= \sum_{k=1}^{K} a_k(\mathbf{p}) \left(H_k + \sum_{i=1}^{N} \mathbf{w}_k(i) g(\mathbf{p}, \mathbf{p}_i) \right).$$
(13)

This formulation mirrors that developed in our previous conference paper [17]. For applications of this global regression approach to data fitting and image warping, we encourage interested readers to refer to [17].

2.2 Accelerated Coherence Based Global Regression

Examining the formulation presented in Sect. 2.1, one can find that the length of \mathbf{w}_k vectors and hence the number of variables in Eqns. (8) and (12) increase linearly with N, the number of observed data points under consideration. This creates a computational burden when N is large. Motivated to address this challenge, we propose an approximation that decouples this linear computational dependency, with its full derivation given in Appendix A.

Consider $f_k(\mathbf{p})$ in Eqn. (13), the k-th smooth function to be estimated. If many data points in the input set $\{\mathbf{p}_i\}$

2. This may partially explain the success of radial basis functions in learning networks [39].

are adjacent, a good approximation of $f_k(\mathbf{p})$ can be given as follows.

$$f_k(\mathbf{p}) \approx \widetilde{f}_k(\mathbf{p}) = H_k + \sum_{j=1}^M \widetilde{\mathbf{w}}_k(j) g(\mathbf{p}, \widetilde{\mathbf{p}}_j) ,$$
 (14)

where $\tilde{\mathbf{w}}_k$ is an M-dimensional vector ($M \ll N$), and $\{\tilde{\mathbf{p}}_j\}$ are M representative cluster centroids distributed across the space occupied by the original points $\{\mathbf{p}_j\}$. These centroids are usually obtained by K-means clustering [40] on $\{\mathbf{p}_j\}$. If most of the \mathbf{p}_j points have near duplicates in the set $\{\tilde{\mathbf{p}}_j\}$, the function $f_k(\mathbf{p})$ will be able to closely approximate all possible values of the original $f_k(\mathbf{p})$ function in Eqn. (13).

Interestingly, the approximated regression functions $\{\tilde{f}_k(\mathbf{p})\}$ can be easily integrated into the global energy function (12), which becomes

$$\arg\min_{\{\widetilde{\mathbf{w}}_k, H_k\}} \sum_{j=1}^{N} C\left(\hat{q}_j - \sum_{k=1}^{K} a_k(\mathbf{p}_j) \widetilde{f}_k(\mathbf{p}_j)\right) + \lambda \sum_{k=1}^{K} \widetilde{\mathbf{w}}_k^T \widetilde{G} \widetilde{\mathbf{w}}_k ,$$

where \widetilde{G} is an $M \times M$ matrix, with $\widetilde{G}(i,j) = g(\widetilde{\mathbf{p}}_i, \widetilde{\mathbf{p}}_j)$ and the overall cost remains convex. Note that in this approximation, the values of λ and M are not coupled, i.e., the same λ can be properly used for different settings of M.

This clustering-based approximation is actually useful in accelerating many computer vision tasks, where closely located pixels create many redundant variables. Apart from this paper, we believe our approximation can also aid the original motion coherence [22] and its derivations [23], [24].

2.3 Summary of the Coherence Based Regression

To provide a succinct summary of the regression process, we give a recap of it here. Given N noisy scalar values $\{\hat{q}_j\}$ at corresponding locations $\{\mathbf{p}_j\}$, where \mathbf{p}_j is a D-dimensional vector, we can explain the observed data $\{\mathbf{p}_j, \hat{q}_j\}$ with a continuous function $\tilde{f}(\mathbf{p})$. The function $\tilde{f}(\mathbf{p})$ is composed of a linear combination of K smooth functions, which returns a value for each query point $\{\mathbf{p}\}$

$$q = \widetilde{f}(\mathbf{p}) = \sum_{k=1}^{K} a_k(\mathbf{p}) \widetilde{f}_k(\mathbf{p}) , \qquad (16)$$

where $\{a_k(\mathbf{p})\}$ are known weighting functions over the D-dimensional domain \mathbf{p} . The function $\widetilde{f}_k(\mathbf{p})$ is in turn parametrized by $\{\widetilde{\mathbf{w}}_k, H_k\}$ variables:

$$\widetilde{f}_{k}(\mathbf{p}) = H_{k} + \sum_{j=1}^{M} \widetilde{\mathbf{w}}_{k}(j)g(\mathbf{p}, \widetilde{\mathbf{p}}_{j})$$

$$= H_{k} + \sum_{j=1}^{M} \widetilde{\mathbf{w}}_{k}(j)e^{-\|\mathbf{p} - \widetilde{\mathbf{p}}_{j}\|^{2}/\gamma^{2}},$$
(17)

which can be estimated by a gradient descent minimization of the convex cost

$$\arg \min_{\{\widetilde{\mathbf{w}}_{k}, H_{k}\}} \sum_{j=1}^{N} C\left(\widehat{q}_{j} - \widetilde{f}(\mathbf{p}_{j})\right) + \lambda \sum_{k=1}^{K} \widetilde{\mathbf{w}}_{k}^{T} \widetilde{G} \widetilde{\mathbf{w}}_{k} =$$

$$\arg \min_{\{\widetilde{\mathbf{w}}_{k}, H_{k}\}} \sum_{j=1}^{N} C\left(\widehat{q}_{j} - \sum_{k=1}^{K} a_{k}(\mathbf{p}_{j}) \widetilde{f}_{k}(\mathbf{p}_{j})\right) + \lambda \sum_{k=1}^{K} \widetilde{\mathbf{w}}_{k}^{T} \widetilde{G} \widetilde{\mathbf{w}}_{k}.$$
(18)

A Huber loss function C(.) is adopted to evaluate the data fitting quality. A summary of the estimated variables and user-defined parameters are given in Table 1.

Estimated	$\{H_k\}$	optional bias variable in $f_k(\mathbf{p})$			
Estimatea	$\{\widetilde{\mathbf{w}}_k\}$	function variables in $f_k(\mathbf{p})$			
	λ	weight of smoothness terms			
	γ	std. deviation of smoothing Gaussian			
	ϵ	threshold in the Huber function $C(.)$			
User-defined	K	number of smooth functions for $\widetilde{f}(\mathbf{p})$			
	$\{a_k(\mathbf{p})\}$	user defined weighting function			
	M	length of $\widetilde{\mathbf{w}}_k$			
	$\{\widetilde{\mathbf{p}}_j\}$	representative subsampled points			
	\tilde{N}	number of data points			
Others	$\widetilde{f}(\mathbf{p})$	regression function in Eqn. (16)			
Others	cost	given in Eqn. (18)			

TABLE 1: Summary of parameters and variables in our multi-function regression in Eqns. (16), (17), and (18).

