DISTANCE TRANSFORM GRADIENT DENSITY ESTIMATION USING THE STATIONARY PHASE APPROXIMATION

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Abstract. The complex wave representation (CWR) converts unsigned 2D distance transforms into their corresponding wave functions. Here, the distance transform S(X) appears as the phase of the wave function $\phi(X)$ —specifically, $\phi(X) = \exp\left(\frac{iS(X)}{\tau}\right)$ where τ is a free parameter. In this work, we prove a novel result using the higher-order stationary phase approximation: we show convergence of the normalized power spectrum (squared magnitude of the Fourier transform) of the wave function to the density function of the distance transform gradients as the free parameter $\tau \to 0$. In colloquial terms, spatial frequencies are gradient histogram bins. Since distance transform gradients carry only orientation information (as their magnitudes are identically equal to one almost everywhere), the 2D Fourier transform values mainly lie on the unit circle in the spatial frequency domain as $\tau \to 0$. The proof of the result involves standard integration techniques and requires proper ordering of limits. Our mathematical relation indicates that the CWR of distance transforms is an intriguing, new representation.

Key words. Stationary phase approximation; distance transform; gradient density; Fourier transform, complex wave representation

AMS subject classifications. 42B10; 41A60

1. Introduction. Euclidean distance functions (more popularly referred to as distance transforms) are widely used in many domains [17, 22]. An important subset—point-set based distance functions—also finds application in many domains with computer vision being a prominent example [22, 11, 20]. Since distance transforms allow us to transition from shapes to a scalar field, problems such as shape registration are often couched in terms of rigid, affine or nonrigid alignment of distance transform fields, where the shapes are parameterized as a set of points [19]. In medical imaging, they are used in the construction of neuroanatomical shape complex atlases based on an information geometry framework [2].

Even when one begins with a set of closed curves (as a shape template for example), the curves are often discretized to yield a point-set prior to the application of fast sweeping [25] and other distance transform estimation methods. Signed and unsigned distance transforms are deployed in 3D as well with their zero level-sets corresponding to surfaces. Furthermore, medial axis methods and skeletonization often involve distance transform representations [13]. In the domain of computer vision, the gradient density function is popularly known as the histogram of oriented gradients (HOG). Since the advent of HOG a few years ago, gradient density estimation has risen in prominence and is employed in human recognition systems [4].

The distance transform for a set of K discrete points $Y = \{Y_k \in \mathbb{R}^D\}, k \in \{1, \ldots, K\}$ where D is the dimensionality of the point-set is defined as

(1.1)
$$S(X) \equiv \min_{k} \|X - Y_k\|,$$

where $X \in \Omega$ is a closed bounded domain in \mathbb{R}^D . In this article, we are only concerned with D = 2.

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In computational geometry, Euclidean distance functions correspond to the Voronoi problem [5] and the solution S(X) can be visualized as a set of cones (with the centers being the point-set locations $\{Y_k\}$). The distance transform satisfies the static, non-linear Hamilton-Jacobi equation

$$(1.2) \|\nabla S\| = 1$$

almost everywhere, barring the point-set locations and the Voronoi boundaries where it is not differentiable [17, 18, 21]. Here $\nabla S = (S_x, S_y)$ denotes the gradients of Sand $\|\cdot\|$ represents its Euclidean magnitude. Furthermore S(X) = 0 at the point-set locations. Following the wave optics literature, one can envisage light waves simultaneously emanating from the given point sources and propagating with a velocity of one in all directions. The value of S at a grid point X_0 , namely $S(X_0)$, corresponds to the *time taken* by the first light wave (out of the K light waves) to reach the grid location X_0 . Driven by this optics analogy, when we express S as the phase of a wave function ϕ as in

(1.3)
$$\phi = \exp\left(\frac{iS}{\tau}\right),$$

we made an intriguing empirical observation. The power spectrum of the wave function approximates the density function of the gradients of the distance transform as the parameter τ in Equation 1.3 tends to zero. In this paper, we formally prove this result. We refer to this wave function ϕ which satisfies the phase relation with S as the *Complex Wave Representation* (CWR) of distance transforms.

2. Main Contribution. The centerpiece of this work is to provide a useful application of the stationary phase method, wherein we show an equivalence between the density function of the gradients of the distance function $\nabla S = (S_x, S_y)$ and the power spectrum (squared magnitude of the Fourier transform) of the CWR (ϕ) as the free parameter τ (in Equation 1.3) approaches zero. Here, the density function of the gradients is obtained via a random variable transformation of a uniformly distributed random variable W (over the bounded domain Ω) using the gradients $\nabla S = (S_x, S_y)$ as the transformation functions. In other words, if we define a random variable $Z = \nabla S(W)$ where the random variable W has a uniform distribution on a closed bounded domain $\Omega \subset \mathbb{R}^2$, the density function of Z represents the density function of the gradients of the distance transform.

As the norm of the gradients ∇S is defined to be 1 almost everywhere (from Equation 1.2), we observe that the density function of the gradients is one-dimensional and defined over the space of orientations. Section 3 provides a closed-form expression for this density function. As the gradients are unit vectors, we notice that the Fourier transform values of the CWR (ϕ) lie mainly on the unit circle and this behavior tightens as $\tau \to 0$. Specifically, if $F_{\tau}(\tilde{r}, \omega)$ represents the Fourier transform of ϕ in the polar coordinate system at a given value of τ , Theorem 4.2 demonstrates that if $\tilde{r} \neq 1$, then $\lim_{\tau \to 0} F_{\tau}(\tilde{r}, \omega) = 0$.

Our main result is established in Theorem 5.1 where we show that the power spectrum of the wave function ϕ when polled close to the unit circle, is approximately equal to the density function of the distance transform gradients, with the approximation becoming increasingly exact as $\tau \to 0$. In other words, if $P(\omega)$ denotes the closed-form density of the gradients defined over the orientation ω and if $P_{\tau}(\tilde{r}, \omega)$ corresponds to the power spectrum of ϕ represented in the polar coordinate system at a given value of τ , Theorem 5.1 constitutes the following relation

(2.1)
$$\lim_{\delta \to 0} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \left\{ \int_{1-\delta}^{1+\delta} P_{\tau}(\tilde{r},\omega) \tilde{r} d\tilde{r} \right\} d\omega = \int_{\omega_0}^{\omega_0 + \Delta} P(\omega) d\omega$$

for any (small) value of the interval measure Δ on ω . We show this result using the higher-order stationary phase approximation, a well known technique in asymptotic analysis [23]. Through the pioneering works of Jones and Kline [12], Olver [15], Wong [23], McClure and Wong [14], among others, the stationary phase approximation has become a widely deployed tool in the approximation of oscillatory integrals. Our work showcases a novel application of the stationary phase method for estimating the probability density function of distance transform gradients. The significance of our mathematical result is that spatial frequencies become histogram bins and hence the power spectrum P_{τ} can serve as a gradient density estimator at small, non-zero values of τ . We would like to emphasize that our work is fundamentally different from estimating the gradients of a density function [8] and should not be semantically confused with it.