3 COHERENCE BASED DECISION BOUNDARIES

With the general formulation for coherence based data regression presented in Sect. 2, we now apply the regression techniques to the correspondence problem. The key idea is to estimate a coarse coherence based model from a sparse set of very noisy feature matches. This model quantifies the coherent matching evidence for every possible motion, providing a coherence based decision boundaries for verifying correspondence hypotheses. More specifically, we achieve this through two regression functions which act as a chain of cascaded filters for eliminating wrong feature matches. We describe individual modules below and elaborate on design decisions and properties in Sect. 3.1.

Matching likelihood boundaries: What constitutes coherent motion? We consider a local motion is coherent if it satisfies either of these criteria: a) a concentrated cluster of local features making the motion; and b) many features over a large spatial extent making the motion. A visual illustration is shown in Fig. 3.

To enforce this coherence, we formulate a regression based likelihood function in which each given feature match is a noisy *observed data* of the form:

$$\{\mathbf{p}_j = [\mathbf{x}_j, \mathbf{m}_j, \mathbf{x}_j + \mathbf{m}_j, \mathbf{o}_j], \ \hat{q}_j = 1\}$$
, (19)

where each match indexed by j hypothesizes a "1" value at location \mathbf{p}_j . Here $\mathbf{x}_j = [x_j, y_j]$ and $\mathbf{m}_j = [u_j, v_j]$ are two-dimensional vectors representing image coordinates and motion vectors respectively, while \mathbf{o}_j is a 4×1 vector representing the relative affine orientation of the matched feature descriptor. From Eqn. (17), we choose a single function K = 1 without the H_k bias variable, and set $a_k(\mathbf{p}) = 1, \forall \mathbf{p}$. This leads to a regression function

$$\widetilde{f}(\mathbf{p}) = \sum_{j=1}^{M} \widetilde{\mathbf{w}}(j) g(\mathbf{p}, \widetilde{\mathbf{p}}_j) .$$
 (20)

Substituting the regression function (20) into the cost (18), we attain

$$\arg\min_{\tilde{\mathbf{w}}} \sum_{j=1}^{N} C(1 - \tilde{f}(\mathbf{p}_{j})) + \lambda \tilde{\mathbf{w}}^{T} \tilde{G} \tilde{\mathbf{w}}, \qquad (21)$$

whose minimization estimates the parameters of $\widetilde{f}(\mathbf{p})$.

Observe that the smoothness cost $\tilde{\mathbf{w}}^T \tilde{G} \tilde{\mathbf{w}}$ in Eqns. (4) and (10) not only penalizes un-smooth motions but penalizes all motions by encouraging $\tilde{\mathbf{w}}$ to be zero. Thus, the lack

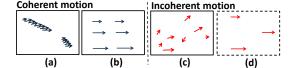


Fig. 3: Coherence based separation of true and false matches. Motions are considered coherent if (a) many local points make similar motions or (b) there is broad spatial support for the motion. This is enforced via the likelihood function in Eqn. (21). In contrast, feature matches in (c) and (d) do not give coherent motions, as the matches are not consistent in (c), while there are insufficient smoothly moving points to justify a long-range motion coherence model in (d).

of a bias term in $\widetilde{f}(\mathbf{p})$ given in Eqn. (20) means the likelihood function stays at zero, unless forced upward by the data. The likelihood function rises towards one if there exist a cluster of adjacent matches whose collective pull justifies an "un-smooth" peak. Alternatively, if a motion has significant supports across a wide area but with no strong local pull, the likelihood function can also rise to one, since a smooth surface over a large extent can be fitted for these matches by incurring a low smoothness penalty. In contrast, randomly scattered, incorrect matches do not exert sufficient collective pull and are ignored by the regression function. A simplified matching likelihood function is illustrated in Fig. 4.

From noisy feature correspondences, we use Eqn. (21) to fit a robust likelihood function given by $\widetilde{f}(\mathbf{p})$. The acceptance condition for a feature matching hypothesis \mathbf{p}_i is

$$accept(\mathbf{p}_i) = \text{true} \quad \text{if } \widetilde{f}(\mathbf{p}_i) > \epsilon_{likelihood}$$
 (22)

This mechanics can take very noisy data and remove most gross matching errors. However, it lacks fine spatial awareness and some erroneous matches will remain.

Bilaterally varying affine motion boundaries: The matching likelihood function can be considered as a blind cluster discovery mechanics. However, we know that correct motions not only cluster but tend to approximate a piecewise smoothly varying affine model [23]. We enforce this knowledge through a bilaterally varying affine regression function, which computes a motion likelihood by checking a hypothetical motion's consistency against its estimated model.

We focus on the \boldsymbol{X} motion direction first. Observed data take the form:

observed data =
$$\{\mathbf{p}_j, \hat{q}_j = x_j + u_j\}$$
, (23)

where x_j , u_j are obtained from the given feature correspondence. The definition of \mathbf{p}_j remains unchanged from the likelihood function formulation in Eqn. (19).

By setting K=3 in Eqn. (16) and setting $a_1(\mathbf{p})=x, a_2(\mathbf{p})=y, a_3(\mathbf{p})=1$, we have

$$\tilde{f}_x(\mathbf{p}) = \tilde{f}_1(\mathbf{p})x + \tilde{f}_2(\mathbf{p})y + \tilde{f}_3(\mathbf{p}),$$
 (24)

where each $\widetilde{f}_k(\mathbf{p}) = H_k + \sum_{j=1}^M \widetilde{\mathbf{w}}_k(j) g(\mathbf{p}, \widetilde{\mathbf{p}}_j)$ represents an affine motion parameter for the location \mathbf{p} . To estimate

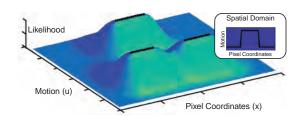


Fig. 4: Inset: An example set of motion hypotheses with discontinuities over the one-dimensional X axis. Main figure: The same data (black dots) over the bilateral domain u and X gives rise to a motion likelihood model built from the proposed regression technique. Observe that estimating a regression function over the bilateral domain involves estimating a value for every possible motion at every spatial location. Thus, a cross section along the specified X-position actually yields a likelihood function over the admissible motion range that a query X-position can take from. Note also that a smooth likelihood function computed for a motion-augmented bilateral domain can explain motion data with discontinuities.

 $f_k(\mathbf{p})$, from Eqn. (17), we optimize the following cost function

$$\arg\min_{\{\tilde{\mathbf{w}}_k, \tilde{H}_k\}} \sum_{i=1}^{N} C(x_i + u_i - \tilde{f}_x(\mathbf{p}_i)) + \lambda \sum_{k=1}^{3} \tilde{\mathbf{w}}_k \tilde{G} \tilde{\mathbf{w}}_k .$$
(25)

Similarly for the Y direction, we have a bilaterally varying affine model as follows,

$$\tilde{f}_y(\mathbf{p}) = \tilde{f}_4(\mathbf{p})x + \tilde{f}_5(\mathbf{p})y + \tilde{f}_6(\mathbf{p}),$$
 (26)

which can be computed from the cost

$$\arg\min_{\{\tilde{\mathbf{w}}_k, \tilde{H}_k\}} \sum_{i=1}^{N} C(y_i + v_i - \tilde{f}_y(\mathbf{p}_i)) + \lambda \sum_{k=4}^{6} \tilde{\mathbf{w}}_k \tilde{G} \tilde{\mathbf{w}}_k .$$
(27)

While not a direct likelihood function, the bilaterally varying affine model can be applied to distinguish correct and wrong feature matches through a thresholding step:

$$\begin{split} &accept(\mathbf{p}_i) = \text{true} \\ &\text{if } \sqrt{(\tilde{f}_x(\mathbf{p}_i) - x_i - u_i)^2 + (\tilde{f}_y(\mathbf{p}_i) - y_i - v_i)^2} < \epsilon_{spatial} \ . \end{split}$$

In practice, we find this stage gives more refined estimates but lacks the robustness of the preceding likelihood module. Thus, we cascade the affine motion boundary estimation after the likelihood boundary, which has removed most of the grossly incorrect matches.

3.1 Important Design Considerations and Properties

Having outlined our basic formulation, we will now delve into the properties and design considerations of our coherence based regression functions.

Global smoothness: It is enforced in all the functions. Unlike a local smoothness penalization which directly penalizes (motion) differences between neighbors, global smoothness enforced in our formulation seeks the globally smoothest function. By integrating information from across

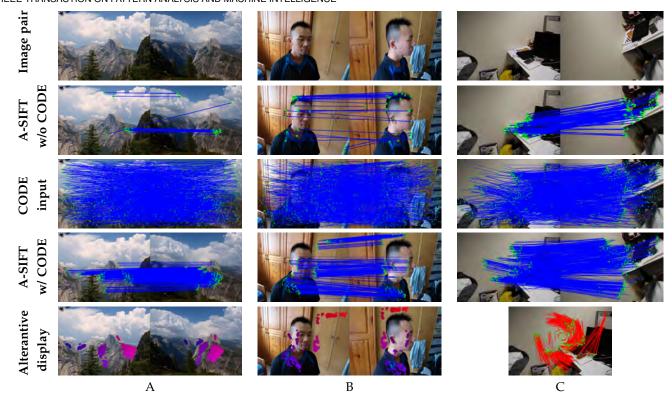


Fig. 5: *CODE* applied to A-SIFT features [6]. A and B demonstrate *CODE*'s ability to handle large motion discontinuities and occlusions, while C demonstrates its ability to handle large rotations. In alternative display, we color code feature matches according to their spatial locations on the left image. The matching transfers these color codes to the right image. To better illustrate the large rotation, C uses a vector display.

the image, global functions provide a high level of robustness, allowing the regression functions to be computed from very noisy input as shown in Fig. 5.

Domain choice: The choice of the domain for **p** in the assmooth-as-possible regression formulation can be arbitrary. We are motivated to choose the domain shown in Eqn. (19) out of three concerns:

Discontinuity preservation: Typically, global smoothness interferes with discontinuity preservation. Consider a function $f_k(\mathbf{p})$ where \mathbf{p} represents pixel coordinates, while the range of $f_k(\mathbf{p})$ represents motion. The smoothness (or infinite differentiability) implies a continuity constraint:

$$\lim_{\Delta \mathbf{p} \to 0} f_k(\mathbf{p} + \Delta \mathbf{p}) - f_k(\mathbf{p}) = 0, \quad \forall \mathbf{p} = [x, y]^T \in \mathbb{R}^2, \quad (29)$$

which forces the function value in the neighborhood of ${\bf p}$ to be similar. This means smooth functions must incur large errors at discontinuous motion boundaries.

We tackle this problem by using the bilateral domain spanning both the spatial and motion dimensions, i.e., $\mathbf{p} = [x,y,u,v]^T$, for conceptual understanding, we use a simpler form than that in (19). This redefines neighbors such that spatially neighboring points with different velocities are no longer adjacent. Therefore, we can assign very different function values to points with adjacent spatial coordinates, while retaining the constraint that $f_k(\mathbf{p})$ must be smooth. If the motion difference $(\Delta u, \Delta v)$ between two points in the bilateral domain \mathbf{p} tends to infinity, the point separation also tends to infinity, significantly reducing their influence on each other. This separation holds irrespective of the spatial

coordinates, and allows motion discontinuity to be accommodated by a smooth function shown in Fig. 4. Examples of this discontinuity handling are shown in Fig. 5 A, B).

Affine smoothness: It has been noted in [23] that because of the local generalization property of affines, motions are smoother when over-parameterized in the affine domain. Thus, we concatenate the affine parameters o to \mathbf{p} in (19). This allows us to better handle circular motions such as in Fig. 5 C).

Symmetry: Finally, for the sake of mathematical symmetry between left and right images, we concatenate the arguably redundant $\mathbf{x}_j + \mathbf{m}_j$ term to \mathbf{p} in (19).

Note that we find many different domain choices give good results and the domain's chosen in this paper serve as guidelines rather than canonical selections.

Multiple motion hypotheses: Computing regressions on the bilateral domain means estimating a function value for every $[x,y,u,v]^T$. This forces the function to consider every possible value that each pixel can take. Thus, unlike bilateral filters which estimate a single value for each pixel, CODE regression functions encode the likelihood for every possible motion each pixel can take. This soft-modeling adds a significant layer of robustness, as mistakenly giving a high score to an incorrect motion need not affect the score of the true motion. The computed regression functions and the resulting decision boundaries of CODE are subsequently used to verify feature correspondence hypotheses, with the hard, sub-pixel accurate feature matches forming a natural complement to the soft likelihood models.

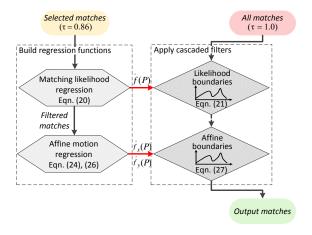


Fig. 6: Workflow of the proposed *CODE* technique. It consists of two major steps: building regression functions and applying cascaded filters, which take noisy feature matches obtained with different threshold values τ as the input. See the text for details.

Convex cost: The cost functions for estimating likelihood and bilaterally varying affine boundaries are convex. This avoids the need for initialization and allows our algorithm to handle challenging motions such as the large rotation shown in Fig. 5C).

3.2 Implementation

Our algorithm begins with feature matching. View-invariant features are computed through a GPU A-SIFT [15] with nearest-neighbor matching computed by FLANN [16]. Typically, feature matchers choose to accept matches based on a ratio test with a threshold setting $\tau=0.6$. i.e., a match is accepted if its nearest-neighbor difference is at least 0.6 times smaller than its second nearest neighbor. However, this can lead to few matches as shown in Fig. 1A). In this paper, we take *selected matches* obtained with a weaker threshold setting $\tau=0.86$ as the input. While matching results appear like a mess in Fig. 1B), the weaker threshold provides many true matches, from which information can be mined.