2.1. Motivation from quantum mechanics. Our new mathematical relationship is motivated by the classical-quantum relation, wherein classical physics is expressed as a limiting case of quantum mechanics [10, 6]. When S is treated as the Hamilton-Jacobi scalar field, the gradients of S correspond to the classical momentum of a particle [9]. In the parlance of quantum mechanics, the squared magnitude of the wave function expressed either in its position or momentum basis corresponds to its position or momentum density respectively. Since these representations (either in the position or momentum basis) are simply (suitably scaled) Fourier transforms of each other, the squared magnitude of the Fourier transform of the wave function expressed in its position basis is its quantum momentum density. However, the time independent Schrödinger wave function $\phi(x, y)$ (expressed in its position basis) can be approximated by $\exp\left(\frac{iS(x,y)}{\tau}\right)$ as $\tau \to 0$ [6]. Here τ (treated as a free parameter in our work) represents Planck's constant. Hence the squared magnitude of the Fourier transform $\exp\left(\frac{iS(x,y)}{\tau}\right)$ corresponds to the quantum momentum density of S. The principal results proved in the article (Theorem 5.1 and Proposition 6.5) state that the classical momentum density (denoted by P) can be expressed as a limiting case (as $\tau \to 0$) of its corresponding quantum momentum density (denoted by P_{τ}), in agreement with the correspondence principle.

3. The Distance Transform Gradient Density Function. As mentioned above, the geometry of the distance transform corresponds to a set of intersecting cones with the origins at the Voronoi centers [5]. The gradients of the distance transform (which exist globally except at the cone intersections and origins) are unit vectors and satisfy Equation 1.2. Therefore the gradient density function is one-dimensional and defined over the space of *orientations*. The orientations are constant and unique along each ray of each cone. Its probability distribution function is given by

(3.1)
$$\mathcal{F}(\theta \le \Theta \le \theta + \Delta) \equiv \frac{1}{L} \int \int_{\theta \le \arctan\left(\frac{S_y}{S_x}\right) \le \theta + \Delta} dx dy$$

where L is the area of the bounded domain Ω . We have expressed the orientation random variable as $\Theta = \arctan\left(\frac{S_y}{S_x}\right)$. The probability distribution function also induces a *closed-form expression* for its density function as shown below.

Let $\Omega \subset \mathbb{R}^2$ denote a polygonal grid such that its boundary $\partial\Omega$ is composed of a finite sequence of straight line segments. The reason for restricting only to polygonal domains with boundaries made of line segments will become clear when we discuss Theorem 4.2. Let the set $Y = \{Y_k \in \mathbb{R}^2, k \in \{1, \ldots, K\}\}$ be the given point-set locations and let $Y_k = (x_k, y_k)$. Then the Euclidean distance transform at a point $X = (x, y) \in \Omega$ is given by

(3.2)
$$S(X) \equiv \min_{k} \|X - Y_k\| = \min_{k} (\sqrt{(x - x_k)^2 + (y - y_k)^2}).$$

Let \mathcal{D}_k , centered at Y_k , denote the k^{th} Voronoi region corresponding to the input point Y_k . \mathcal{D}_k can be represented by a Cartesian product $[0, 2\pi) \times [0, R_k(\theta)]$ where $R_k(\theta)$ is the length of the ray of the k^{th} cone at an orientation θ . If a grid point $X = (x, y) \in (Y_k + \mathcal{D}_k)$, then $S(X) = ||X - Y_k||$. Each \mathcal{D}_k is a convex polygon whose boundary $\partial \mathcal{D}_k$ is also composed of a finite sequence of straight line segments as shown in Figure 3.1.



FIGURE 3.1. Voronoi diagram of the given K points. Each Voronoi boundary is composed of straight line segments.

Note that even for points that lie on the Voronoi boundary where the radial length equals $R_k(\theta)$, the distance transform is well defined. The area L of the polygonal grid Ω is given by

(3.3)
$$L \equiv \sum_{k=1}^{K} \int_{0}^{2\pi} \int_{0}^{R_{k}(\theta)} r dr d\theta = \sum_{k=1}^{K} \int_{0}^{2\pi} \frac{R_{k}^{2}(\theta)}{2} d\theta.$$

With the above set-up in place, after recognizing the cone geometry at each Voronoi center Y_k , Equation 3.1 can be simplified as

(3.4)
$$\mathcal{F}(\theta \le \Theta \le \theta + \Delta) \equiv \frac{1}{L} \sum_{k=1}^{K} \int_{\theta}^{\theta + \Delta} \int_{0}^{R_{k}(\theta)} r dr d\theta = \frac{1}{L} \sum_{k=1}^{K} \int_{\theta}^{\theta + \Delta} \frac{R_{k}^{2}(\theta)}{2} d\theta.$$

Following this drastic simplification, we can write the closed-form expression for the density function of the unit vector distance transform gradients as

(3.5)
$$P(\theta) \equiv \lim_{\Delta \to 0} \frac{\mathcal{F}(\theta \le \Theta \le \theta + \Delta)}{\Delta} = \frac{1}{L} \sum_{k=1}^{K} \frac{R_k^2(\theta)}{2}.$$

Based on the expression for L in Equation 3.3, it is easy to see that

(3.6)
$$\int_0^{2\pi} P(\theta) d\theta = 1.$$

Since the Voronoi cells are convex polygons [5], each cell contributes exactly one conical ray to the density function on orientation.

4. Properties of the Fourier Transform of the CWR. Since the distance transform is not differentiable at the point-set locations $\{Y_k\}_{k=1}^K$ and also along the Voronoi boundaries $\partial \mathcal{D}_k, \forall k$ (a measure zero set in 2D), we restrict ourselves to the region which excludes both of them. To this end, let $0 < \epsilon < \frac{1}{2}$ be given. Let the region \mathcal{D}_k^{ϵ} centered at Y_k be represented by the Cartesian product $[0, 2\pi) \times [R_k^{(1)}(\theta), R_k^{(2)}(\theta)]$ where,

(4.1)
$$\begin{aligned} R_k^{(1)}(\theta) &= \epsilon R_k(\theta) \quad \text{and} \\ R_k^{(2)}(\theta) &= (1-\epsilon) R_k(\theta). \end{aligned}$$

The length of the ray at the orientation θ in \mathcal{D}_k^{ϵ} equals $R_k^{(2)}(\theta) - R_k^{(1)}(\theta)$. Note that in the definition of \mathcal{D}_k^{ϵ} , we have explicitly removed the source point Y_k where the ray length $r(\theta) = 0$ and the boundary of the Voronoi cell where $r(\theta) = R_k(\theta)$ as shown in Figure 4.1.



FIGURE 4.1. Region that excludes both the source point and the Voronoi boundary.