The selected matches are used to compute likelihood boundaries through the regression step in Eqn. (21). The current set is filtered by the likelihood boundaries and acceptable matches are used to compute bilateral affine motion boundaries in Eqns. (25) and (27). Once the regression functions and the resulting decision boundaries of CODE are computed, all matches obtained by setting $\tau=1.0$ are passed through the cascaded filters. The finally accepted feature matches are output matches. The minimizer used is the Ceres solver [41] and the process is illustrated in Fig. 6.

In theory, better results can be obtained through iterations, where the *output matches* form a new set of *selected matches*. In practice, we find the gain is small and hence restrict the study to a single iteration in this paper.

Parameter settings: To accommodate images of different sizes, the Hartley normalization is applied to the *selected matches* such that they have zero mean and average distance from the center of $\sqrt{2}$. The user-defined parameters in

Table 1 for the likelihood boundaries are $\lambda=1, \gamma=1, \epsilon=0.1, K=1, M=100, N=30,000$, with the set $\{\widetilde{\mathbf{p}}_j\}$ being centroids obtained from K-means clustering of the selected data. Matches are accepted with an $\epsilon_{likelihood}=0.6$ in Eqn. (22). The user-defined parameters for the bilateral affine boundaries are $\lambda=1, \gamma=1, \epsilon=0.1, K=3, M=100, N=1000$. Matches are accepted with an $\epsilon_{spatial}=0.01$ in Eqn. (28). The design details of the regression function $f(\mathbf{p})$ and weighting functions $\{a_k(\mathbf{p})\}$ are provided in Eqns. (20), (24) and (26).

4 EXPERIMENTAL RESULTS

To evaluate the effectiveness of the proposed *CODE* technique, we performed a series of experiments on different feature matching tasks and applications. Experimental results are reported in four subsections. Sect. 4.1 shows how the proposed regression technique *CODE* significantly improves the A-SIFT feature matcher [6]'s performance. Sect. 4.2 applies an A-SIFT with *CODE* system (A-SIFT w CODE) to the Structure from Motion problem. Sect. 4.3 compares our A-SIFT w CODE to other correspondence algorithms. Finally Sect. 4.4 discusses relative computational time of various correspondence approaches.

4.1 A-SIFT and CODE

CODE's central purpose is true-false feature match differentiation to retain ambiguous matches, which are often discarded in existing methods due to a conservative thresholding step. To evaluate effectiveness on real scenes, we computed A-SIFT, with and without CODE on twenty image pairs shown in Appendix D in the supplementary material. These images are grouped into sets A and B. Set A has moderate viewpoint changes but some image pairs with large illumination changes, while set B has very large viewpoint changes. All images are resized to a height of 600 pixels (to accommodate the original A-SIFT's limitation in image resolution).

We show feature matching results in Fig. 7. **Av. Precision** shows A-SIFT w CODE is more accurate than the original A-SIFT. A match is considered incorrect if its distance exceeds 7 pixels (on an image normalized to 640×480)³ from the fundamental matrix, and we used the implementation of RANSAC algorithms from Peter Kovesi [42]. As reflected from a comparison on **Av. Match Num**, CODE achieves higher precision while at the same time providing a few orders of magnitude more matches than the A-SIFT matcher. This illustrates the sheer number of potential correspondences that are discarded due to the inability to differentiate between true and false matches. The improvements are more marked at wide-baselines as can be seen by comparing sets A and B.

Apart from quantitative improvements, Fig. 7 also shows the proposed *CODE* approach produces qualitatively better results, yielding quasi-dense feature correspondences compared to the scattered matches of A-SIFT. **Area** % is an alternative measure of feature match numbers. Raw matching numbers are not so informative in themselves since an

3. We consistently set a threshold value of 7 as a compromise between the thresholds of 12 and 5 pixels used in [36] and [29], respectively.

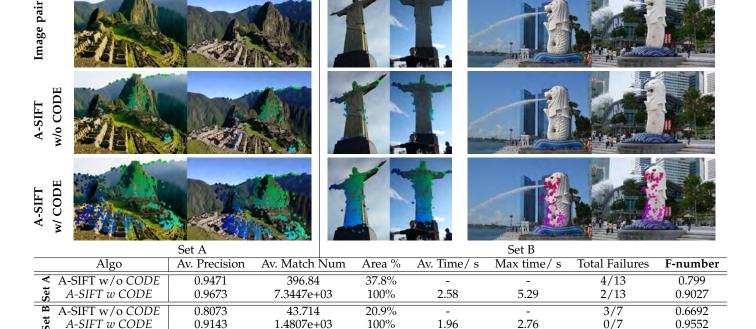


Fig. 7: Comparison between the original A-SIFT matcher [6] and *A-SIFT w CODE*, applying the proposed *CODE* regression to A-SIFT. For this test, we constructed a challenging dataset of static scene images. It includes representative images from a variety of prior papers [11], [13], [24], [29]. Set A has limited out-of-plane rotation, while Set B features very large viewpoint changes. *CODE* recovers many true feature matches that A-SIFT discards while maintaining high precision. This gain is especially noticeable for wide-baseline image pairs contained in Set B.

algorithm can interpolate and extrapolate to obtain more correspondences. To simulate this possibility, each match is dilated with a radius of 12 pixels (on its 640×480 rescaled image) to define a matching area. The total matching area covered by the A-SIFT is then measured against that of A-SIFT w CODE, and the resulting area ratio is reported as Area %. Even on this measure, CODE improves A-SIFT quite significantly and allows to cover an appreciably larger area. Note that this measure under-weights more difficult scenes which have small areas of overlap. Therefore, we also report Total Failures which gives the fraction of the tested scene pairs with no matches returned from an algorithm. In the absence of ground-truth, an approximate Recall is given as $(1 - Total\ Failures)$. Finally, F-Number is calculated as a summary statistic:

$$F - number = \frac{2 \times Recall \times Precision}{Recall + Precision}.$$
 (30)

Av. Time and **Max. Time** of 1.96 and 2.76 seconds reflect the computational overhead of integrating *CODE*. In fact, such a computational overhead is marginal when compared with the runtime of the original A-SIFT matcher (about 20 seconds) and our re-implementation (about 10 seconds). Runtimes are measured on a laptop with Intel i7CPU (2.4GHz), 8GB RAM and NVIDIA GPU GeForce GTX 660M.

4.2 Structure from Motion

In the last subsection, we performed a direct comparison between the A-SIFT matcher and our integration of CODE

in feature matching. The experiments justified *CODE* significantly improves A-SIFT. In this subsection, we choose to evaluate the two different feature matchers by integrating them into a complete end-to-end Structure from Motion (SfM) system, where feature matching is a fundamental early processing module.

Structure from Motion (SfM) seeks to infer 3D models from 2D images. It typically comprises two main stages: a) camera position computation where camera poses are inferred from feature matches such as A-SIFT, and b) a multi-view stereo reconstruction step, where images with known relative positions are fused into a full 3D model. An interesting fact is a test dataset, designed more for one of the stages, can be difficult for the other. For example, images from the multi-view stereo database [43] permit high quality dense multi-view reconstruction, if camera positions are assumed known; but the same set of images are less amenable to camera position recovery, due to difficulty in computing reliable feature correspondences. Hence, most SfM systems fail to reconstruct significant 3D scene sections. We attempted to reconstruct these scenes by integrating A-SIFT w CODE in the well-regarded Visual SfM system [3], [44], [45], [46], [47]. While reconstruction results are not always perfect, we can recover a complete 3D structure for all multi-view scenes. An example is shown in Fig. 8.

This experiment serves two purposes. First, it demonstrates that *CODE*'s high quality feature matching results can translate into meaningful improvements on other computer vision tasks. Second, feature correspondence is indeed critical to the final 3D model's quality. Our experiment



A set of multi-view images [43]



Visual SfM [3], [44], [45], [46], [47]



Agisoft [48]: A commercial 3D reconstruction software



Visual SfM using feature matches returned by A-SIFT w CODE

Fig. 8: Structure-from-motion seeks to fuse 2D images into 3D models. *CODE* recovers large numbers of reliable feature matches, making this task easier and more robust.

provides an indirect visualization of feature correspondence quality on challenging multi-view stereo datasets like [43]. Here each of the six image sequences has twenty-four images, leading to a total of 1656 pairwise matches. [Lu: Since the 3D reconstruction task here is computationally much heavier than all the other tests, we ran this comparative experiment on a desktop PC with an NVIDIA GeForce GTX980-Ti GPU.] The complete 3D reconstruction results for A-SIFT with and without *CODE* are shown in Appendix C in the supplementary material.

4.3 Comparison to Other Correspondence Algorithms

Thus far evaluations have focused on CODE's ability to facilitate A-SIFT feature correspondence. We now turn our attention to comparing our full A-SIFT w CODE to other matching techniques. In this section, we drop the cumbersome name A-SIFT w CODE and call our technique CODE. The most related algorithms are those which seek to remove or refine incorrect feature matches. Two recent examples are bounded distortion (BD) by Yaron et al. [29], and Mode Seeking by Chao et al. [13], [26]. At the time of writing, they represent a fair representation of the state-of-the-art. For experimental completeness, we also evaluate quasi-dense correspondence NRDC by HaCohen et al. [36] and agglomerative clustering ACC by Cho et al. [38]. As NRDC seeks dense correspondence, computation time on large images is prohibitively long. Because of this, unlike other algorithms, NRDC is run on down-sampled images of 640×480 resolution. As stated in [36], NRDC requires different settings for evaluation on the co-recognition dataset [49]. However, this impacts its general performance. Thus, we evaluate NRDC using its default parameters, and also NRDC(t) tuned for the co-recognition dataset. Next, we conduct evaluations for these competing algorithms on three datasets, and we present the evaluation protocol and experimental results for each of these datasets one by one.

1) Co-recognition dataset [49] consists of six image pairs where significant sections of the scene is re-shuffled.

It tests an algorithm's ability to handle large independent motions while providing large numbers of correspondence. Ground-truth segmentation boundaries are provided. A match is considered correct if it lies within a 7-pixel radius of the ground-truth segment boundary. The fraction of true matches is tabulated as Av. Precision in Fig. 9. We compute a correspondence area by dilating each match with a 12-pixel boundary [36]. The percentage of the total area overlapping with ground-truth segmentation is the Precision (area), a stricter measure than Av. Precision. Recall (area) is the percentage of ground-truth segmentation detected after dilation. F-number from Eqn. (30) provides a performance summary. When comparing computational time, we exclude the runtime of feature detection and nearestneighbor matching (our implementation is faster), and focus only on that of a core, correspondence evaluation/filtering algorithm. As NRDC does not use features, we provide its full computational time, and acknowledge this difference by indicating its runtime with a superscript '*'. A more detailed discussion on timing is provided in Sect. 4.4. We report all the results based on the above metrics in Fig. 9.

The area based statistics used for co-segmentation evaluation favors NRDC, a quasi-dense matching technique. However, our sparse feature matching method CODE remains surprisingly competitive, with its Recall (area) performance comparable to NRDC(t). In addition, it retains the computational speed and precision traditionally associated with feature matchers.

2) Our dataset consists of 20 images provided with the code of [13], [26], [29] with some additions. The images chosen have wide variations in scale, illumination, viewpoint change and image resolution. This is also the dataset used in Sect. 4.1, except that images are no longer downsampled. Image pairs are divided in two sets. Set A has no large out-of-plane rotational motion and is in the working range of all matching algorithms. Images in Set B have large out-of-plane rotational motions and cannot be meaningfully handled by [29] which utilizes SIFT features. Results are



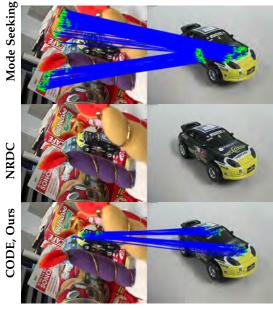
Algo	Av. Precision	Av. Match Num	Av. Time /s	Max time/ s	Total Failures	Recall (area)	Precision (area)	F-number (area)
BD [29]	0.81	31.833	10.17	18.01	0/6	0.0781	0.786	0.1421
Mode Seeking [13]	0.96	915.16	14.09	20.10	0/6	0.4445	0.877	0.5901
NRDC(t) [36]	0.95	3.565e+04	27.30*	30.75*	0/6	0.5096	0.776	0.6155
NRDC [36]	0.80	2.653e+04	25.12*	35.28*	0/6	0.3291	0.716	0.4510
ACC [38]	0.99	68	3.00	3.62	0/6	0.1901	0.934	0.3159
CODE, Ours	0.99	6.668e+03	2.94	3.43	0/6	0.5900	0.939	0.7249

Fig. 9: Evaluation on the co-recognition dataset [49]. We use a Recall (area) metric which favors dense correspondence over feature matchers. Despite this, *CODE* is surprisingly competitive at the co-segmentation task. This is a testament to *CODE*'s number of correspondences and ability to handle multiple motion models. Example results are visualized in the top panel.

tabulated in Fig. 10. For elaboration on individual statistics, please refer to Sect. 4.1. *CODE* works quite well across a wide variety of scenes as shown by the low number of *Total Failures*. Its matches are also of good quality as reflected by the over 90% precision.