Define the grid

(4.2)
$$\Omega^{\epsilon} \equiv \bigcup_{k=1}^{K} \left(Y_k + \mathcal{D}_k^{\epsilon} \right).$$

Its area L^{ϵ} equals

(4.3)
$$L^{\epsilon} \equiv \sum_{k=1}^{K} \int_{0}^{2\pi} \int_{R_{k}^{(1)}(\theta)}^{R_{k}^{(2)}(\theta)} r dr d\theta = (1 - 2\epsilon) \sum_{k=1}^{K} \int_{0}^{2\pi} \frac{R_{k}^{2}(\theta)}{2} d\theta.$$

From Equation 3.3 we get $L^{\epsilon} = (1 - 2\epsilon)L$ and hence $\lim_{\epsilon \to 0} L^{\epsilon} = L$. Let $l^{\epsilon} = \sqrt{L^{\epsilon}}$. Define a function $F^{\epsilon} : \mathbb{R} \times \mathbb{R} \times \mathbb{R} \to \mathbb{C}$ as

(4.4)
$$F^{\epsilon}(u,v,\tau) \equiv \frac{1}{2\pi\tau l^{\epsilon}} \iint_{\Omega^{\epsilon}} \exp\left(\frac{iS(x,y)}{\tau}\right) \exp\left(\frac{-i(ux+vy)}{\tau}\right) dxdy.$$

For a fixed value of τ , define a function $F^{\epsilon}_{\tau}: \mathbb{R} \times \mathbb{R} \to \mathbb{C}$ as

(4.5)
$$F_{\tau}^{\epsilon}(u,v) \equiv F^{\epsilon}(u,v,\tau).$$

Note that F_{τ}^{ϵ} is closely related to the Fourier transform of the CWR, $\phi = \exp\left(\frac{iS}{\tau}\right)$ [1]. The scale factor $\frac{1}{2\pi\tau l^{\epsilon}}$ is the normalization factor such that the ℓ_2 norm of F_{τ}^{ϵ} is 1 as seen in the following Lemma (with the proof given in Appendix A).

LEMMA 4.1. With F^{ϵ}_{τ} defined as above, $F^{\epsilon}_{\tau} \in L^2(\mathbb{R}^2)$ and $||F^{\epsilon}_{\tau}|| = 1$.

Consider the polar representation of the spatial frequencies (u, v) namely $u = \tilde{r} \cos(\omega)$ and $v = \tilde{r} \sin(\omega)$ where $\tilde{r} > 0$. For $(x, y) \in (Y_k + \mathcal{D}_k^{\epsilon})$, let $x - x_k = r \cos(\theta)$ and $y - y_k = r \sin(\theta)$ where $r \in [R_k^{(1)}(\theta), R_k^{(2)}(\theta)]$. Then Equation 4.4 can be rewritten as

(4.6)
$$F_{\tau}^{\epsilon}(\tilde{r},\omega) = \sum_{k=1}^{K} C_k I_k(\tilde{r},\omega)$$

where

(4.7)
$$C_k = \exp\left\{-\frac{i}{\tau}\left[\tilde{r}\cos(\omega)x_k + \tilde{r}\sin(\omega)y_k\right]\right\}$$

and

(4.8)
$$I_k(\tilde{r},\omega) = \frac{1}{2\pi\tau l^\epsilon} \int_0^{2\pi} \int_{R_k^{(1)}(\theta)}^{R_k^{(2)}(\theta)} \exp\left\{\frac{i}{\tau} r \left[1 - \tilde{r}\cos(\theta - \omega)\right]\right\} r dr d\theta.$$

With the above set-up in place, we have the following theorem, namely, THEOREM 4.2. [Circle Theorem] If $\tilde{r} \neq 1$, then,

(4.9)
$$\lim_{\tau \to 0} F_{\tau}^{\epsilon}(\tilde{r}, \omega) = 0,$$

for any $0 < \epsilon < \frac{1}{2}$.

4.1. An Intuitive Examination of Theorem 4.2. Before we furnish a rigorous proof for the aforementioned theorem, we provide an intuitive picture of why the statement is true. Observe that the first exponential $\exp\left(\frac{iS(x,y)}{\tau}\right)$ in Equation 4.4 is a varying complex "sinusoid" and the second exponential $\exp\left(\frac{-i(ux+vy)}{\tau}\right)$ in Equation 4.4 is a fixed complex sinusoid at frequencies $\frac{u}{\tau}$ and $\frac{v}{\tau}$ along the x-and y-coordinate axes respectively. When we multiply these two complex exponentials, at low values of τ , the two sinusoids are usually not "in sync" and cancellations occur in the integral. Exceptions to the cancellation happen at locations where $\nabla S = (S_x, S_y) = (u, v)$, as around these locations, the two sinusoids are in perfect sync. Since $\|\nabla S\| = 1$ for distance transforms, strong resonance occurs only when $u^2 + v^2 = 1$ ($\tilde{r} = 1$). When $\tilde{r} \neq 1$, the two sinusoids tend to cancel each other out as $\tau \to 0$, resulting in F_{τ}^{ϵ} becoming zero at these locations.

4.2. Proof of Theorem 4.2. Having given an intuitive picture of why Theorem 4.2 holds true, we now proceed with the formal proof. As each C_k is bounded, it suffices to show that if $\tilde{r} \neq 1$, then $\lim_{\tau \to 0} I_k(\tilde{r}, \omega) = 0$ for all I_k .

Proof. Consider the integral

(4.10)
$$I(\tilde{r},\omega) = \frac{1}{2\pi\tau l^{\epsilon}} \int_{0}^{2\pi} \int_{R^{(1)}(\theta)}^{R^{(2)}(\theta)} \exp\left\{\frac{i}{\tau} r \left[1 - \tilde{r}\cos(\theta - \omega)\right]\right\} r dr d\theta,$$

where $R^{(1)}(\theta) = \epsilon R(\theta)$ and $R^{(2)}(\theta) = (1-\epsilon)R(\theta)$. Let the region $[0, 2\pi) \times [R^{(1)}(\theta), R^{(2)}(\theta)]$ be denoted by \mathcal{D}^{ϵ} . $R(\theta)$ is defined in such a way that the boundary of D^{ϵ} consists of a finite sequence of straight line segments as in the case of each $\mathcal{D}^{\epsilon}_{k}$. Notice that \mathcal{D}^{ϵ} doesn't contain the origin (0, 0). In order to prove Theorem 4.2, it is sufficient to show that $\lim_{\tau \to 0} I(\tilde{r}, \omega) = 0$.

Let $p(r, \theta; \tilde{r}, \omega) = r(1 - \tilde{r}\cos(\theta - \omega))$ denote the phase term of *I* in Equation 4.10 for a given \tilde{r} and ω . The partial derivatives of $p(r, \theta; \tilde{r}, \omega)$ (with \tilde{r} and ω held fixed) are given by

(4.11)
$$\frac{\partial p}{\partial r} = 1 - \tilde{r}\cos(\theta - \omega), \quad \frac{\partial p}{\partial \theta} = r\tilde{r}\sin(\theta - \omega).$$

Since \mathcal{D}^{ϵ} is bounded away from the origin (0,0), ∇p is well-defined and bounded and equals zero only when $\tilde{r} = 1$ and $\theta = \omega$. Since $\tilde{r} \neq 1$ by assumption, no stationary point exists $(\nabla p \neq 0)$ and hence we can expect $I(\tilde{r}, \omega) \to 0$ as $\tau \to 0$ [3, 12, 24]. Below, we show this result more explicitly.

Below, we show this result more explicitly. Define a vector field $\mathbf{u}(r,\theta;\tilde{r},\omega) = \frac{\nabla p}{\|\nabla p\|^2}r$ at a fixed value of \tilde{r} and ω . Note that

(4.12)

$$\nabla \cdot \left[\mathbf{u}(r,\theta;\tilde{r},\omega) \exp\left(\frac{ip(r,\theta;\tilde{r},\omega)}{\tau}\right) \right] = \left(\nabla \cdot \mathbf{u}(r,\theta;\tilde{r},\omega)\right) \exp\left(\frac{ip(r,\theta;\tilde{r},\omega)}{\tau}\right) + \frac{i}{\tau} \exp\left(\frac{ip(r,\theta;\tilde{r},\omega)}{\tau}\right) r$$

where the gradient operator $\nabla = \left(\frac{\partial}{\partial r}, \frac{1}{r}\frac{\partial}{\partial \theta}\right)$. Inserting Equation 4.12 in Equation 4.10, we get

(4.13)
$$I(\tilde{r},\omega) = I^{(1)}(\tilde{r},\omega) - I^{(2)}(\tilde{r},\omega),$$

where

$$I^{(1)}(\tilde{r},\omega) = \frac{1}{2\pi i l^{\epsilon}} \iint_{\mathcal{D}^{\epsilon}} \nabla \cdot \left(\mathbf{u}(r,\theta;\tilde{r},\omega) \exp\left(\frac{ip(r,\theta;\tilde{r},\omega)}{\tau}\right) \right) dr d\theta, \text{ and}$$

$$(4.14) \quad I^{(2)}(\tilde{r},\omega) = \frac{1}{2\pi i l^{\epsilon}} \iint_{\mathcal{D}^{\epsilon}} \left(\nabla \cdot \mathbf{u}(r,\theta;\tilde{r},\omega) \right) \exp\left(\frac{ip(r,\theta;\tilde{r},\omega)}{\tau}\right) dr d\theta.$$

Consider the integral $I^{(1)}(\tilde{r}, \omega)$. From the divergence theorem, we have

(4.15)
$$I^{(1)}(\tilde{r},\omega) = \frac{1}{2\pi i l^{\epsilon}} \int_{\Gamma} (\mathbf{u}^T \mathbf{n}) \exp\left(\frac{i p(r,\theta;\tilde{r},\omega)}{\tau}\right) ds$$

where Γ is the positively oriented boundary of \mathcal{D}^{ϵ} , s is the arc length of Γ and \mathbf{n} is the unit outward normal of Γ . The boundary Γ consists of two disjoint regions, one along $r(\theta) = R^{(1)}(\theta)$ and another along $r(\theta) = R^{(2)}(\theta)$. If the level curves of $p(r, \theta; \tilde{r}, \omega)$ are tangential to Γ only at a discrete set of locations giving rise to stationary points of the second kind [23, 24, 14]—in other words, if $p(r, \theta; \tilde{r}, \omega)$ is not constant along the boundary Γ for any contiguous interval of θ —then, using the one dimensional stationary phase approximation [15, 16], $I^{(1)}(\tilde{r}, \omega)$ can be shown to be $O(\sqrt{\tau})$ and hence converges to zero as $\tau \to 0$. Since the boundary of \mathcal{D}^{ϵ} is composed of straight line segments (specifically not arc-like), we can show that the level curves of $p(r, \theta; \tilde{r}, \omega)$ cannot overlap with Γ for a non-zero finite interval. (The next paragraph takes care of this technical issue and can be skipped without loss of continuity.)

The level curves of $p(r,\theta;\tilde{r},\omega)$ are given by $R(\theta)(1-\tilde{r}\cos(\theta-\omega)) = c$, where c is a constant. Recall that each of the two disjoint regions of Γ is composed of a finite sequence of line segments. For the level curves of $p(r,\theta;\tilde{r},\omega)$ to coincide with Γ over a non-zero finite interval, $y(\theta) = R(\theta)\sin(\theta) = \frac{c\sin(\theta)}{1-\tilde{r}\cos(\theta-\omega)}$ and $x(\theta) = R(\theta)\cos(\theta) = \frac{c\cos(\theta)}{1-\tilde{r}\cos(\theta-\omega)}$ should satisfy the line equation y = mx + b for some slope m and slope-intercept b, when θ varies over some contiguous interval $\theta \in [\theta_1, \theta_2]$. Plugging in the value of $y(\theta)$ and $x(\theta)$ into the line equation and expanding $\cos(\theta-\omega)$, we have

(4.16)
$$c\sin(\theta) = mc\cos(\theta) + b - b\tilde{r}[\cos(\theta)\cos(\omega) + \sin(\theta)\sin(\omega)].$$

Combining the terms, we get

(4.17)
$$\sin(\theta)[c+b\tilde{r}\sin(\omega)] - \cos(\theta)[mc-b\tilde{r}\cos(\omega)] = b$$

By defining $\lambda_1 \equiv c + b\tilde{r}\sin(\omega)$ and $\lambda_2 \equiv -(mc - b\tilde{r}\cos(\omega))$, we see that $\sin(\theta)$ and $\cos(\theta)$ need to satisfy the linear relation

(4.18)
$$\lambda_1 \sin(\theta) + \lambda_2 \cos(\theta) = b$$

for $\theta \in [\theta_1, \theta_2]$ in order for the level curves of $p(r, \theta; \tilde{r}, \omega)$ to overlap with the piecewise linear boundary Γ . As Equation 4.18 cannot be true for a finite interval of θ , $I^{(1)}(\tilde{r}, \omega) = O(\sqrt{\tau})$ as $\tau \to 0$ and hence converges to zero in the limit.

Now $I^{(2)}(\tilde{r},\omega)$ has a similar form as the original $I(\tilde{r},\omega)$ in Equation 4.10 with r replaced by $g_1(r,\theta;\tilde{r},\omega) = (\nabla \cdot \mathbf{u})$. Letting $\mathbf{u}_1(r,\theta;\tilde{r},\omega) = \frac{\nabla p}{\|\nabla p\|^2} g_1(r,\theta;\tilde{r},\omega)$, from Equation 4.12 and the divergence theorem, we get

(4.19)
$$I^{(2)}(\tilde{r},\omega) = \frac{-\tau}{2\pi l^{\epsilon}} \int_{\Gamma} (\mathbf{u}_{1}^{T}\mathbf{n}) \exp\left(\frac{ip(r,\theta;\tilde{r},\omega)}{\tau}\right) ds + \frac{\tau}{2\pi l^{\epsilon}} \iint_{\mathcal{D}^{\epsilon}} (\nabla \cdot \mathbf{u}_{1}(r,\theta;\tilde{r},\omega)) \exp\left(\frac{ip(r,\theta;\tilde{r},\omega)}{\tau}\right) dr d\theta.$$

As $I^{(2)}(\tilde{r},\omega) = O(\tau)$, it converges to zero as $\tau \to 0$. Applying the obtained results to Equation 4.13, we see that $I(\tilde{r},\omega)$ (and also $I_k(\tilde{r},\omega)$ defined in Equation 4.8) $\to 0$ as $\tau \to 0$ which completes the proof.