3) ETHZ-toys [50], [51] consist of nine toys in different positions making a total of 40 object images. There are another 23 scene images where toys are hidden. In many scenes, the toy is, for instance, partially occluded, at a scale significantly different from that in its object image, distorted due to physical bending of the object, or having a glass encapsulation. We compute correspondence between every scene-toy image pair. If a scene contains a toy and a correspondence algorithm finds matches, it is considered a correct retrieval. If a scene does not contain a toy but an algorithm finds correspondences, it is considered an incorrect retrieval. Recall and precision statistics are computed for these trials and tabulated in Fig. 11. As processing 920 image pairs with all algorithms is prohibitively costly, we only compare to the top performers from the previous tests over the image sets A and B. The results in Fig. 11 show that CODE is general enough to accommodate large scale changes, non-rigid deformations and significant partial occlusion. It does so while avoiding correspondence on image pairs of different scenes. Over 920 image pairs, we had 0 incidence of incorrect retrieval and correctly retrieved 95.4% of the objects. This compares favorably to the next best algorithm, NRDC, with a precision and recall of 75% and 48.8% respectively.

To summarize the experiments, *CODE* demonstrates superior standings on a variety of challenging test cases, as a very competitive feature matching technique. In terms of precision, number of matches and computational time, it is consistently better than alternatives like *BD* [29] and *Mode Seeking* [13], [26]. Its match numbers is exceeded by quasidense matching *NRDC*. However, *CODE* can obtain reliable



Algo	Precision	Recall	F-number	
Mode Seeking [13]	0.2333	0.9767	0.3767	
NRDC [36]	0.7500	0.4884	0.5915	
CODE, Ours	1	0.9535	0.9762	

Fig. 11: The *ETHZ-toys* dataset [50], [51]. We measure the object-retrieval correctness. Mode Seeking [13] tends to find correspondence on image pairs even when there exists no common objects. This produces high recall but low precision. NRDC [36] represents a better balance. In contrast, our algorithm is aggressive in match retrieval but also strict in enforcing correctness, leading to a high F-number.





Algo	Av. Precision	Av. Match Num	Area %	Av. Time/ s	Max time/ s	Total Failures	F-number
BD [29]	0.84	616.76	41.96%	102.85	577.6	6/13	0.657
Mode Seeking [13]	0.86	606.07	40.64%	40.39	270.23	3/13	0.815
NRDC(t) [36]	0.95	4.4946e+04	105.8%	24.9*	29.6*	5/13	0.747
NRDC [36]	0.95	4.8556e+04	112.5%	24.7*	31.4*	4/13	0.803
ACC [38]	0.96	65	18.8%	3.68	6.07	5/13	0.751
CODE, Ours	0.96	4.41e+03	100%	2.97	6.27	1/13	0.940
Algo	Av. Precision	Av. Match Num	Area %	Av. Time/s	Max time/ s	Total Failures	F-number
BD [29]	0.61	14.1	13.5%	27.51	64.88	2/7	0.657
Mode Seeking [13]	0.84	416.4	22.07%	8.14	9.98	3/7	0.680
NRDC(t) [36]	0.94	1.843e+03	14.29%	25.69*	31.17*	5/7	0.439
NRDC [36]	0.96	7.985e+03	48.03%	21.93*	26.59*	2/7	0.822
ACC [38]	0.99	9.28	6.90%	1.87	2.60	5/7	0.444
CODE, Ours	0.92	4.41e+03	100%	1.73	2.05	0/7	0.960

Fig. 10: Evaluation of different feature matching techniques with the proposed *CODE*. A challenging image dataset and a set of metrics as described in Sect. 4.1 are used for this evaluation. The dataset contains full-resolution, representative images primarily from a variety of papers [13], [24], [26], [29]. Set A has limited out-of-plane rotation while Set B has very large viewpoint changes. *CODE* with constant parameters excels in a wide variety of scenarios as can be seen from the low number of *Total Failures* and high *Precision*.

feature correspondences on many scenes where *NRDC* has no matches as seen in *Total Failures* and *Recall* from Fig. 10 and Fig. 11, respectively.

4.4 Computational Time

It is true that computational efficiency is not the most critical issue when the primary concern is on estimating quality feature matches for an image pair. However, applications like Structure from Motion and image stitching typically require all-pair feature correspondence, where the complexity scales quadratically with the dataset size. As a result, an innocuous 30-second runtime scales to 50 minutes for all-pair matching on a small 10-image set. In this respect, feature matchers have a significant advantage over dense matchers as the heavy cost absorbed in the feature computation is only incurred once per image. This leaves a relatively low recurring time shown in Table 2. Observe that our theoretical time on a 10-image dataset is 12.4 minutes, while NRDC's is 45 minutes. In addition, our timings reported in Table 2 include many high resolution images, while NRDC used down-sampled 640×480 images. Such a clear runtime advantage enables a practical algorithm to leverage the enhanced accuracy offered by high resolution imagery to benefit challenging tasks such as Structure from Motion.

Next, we compare our computational time against other match verification or refinement alternatives like *BD* [29] and *Mode Seeking* [13], [26]. From Fig. 9 and Fig. 10, it can be seen that *CODE*'s average time is typically 5 to 6 times lower than alternatives. However, *CODE*'s maximum time can be one to two orders of magnitude lower, suggesting a scalability advantage. The picture is clearer in Fig. 12, which shows *CODE*'s computational cost scales well to handling large numbers of feature matches. This is due to

Time in seconds	Feature computa- tion	Match	CODE	Recur- ring time	Total time
A-SIFT w CODE	8.62	4.37	2.96	7.33	15.95
NRDC [36]	-	26.98	-	26.98	26.98

TABLE 2: Breakdown of average timing of Set A in Fig. 10. We compare our feature matcher with NRDC, a dense matcher. Note that while our algorithm is only about 2 times faster, its recurring time is 3 times faster than NRDC, leading to significant efficiency gains on image sets of moderate or large size. See the text for a detailed discussion.

CODE's generalization ability. As proposed in Sect. 2.2, the computationally expensive regression step can be computed on a subset of M matches, where M is pre-defined. The subsequent match verification step is fast and scales efficiently with the increasing number of input matches N ($N \gg M$). This makes the overall cost O(M).

5 CONCLUSION AND FUTURE WORK

We present a feature correspondence algorithm in which a smoothness based regression is utilized to identify correct matches based on their shared coherence. By reducing the strain on feature uniqueness, it provides many more matches over wider-baselines than previous solutions. These results are especially remarkable for a novel formulation and suggest a promising research direction. Our

4. Executable files and experimental results can be found at http://www.kind-of-works.com/.

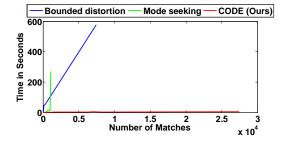


Fig. 12: CODE scales efficiently to high resolution images with many thousands of matches. The runtimes over different numbers of input feature matches N are plotted: red for CODE, blue for BD [29] and green for Mode Seeking [13], respectively.

current experiments are limited to sparse feature matching with an emphasis on speed and reliability. As such our solution ignores fine motion details and small independent motions. More visual illustrations are shown in Appendix B. However, by reliably modeling general motion, we believe our technique lays a solid foundation for subsequent works to enhance the density and fineness of motion estimates.

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APPENDIX A APPROXIMATE REGRESSION

This section provides an approximation that allows for rapid non-linear regression. We begin with a brief overview of the problem. In main body, the regression problem is reduced to an energy minimization given by Eqn. (12). Here, $\{\mathbf{p}_j, \hat{q}_j\}$ are N given data points. The minimization's goal is to fit a curve $f(\mathbf{p})$, which best explains the given data. As $f(\mathbf{p})$ is parametrized by $\{\mathbf{w}_k, H_k\}$ terms, the minimization in Eqn. (12) is

$$\arg\min_{\{\mathbf{w}_k, H_k\}} \sum_{j=1}^{N} C(\hat{q}_j - f(\mathbf{p}_j)) + \lambda \sum_{k=1}^{K} \mathbf{w}_k^T G \mathbf{w}_k , \quad (31)$$

where

$$G(i,j) = g(\mathbf{p}_i, \mathbf{p}_j) = \exp^{-\frac{(\mathbf{p}_i - \mathbf{p}_j)^2}{\gamma^2}},$$

$$f(\mathbf{p}) = \sum_{k=1}^K a_k(\mathbf{p}) f_k(\mathbf{p}),$$

$$f_k(\mathbf{p}) = H_k + \sum_{i=1}^N \mathbf{w}_k(i) g(\mathbf{p}, \mathbf{p}_i).$$
(32)

Here H_k is a scalar and \mathbf{w}_k is an N-dimensional vector.

We seek to approximate the cost of Eqn. (31) with a significantly shorter $\tilde{\mathbf{w}}_k$ vector by leveraging the fact that there are many similar $\mathbf{p}_i, \mathbf{p}_j$ data points. We assume without loss of generality that the last $\mathbf{p}_{N-t}, ..., \mathbf{p}_N$ data points are similar, and remove t-1 of them to create a shrunk dataset $\{\tilde{\mathbf{p}}_j\}$, with M=N-t+1 elements. We then approximate the curve $f_k(\mathbf{p})$ in Eqn. (32) with the shorter M-dimensional vector $\tilde{\mathbf{w}}_k$:

$$f_k(\mathbf{p}) \approx \tilde{f}_k(\mathbf{p}) = H_k + \sum_{i=1}^{M} \tilde{\mathbf{w}}_k(i) g(\mathbf{p}, \mathbf{p}_i)$$
 (33)

where

$$\begin{split} \tilde{\mathbf{w}}_k(i) &= \begin{cases} \mathbf{w}_k(i), & \text{if } i < M \ , \\ \sum_{j=M}^N \mathbf{w}_k(j), & \text{if } i = M \ . \end{cases} \\ \tilde{G}(i,j) &= g(\tilde{\mathbf{p}}_i, \tilde{\mathbf{p}}_j) & \text{where } \tilde{G} \text{ is a } M \times M \text{ matrix.} \end{cases} \tag{34}$$

The shorter $\tilde{\mathbf{w}}_k$ vector will approximate the original $f_k(\mathbf{p})$ function better if the values of $\mathbf{p}_{N-t},...,\mathbf{p}_N$ are more similar, and will do so perfectly if $\mathbf{p}_{N-t},...,\mathbf{p}_N$ are identical.

Replacing \mathbf{w}_k with $\widetilde{\mathbf{w}}_k$ not only affects the regression surface $f(\mathbf{p})$ but also the smoothness cost $\mathbf{w}_k^T G \mathbf{w}_k$ in Eqn. (31). This can be theoretically accommodated by changing the smoothness weight λ appropriately. Thus we seek to discover the interaction between λ and \mathbf{w}_k 's vector length.

We observe that because of the values taken by $\tilde{\mathbf{w}}_k$ in Eqn. (34),

$$\tilde{G}(i,:)\tilde{\mathbf{w}}_k \approx \begin{cases} G(i,:)\mathbf{w}_k, & \text{if } i < M, \\ \sum_{j=M}^{N} \mathbf{w}_k(j)G(i,:), & \text{if } i = M, \end{cases}$$
(35)

this leads to the following convenient equivalence

$$\mathbf{w}_k^T G \mathbf{w}_k \approx \tilde{\mathbf{w}}_k^T \tilde{G} \tilde{\mathbf{w}}_k . \tag{36}$$

In fact, the changes in $\tilde{\mathbf{w}}_k$ values counter balance the smaller \tilde{G} matrix, without need to modify λ in the end.

As in Eqn. (33), this approximation becomes perfect, if $\mathbf{p}_{N-t},...,\mathbf{p}_N$ are identical.

The result of the approximations in Eqn. (33) and Eqn. (36) is that the curve $f(\mathbf{p})$ in Eqn. (32) can be approximated as

$$f(\mathbf{p}) \approx \tilde{f}(\mathbf{p}) = \sum_{k=1}^{K} a_k(\mathbf{p}) \tilde{f}_k(\mathbf{p}) ,$$
 (37)

and the cost function in Eqn. (31) can be approximated as

$$\arg\min_{\{\tilde{\mathbf{w}}_k, H_k\}} \sum_{j=1}^{N} C(\hat{q}_j - \tilde{f}(\mathbf{p}_j)) + \lambda \sum_{k=1}^{K} \tilde{\mathbf{w}}_k^T \tilde{G} \tilde{\mathbf{w}}_k .$$
 (38)

While the above proof is for the consolidation of a single cluster, the same proof allows $\tilde{\mathbf{w}}_k$ to be formed from recursive consolidation of multiple clusters. This makes $\tilde{\mathbf{w}}_k$ substantially shorter than \mathbf{w}_k , and hence Eqn. (38) is much cheaper to minimize compared to Eqn. (31).

Note that recursive consolidation presented here is mainly for theoretical understanding. In practice, $\{\tilde{\mathbf{p}}_j\}$ is directly computed through K-means clustering [40] of $\{\mathbf{p}_j\}$ data points. Due to the finding in Eqn. (36), the size of the match clusters need not to be explicitly kept and only their centroid positions are used.

APPENDIX B NON-RIGID MOTION

One of the strengths of our *CODE* formulation is that it naturally accommodates handling large non-rigid motion scenes. However, *CODE*'s formulation means that small independent motions are invariably classified as incorrect and hence are ignored. Quantitative evaluation of this property is shown in Fig. 9, and more visual illustration is shown in Fig. 13.