Since Theorem 4.2 is true for any $0 < \epsilon < \frac{1}{2}$, it also holds as $\epsilon \to 0$. As a corollary, we have the following result:

COROLLARY 4.3. If $\tilde{r} \neq 1$, then

(4.20)
$$\lim_{\epsilon \to 0} \lim_{\tau \to 0} F^{\epsilon}_{\tau}(\tilde{r}, \omega) = 0.$$

5. Spatial Frequencies as Gradient Histogram Bins. We now show that the squared magnitude of the Fourier transform of the CWR (ϕ) when polled close to the unit circle ($\tilde{r} = 1$) is approximately equal to the density function of the distance transform gradients (P) with the approximation becoming increasingly tight as $\tau \to 0$.

The squared magnitude of the Fourier transform—also called its *power spectrum* [1]—is given by

(5.1)
$$P_{\tau}^{\epsilon}(\tilde{r},\omega) \equiv |F_{\tau}^{\epsilon}(\tilde{r},\omega)|^2 = F_{\tau}^{\epsilon}(\tilde{r},\omega)\overline{F_{\tau}^{\epsilon}(\tilde{r},\omega)}.$$

By definition, $P^{\epsilon}_{\tau}(\tilde{r}, \omega) \geq 0$. From Lemma 4.1, we have

(5.2)
$$\int_{0}^{2\pi} \int_{0}^{\infty} P_{\tau}^{\epsilon}(\tilde{r},\omega) \tilde{r} d\tilde{r} d\omega = 1$$

independent of τ . Hence, $P_{\tau}^{\epsilon}(\tilde{r}, \omega)$ can be treated as a density function for all values of τ . We earlier observed that the gradient density function of the unit vector distance transform gradients is one-dimensional and defined over the space of orientations ω . For $P_{\tau}^{\epsilon}(\tilde{r}, \omega)$ to behave as an orientation density function, it needs to be integrated along the radial direction \tilde{r} . Since Theorem 4.2 states that the Fourier transform values are concentrated only on the unit circle $\tilde{r} = 1$ and converges to zero elsewhere as $\tau \to 0$, it should be sufficient if the integration for \tilde{r} is done over a region very close to $\tilde{r} = 1$. The following theorem—the principal result in this paper—confirms our observation.

THEOREM 5.1. For any given $0 < \epsilon < \frac{1}{2}$, $0 < \delta < 1$, $\omega_0 \in [0, 2\pi)$ and $0 < \Delta < 2\pi$,

(5.3)
$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \int_{1-\delta}^{1+\delta} P_{\tau}^{\epsilon}(\tilde{r}, \omega) \tilde{r} d\tilde{r} d\omega = \int_{\omega_0}^{\omega_0 + \Delta} P(\omega) d\omega.$$

5.1. An Intuitive Examination of Theorem 5.1. Before we proceed with the formal proof, we again try and give an intuitive explanation of why the theorem statement is true. The Fourier transform of the CWR defined in Equation 4.4 involves two spatial integrals (over x and y) which are converted into polar coordinate integrals. The squared magnitude of the Fourier transform (power spectrum), $P_{\tau}^{\epsilon}(\tilde{r},\omega)$, involves multiplying the Fourier transform with its complex conjugate. The complex conjugate is yet another 2D integral which we will perform in polar coordinates. As the gradient density function is one-dimensional and defined over the space of orientations, we integrate the power spectrum along the radial direction close to the unit circle $\tilde{r} = 1$ (as $\delta \to 0$). This is a fifth integral. When we poll the power spectrum $P_{\tau}^{\epsilon}(\tilde{r},\omega)$ close to $\tilde{r} = 1$, the two sinusoids, namely, $\exp\left(\frac{iS(x,y)}{\tau}\right)$ and $\exp\left(\frac{-i(ux+vy)}{\tau}\right)$ in Equation 4.4 are in resonance only when there is a *perfect match* between the orientation of each ray of the distance transform S(x, y) and the angle of the 2D spatial frequency ($\omega = \arctan\left(\frac{v}{u}\right)$). All the grid locations (x, y) having the same gradient orientation

(5.4)
$$\arctan\left(\frac{S_y}{S_x}\right) = \arctan\left(\frac{v}{u}\right)$$

cast a vote only at their corresponding spatial frequency "histogram" bin ω . Since the histogram bin is generally populated by votes from multiple grid locations, this leads to cross phase factors. Integrating the power spectrum over a small range on the orientation (constituting the sixth integral) helps in canceling out these phase factors giving us the desired result when we take the limit as $\tau \to 0$. This integral and limit cannot be exchanged because the phase factors will not otherwise cancel. The proof mainly deals with managing these six integrals.

5.2. Proof of Theorem 5.1. We now provide the formal proof of Theorem 5.1. For the sake of readability, we divide the proof into smaller subsections. To achieve a good flow, we state major portions of our proof as lemmas whose proofs are given in the appendix. We would like to emphasize that these lemmas are meaningful only within the context of the proof and do not have much significance as stand-alone statements. Important symbols used in the proof are adumbrated in Table 5.1.

Symbol	Comments
Ι	Integral of P_{τ}^{ϵ} over the radial length $[1 - \delta, 1 + \delta]$.
g_{jk}	Integral over the variables r , θ and \tilde{r} after symmetry breaking.
γ_{jk}	Phase term in the integral for g_{jk} .
$I_{jk}^{(1)}, I_{jk}^{(2)}$	Integrals for the main and the error terms of g_{jk} respectively.
	I is the sum of $I_{jk}^{(1)}$ and $I_{jk}^{(2)}$.
p,q	Functions used in the definition of $I_{ik}^{(1)}$.
	p represents the phase term.
$J_{jk}^{(1)}, J_{jk}^{(2)}, J_{jk}^{(3)}$	Integrals obtained when $I_{jk}^{(1)}$ is split over the integral range for θ' .
β	Symbol used to divide the integral range for θ' into three integrals.
	The limit as $\beta \to 0$ is considered in the proof.
G_{jk}	Result of integrating over θ' while evaluating $J_{jk}^{(1)}$.
ψ_{jk}, χ	Integrals for the main and the error terms of G_{jk} respectively.
	$J_{jk}^{(1)}$ is the sum of ψ_{jk} and χ .
ϵ_3,ϵ_4	Error terms used to define χ .

TABLE 5.1Table of important symbols

First, observe that

(5.5)
$$\overline{F_{\tau}^{\epsilon}(\tilde{r},\omega)} \equiv \sum_{k=1}^{K} \frac{\overline{C_k}}{2\pi\tau l^{\epsilon}} \int_0^{2\pi} \int_{R_k^{(1)}(\theta')}^{R_k^{(2)}(\theta')} \exp\left(-\frac{ir'}{\tau} \left[1 - \tilde{r}\cos(\theta' - \omega)\right]\right) r' dr' d\theta'$$

Define

(5.6)
$$I(\omega) \equiv \int_{1-\delta}^{1+\delta} P_{\tau}^{\epsilon}(\tilde{r},\omega)\tilde{r}d\tilde{r} = \int_{1-\delta}^{1+\delta} F_{\tau}^{\epsilon}(\tilde{r},\omega)\overline{F_{\tau}^{\epsilon}(\tilde{r},\omega)}\tilde{r}d\tilde{r}.$$

As $\tau \to 0$, we show that $I(\omega)$ approaches the density function of the gradients of S(x, y). Note that the integral in Equation 5.6 is over the interval $[1 - \delta, 1 + \delta]$, where $\delta > 0$ can be made arbitrarily small (as $\tau \to 0$) and this is due to Theorem 4.2.

Recall that in order to evaluate $I(\omega)$, we need to perform five integrals, four to obtain the power spectrum $P_{\tau}^{\epsilon}(\tilde{r},\omega)$ and a fifth along the radial direction \tilde{r} over $[1-\delta, 1+\delta]$ which is close to the unit circle $\tilde{r} = 1$. An easy way to compute $I(\omega)$ in the limit $\tau \to 0$ would be to apply a 5D stationary phase approximation [23]. Unfortunately, the 5D stationary phase approximation *cannot* be directly employed in our case for reasons detailed in Appendix B.

Breaking the symmetry of the integral. As described in Section B.1, we propose to solve for $I(\omega)$ in Equation 5.6 by breaking the symmetry of the integral. We fix the conjugate variables r' and θ' and perform the integration only with respect to the other three variables namely r, θ and \tilde{r} . To this end, let

(5.7)
$$I(\omega) = \sum_{j=1}^{K} \sum_{k=1}^{K} \frac{1}{(2\pi\tau l^{\epsilon})^2} \int_{0}^{2\pi} \int_{R_k^{(1)}(\theta')}^{R_k^{(2)}(\theta')} \exp\left(\frac{-ir'}{\tau}\right) g_{jk}(r',\theta';\omega)r'dr'd\theta',$$

where

$$(5.8) g_{jk}(r',\theta';\omega) = \int_{1-\delta}^{1+\delta} \int_0^{2\pi} \int_{R_j^{(1)}(\theta)}^{R_j^{(2)}(\theta)} \exp\left\{\frac{i}{\tau}\gamma_{jk}(r,\theta,\tilde{r};r',\theta',\omega)\right\} f_2(r,\tilde{r}) dr d\theta d\tilde{r}.$$

Here,

(5.9)
$$\gamma_{jk}(r,\theta,\tilde{r};r',\theta',\omega) = r \left[1 - \tilde{r}\cos(\theta - \omega)\right] + r'\tilde{r}\cos(\theta' - \omega) -\tilde{r}\left[\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)\right]$$

and

$$(5.10) f_2(r,\tilde{r}) = r\tilde{r}.$$

In the definition of $\gamma_{jk}(r, \theta, \tilde{r}; r', \theta', \omega)$ in Equation 5.9, ω, r' and θ' are held fixed. Similarly, in the definition of $g_{jk}(r', \theta'; \omega)$ in Equation 5.8, ω is a constant. The phase term of the quantity $C_j \overline{C_k}$ (Equation B.5) is absorbed in γ_{jk} and pursuant to Fubini's theorem [7], the integration with respect to \tilde{r} can be considered before the integration over r' and θ' . Define

(5.11)
$$r_{jk}(r',\theta';\omega) \equiv r'\cos(\theta'-\omega) - [\cos(\omega)(x_j-x_k) + \sin(\omega)(y_j-y_k)].$$

This leads to the following lemma.

LEMMA 5.2. If $r_{jk}(r', \theta'; \omega) > 0$, then as $\tau \to 0$,

(5.12)
$$g_{jk}(r',\theta';\omega) = (2\pi\tau)^{\frac{3}{2}} \sqrt{r_{jk}(r',\theta';\omega)} \exp\left(\frac{ir_{jk}(r',\theta';\omega)}{\tau} + \frac{i\pi}{4}\right) + \tau^{\kappa}\xi_{jk}(r',\theta';\omega)$$

where $\kappa \geq 2$ and $\xi_{jk}(r', \theta'; \omega)$ is some bounded continuous function which includes the contributions from the boundary. If $r_{jk}(r', \theta'; \omega) \leq 0$, then as $\tau \to 0$, $g_{jk}(r', \theta'; \omega) = 0$.

The proof of Lemma 5.2—obtained using a three dimensional stationary phase approximation—is available in Appendix C. Note that for j = k and θ' close to ω , $r_{jk}(r', \theta'; \omega) > 0$ and hence $g_{jk}(r', \theta'; \omega) \neq 0$. Below, we show that the only pertinent scenarios that need consideration are θ' close to ω and j = k. When θ' is away from ω or $j \neq k$, the integral $g_{jk}(r', \theta'; \omega)$ vanishes. Hence, for the sake of readability of our proof, we let $g_{jk}(r', \theta'; \omega)$ take the most general form given in Equation 5.12 for all values of r' and θ' .

Determining $I(\omega)$. Substituting the value of $g_{jk}(r', \theta'; \omega)$ into Equation 5.7, as $\tau \to 0$, we get

(5.13)
$$I(\omega) = \sum_{j=1}^{K} \sum_{k=1}^{K} \left\{ \frac{\eta_{jk}(\omega)}{L^{\epsilon}} I_{jk}^{(1)}(\omega) + I_{jk}^{(2)}(\omega) \right\}$$

where

$$\begin{split} I_{jk}^{(1)}(\omega) &= \frac{1}{\sqrt{2\pi\tau}} \int_{0}^{2\pi} \int_{R_{k}^{(1)}(\theta')}^{R_{k}^{(2)}(\theta')} \exp\left(\frac{ip(r',\theta';\omega)}{\tau}\right) q(r',\theta';\omega) dr' d\theta',\\ I_{jk}^{(2)}(\omega) &= \int_{0}^{2\pi} \int_{R_{k}^{(1)}(\theta')}^{R_{k}^{(2)}(\theta')} \exp\left(\frac{-ir'}{\tau}\right) r' \frac{1}{(2\pi l^{\epsilon})^{2}} \tau^{\kappa-2} \xi_{jk}(r',\theta';\omega) dr' d\theta' \end{split}$$

$$\eta_{jk}(\omega) = \exp\left(\frac{-i\alpha_{jk}(\omega)}{\tau} + \frac{i\pi}{4}\right),$$

$$\alpha_{jk}(\omega) = \cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k),$$

$$p(r', \theta'; \omega) = -r'[1 - \cos(\theta' - \omega)] \text{ and}$$

$$(5.14) q(r', \theta'; \omega) = r'\sqrt{r'\cos(\theta' - \omega) - \alpha_{jk}(\omega)}.$$

In the definition of the functions $p(r', \theta'; \omega)$ and $q(r', \theta'; \omega)$, ω is held fixed. Since $\kappa \geq 2$, by the Riemann-Lebesgue lemma, we have $\lim_{\tau \to 0} I_{jk}^{(2)} = 0$ and from the Lebesgue dominated convergence theorem, it follows that

(5.15)
$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} I_{jk}^{(2)}(\omega) = \int_0^{2\pi} \lim_{\tau \to 0} I_{jk}^{(2)}(\omega) = 0.$$

Using the above result in Equation 5.13, we get

(5.16)
$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} I(\omega) d\omega = \sum_{j=1}^K \sum_{k=1}^K \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} I_{jk}^{(1)}(\omega) d\omega.$$