APPENDIX C STRUCTURE FROM MOTION RESULTS

This sections tabulates all results for the 3-D reconstruction dataset [43], which are given in Fig. 14, Fig. 15, Fig. 16, Fig. 17, Fig. 18, and Fig. 19. For each dataset, reconstruction was achieved with VisualSfM [3], with A-SIFT and A-SIFT w CODE as a feature matcher. Apart from A-SIFT, we also applied other feature matchers to the dataset, but we find that nearly all previous techniques leave large gaps in the reconstruction. Our results are not perfect, still with notable errors in the Dinosaur and Matrix multi-view stereo image sequences in Fig. 16 and Fig. 17, but they represent a significant improvement over the prior state-of-the-art. For readers versed in Structure from Motion, we highlight that no prior camera calibration was assumed. Better reconstruction results can be obtained by using the focal length parameters provided with the dataset [43].

APPENDIX D OUR DATASET

Thumbnails of the twenty image pairs from our dataset are shown in Fig. 20.

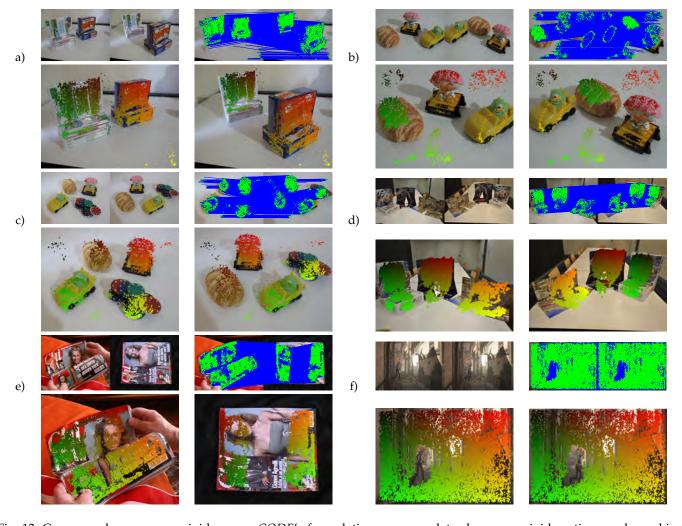


Fig. 13: Correspondence on non-rigid scenes. *CODE*'s formulation accommodates large non-rigid motions such as object permutation or folding shown in (a-e). However, *CODE* ignores small independent motions, because *CODE*'s formulation invariably considers them as wrong matches with low motion coherence. This is illustrated in (f). Depending upon the application, this can be an asset that facilitates motion segmentation or a limitation as small objects cannot be well tracked.



Fig. 14: Predator image sequence.

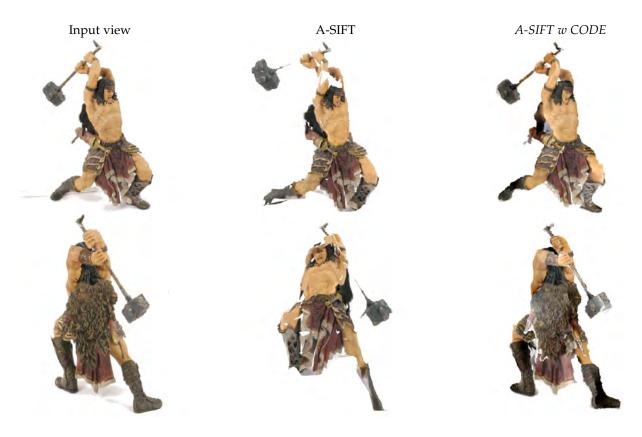


Fig. 15: Warrior image sequence.

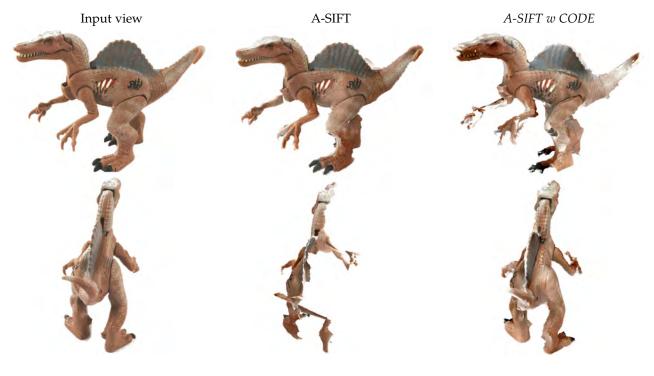


Fig. 16: *Dinosaur* image sequence.



Fig. 17: *Matrix* image sequence.

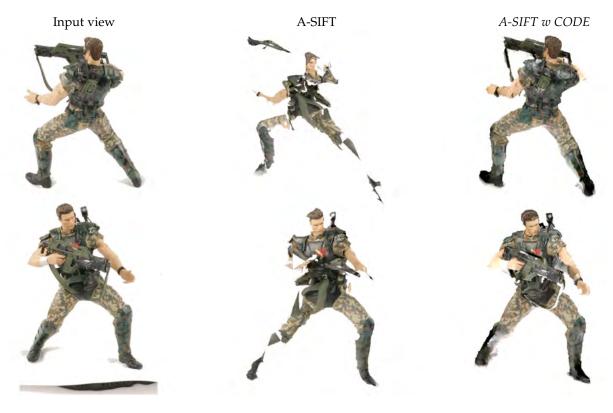


Fig. 18: *Soldier* image sequence.



Fig. 19: Mummy image sequence.

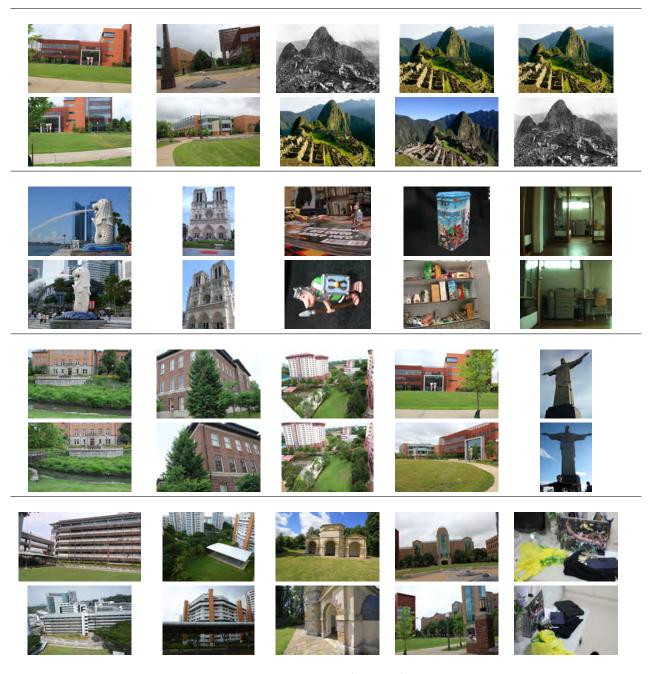


Fig. 20: Image pairs used in our dataset.