Splitting the integral over θ' into three disconnected regions. Consider the integral $I_{jk}^{(1)}(\omega)$. As essential contributions to it come only from the stationary points of $p(r', \theta'; \omega)$ [12, 24, 3] (with ω held fixed), we first determine its critical (stationary) point(s). The partial derivatives of $p(r', \theta'; \omega)$ at a fixed value of ω are given by

(5.17)
$$\frac{\partial p}{\partial r'} = -1 + \cos(\theta' - \omega), \quad \frac{\partial p}{\partial \theta'} = -r' \sin(\theta' - \omega).$$

For $\nabla p = 0$, we must have $\theta' = \omega$. Hence, in order to evaluate $I_{jk}^{(1)}(\omega)$, we find it useful to divide the integral range $[0, 2\pi)$ for θ' into three disjoint regions namely $[0, \omega - \beta)$, $[\omega - \beta, \omega + \beta]$ and $(\omega + \beta, 2\pi)$ for a fixed $\beta > 0$, and write

(5.18)
$$I_{jk}^{(1)}(\omega) = J_{jk}^{(1)}(\beta,\omega) + J_{jk}^{(2)}(\beta,\omega) + J_{jk}^{(3)}(\beta,\omega)$$

where

$$J_{jk}^{(1)}(\beta,\omega) = \frac{1}{\sqrt{2\pi\tau}} \int_{\omega-\beta}^{\omega+\beta} \int_{R_k^{(1)}(\theta')}^{R_k^{(2)}(\theta')} \exp\left(\frac{ip(r',\theta';\omega)}{\tau}\right) q(r',\theta';\omega) dr' d\theta',$$

$$J_{jk}^{(2)}(\beta,\omega) = \frac{1}{\sqrt{2\pi\tau}} \int_0^{\omega-\beta} \int_{R_k^{(1)}(\theta')}^{R_k^{(2)}(\theta')} \exp\left(\frac{ip(r',\theta';\omega)}{\tau}\right) q(r',\theta';\omega) dr' d\theta', \text{ and}$$

$$(5.19) J_{jk}^{(3)}(\beta,\omega) = \frac{1}{\sqrt{2\pi\tau}} \int_{\omega+\beta}^{2\pi} \int_{R_k^{(1)}(\theta')}^{R_k^{(2)}(\theta')} \exp\left(\frac{ip(r',\theta';\omega)}{\tau}\right) q(r',\theta';\omega) dr' d\theta'.$$

Since the above relation is true for any $\beta > 0$, we can let $\beta \to 0$ (after we take the limit $\tau \to 0$). Fix a β close to zero and consider the above integrals as $\tau \to 0$. Then we obtain:

Lemma 5.3.

(5.20)
$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} I_{jk}^{(1)}(\omega) d\omega = \lim_{\beta \to 0} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} J_{jk}^{(1)}(\beta, \omega) d\omega.$$

The proof is available in Appendix D.

Interchanging the order of integration between r' and θ' . We now evaluate $J_{jk}^{(1)}(\beta,\omega)$ by interchanging the order of integration between r' and θ' which requires us to rewrite θ' as a function of r'. Recall that the boundaries of each \mathcal{D}_k^{ϵ} along $r(\theta') = R_k^{(1)}(\theta')$ and $r(\theta') = R_k^{(2)}(\theta')$ respectively are composed of a finite sequence of straight line segments. In order to evaluate $J_{jk}^{(1)}(\beta,\omega)$, we need to consider these boundaries only within the precincts of the angles $[\omega - \beta, \omega + \beta]$ at each \mathcal{D}_k^{ϵ} . But for sufficiently small β , we observe that for every $\omega \in [0, 2\pi)$, when we consider these boundaries (along $R_k^{(1)}(\theta')$ and $R_k^{(2)}(\theta')$ respectively) within the angles $[\omega - \beta, \omega + \beta]$, they are composed of at most two line segments as portrayed in Figure 5.1.



FIGURE 5.1. Boundary considered within the angles $[\omega - \beta, \omega + \beta]$ is comprised of at most two line segments L_1 and L_2 .

Over each line segment, $r'(\theta')$ is either strictly monotonic (strictly increases or strictly decreases) or has exactly one critical point (strictly decreases, attains a minimum and then strictly increases) as described in Figure 5.2.



FIGURE 5.2. Plot of radial length (r) vs angle (θ) .

Hence, it follows that for sufficiently small β , θ' rewritten as a function of r' is composed of at most three disconnected regions (as seen in Figure 5.3).

Let $\mathcal{B}(r') \subseteq [\omega - \beta, \omega + \beta]$ denote the integration region for $\theta'(r')$. Treating θ' as



FIGURE 5.3. Three disconnected regions for the angle (θ) .

a function of r', the integral $J_{jk}^{(1)}(\beta,\omega)$ can be rewritten as

(5.21)
$$J_{jk}^{(1)}(\beta,\omega) = \int_{r_k^{(1)}(\beta,\omega)}^{r_k^{(2)}(\beta,\omega)} G_{jk}(r',\omega) dr',$$

where

(5.22)
$$r_{k}^{(1)}(\beta,\omega) = \inf\{R_{k}^{(1)}(\theta')\}, \text{ and }$$
$$r_{k}^{(2)}(\beta,\omega) = \sup\{R_{k}^{(2)}(\theta')\}$$

(1)

with $\theta' \in [\omega - \beta, \omega + \beta]$ and

(5.23)
$$G_{jk}(r',\omega) = \frac{1}{\sqrt{2\pi\tau}} \int_{\mathcal{B}(r')} \exp\left(\frac{ip(r',\theta',\omega)}{\tau}\right) q(r',\theta',\omega) d\theta'.$$

Note that while evaluating the integral $G_{jk}(r',\omega)$, r' and ω are held fixed. As contributions to $G_{jk}(r',\omega)$ come only from the stationary points of $p(r',\theta',\omega)$ (with r' and ω held fixed) as $\tau \to 0$, we evaluate $\frac{\partial p}{\partial \theta'} = -r' \sin(\theta' - \omega)$ and for it to vanish, we require $\theta' = \omega$. Moreover

(5.24)
$$\begin{aligned} \frac{\partial^2 p}{\partial \theta'^2} \Big|_{\omega} &= -r', \\ p(r', \omega, \omega) &= 0, \text{ and} \\ q(r', \omega, \omega) &= r' \sqrt{r' - \alpha_{jk}(\omega)}. \end{aligned}$$

For the given r', if $\omega \notin \mathcal{B}(r')$, no stationary points exist. Using integration by parts, $G_{jk}(r',\omega)$ can be shown to be $\epsilon_3(r',\omega,\tau) = O(\sqrt{\tau})$, which can be uniformly bounded by a function of r' and ω for small values of τ .

If $\omega \in \mathcal{B}(r')$, then using the one dimensional stationary phase approximation [15, 16], it can be shown that

(5.25)
$$G_{jk}(r',\omega) = \exp\left(\frac{-i\pi}{4}\right)\sqrt{r'}\sqrt{r'-\alpha_{jk}(\omega)} + \epsilon_4(r',\omega,\tau),$$

where $\epsilon_4(r', \omega, \tau)$ can be uniformly bounded by a function of r' and ω for small values of τ and converges to zero as $\tau \to 0$. Here, we assume that the stationary point $\theta' = \omega$ lies in the *interior* of $\mathcal{B}(r')$ and not on the boundary as there can be at most finite (actually 2) values of r' (with Lebesgue measure zero) for which $\theta' = \omega$ can lie on the boundary of $\mathcal{B}(r')$. Computing the integral over ω and r'. Let $r_k^{(-)}(\beta,\omega) \geq r_k^{(1)}(\beta,\omega)$ and $r_k^{(+)}(\beta,\omega) \leq r_k^{(2)}(\beta,\omega)$ be the values of r' such that when $r_k^{(-)}(\beta,\omega) < r' < r_k^{(+)}(\beta,\omega)$, the stationary point $\theta' = \omega$ lies in the interior of $\mathcal{B}(r')$. Substituting the value of $G_{jk}(r',\omega)$ into Equation 5.21 and using the definitions of $\eta_{jk}(\omega)$ and $\alpha_{jk}(\omega)$ from Equation 5.14, we get

$$\int_{\omega_0}^{\omega_0+\Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} J_{jk}^{(1)}(\beta,\omega) d\omega = \psi_{jk}(\beta) + \int_{\omega_0}^{\omega_0+\Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} \left\{ \int_{r_k^{(1)}(\beta,\omega)}^{r_k^{(2)}(\beta,\omega)} \chi(r',\omega,\tau) \, dr' \right\} d\omega,$$
(5.26)

where

(5.27)
$$\psi_{jk}(\beta) = \frac{1}{L^{\epsilon}} \int_{\omega_0}^{\omega_0 + \Delta} \exp\left(\frac{-i\alpha_{jk}(\omega)}{\tau}\right) \int_{r_k^{(-)}(\beta,\omega)}^{r_k^{(+)}(\beta,\omega)} \sqrt{r'} \sqrt{r' - \alpha_{jk}(\omega)} dr' d\omega$$

and

$$\chi(r',\omega,\tau) = \begin{cases} \epsilon_4(r',\omega,\tau), & r_k^{(-)}(\beta,\omega) < r' < r_k^{(+)}(\beta,\omega), \\ \epsilon_3(r',\omega,\tau), & r' < r_k^{(-)}(\beta,\omega) \text{ or } r_k^{(+)}(\beta,\omega) < r'. \end{cases}$$

Since $|\eta_{jk}(\omega)| = 1$ and $\chi(r', \omega, \tau)$ can be uniformly bounded by a function r' and ω for small values of τ , by the Lebesgue dominated convergence theorem we have

(5.28)
$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} \left\{ \int_{r_k^{(1)}(\beta,\omega)}^{r_k^{(2)}(\beta,\omega)} \chi(r',\omega,\tau) \, dr' \right\} d\omega$$
$$= \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} \left\{ \int_{r_k^{(1)}(\beta,\omega)}^{r_k^{(2)}(\beta,\omega)} \lim_{\tau \to 0} \chi(r',\omega,\tau) \, dr' \right\} d\omega = 0.$$

This leaves us having to prove the following result:

Lemma 5.4.

(5.29)
$$\sum_{j=1}^{K} \sum_{k=1}^{K} \lim_{\beta \to 0} \lim_{\tau \to 0} \psi_{jk}(\beta) = \int_{\omega_0}^{\omega_0 + \Delta} P(\omega) d\omega.$$

The proof of this lemma is given in Appendix E. This completes the proof of Theorem 5.1.

We would like to give a short recap of our proof. Beginning with the definition of $I(\omega)$ in Equation 5.6, Lemma 5.2 and the statements following it lead to the relation 5.16, namely,

(5.30)
$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} I(\omega) d\omega = \sum_{j=1}^K \sum_{k=1}^K \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} I_{jk}^{(1)}(\omega) d\omega.$$

From Lemma 5.3, it follows that

(5.31)
$$\sum_{j=1}^{K} \sum_{k=1}^{K} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} I_{jk}^{(1)}(\omega) d\omega = \sum_{j=1}^{K} \sum_{k=1}^{K} \lim_{\beta \to 0} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} J_{jk}^{(1)}(\beta, \omega) d\omega.$$

Interchanging the order of integration between r' and $\theta',$ we showed that

(5.32)
$$\sum_{j=1}^{K} \sum_{k=1}^{K} \lim_{\beta \to 0} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}(\omega)}{L^{\epsilon}} J_{jk}^{(1)}(\beta, \omega) d\omega = \sum_{j=1}^{K} \sum_{k=1}^{K} \lim_{\beta \to 0} \lim_{\tau \to 0} \psi_{jk}(\beta)$$

Finally, the application of Lemma 5.4 gives the desired result of Theorem 5.1, namely,

(5.33)

$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \int_{1-\delta}^{1+\delta} P_{\tau}^{\epsilon}(\tilde{r}, \omega) \tilde{r} d\tilde{r} d\omega = \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} I(\omega) d\omega$$

$$= \sum_{j=1}^{K} \sum_{k=1}^{K} \lim_{\beta \to 0} \lim_{\tau \to 0} \psi_{jk}(\beta)$$

$$= \int_{\omega_0}^{\omega_0 + \Delta} P(\omega) d\omega.$$

6. Results Stemming from the Main Theorem. As an implication of Theorem 5.1, we have the following corollary.

COROLLARY 6.1. For any given $0 < \delta < 1, \omega_0 \in [0, 2\pi)$,

(6.1)
$$\lim_{\epsilon \to 0} \lim_{\Delta \to 0} \frac{1}{\Delta} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \left\{ \int_{1-\delta}^{1+\delta} P_{\tau}^{\epsilon}(\tilde{r}, \omega) \tilde{r} \, d\tilde{r} \right\} d\omega = P(\omega_0).$$

Proof. From Equation 3.5, we have

(6.2)
$$\lim_{\Delta \to 0} \frac{1}{\Delta} \int_{\omega_0}^{\omega_0 + \Delta} P(\omega) d\omega = \lim_{\Delta \to 0} \frac{F(\omega_0 \le \omega \le \omega_0 + \Delta)}{\Delta} = P(\omega_0).$$

Since Theorem 5.1 is true for any $0 < \epsilon < \frac{1}{2}$, it also holds as $\epsilon \to 0$. The result then follows immediately.

Theorem 5.1 also entails the following lemma. LEMMA 6.2. For any given $0 < \epsilon < \frac{1}{2}$, $0 < \delta < 1$,

(6.3)
$$\lim_{\tau \to 0} \int_0^{2\pi} \int_{1-\delta}^{1+\delta} P_{\tau}^{\epsilon}(\tilde{r},\omega) \tilde{r} d\tilde{r} d\omega = 1.$$

Proof. Since the result shown in Theorem 5.1 holds good for any ω_0 and Δ , we may choose $\omega_0 = 0$ and $\Delta = 2\pi$. Using Equation 3.6 the result follows immediately as

(6.4)
$$\lim_{\tau \to 0} \int_0^{2\pi} \int_{1-\delta}^{1+\delta} P_{\tau}^{\epsilon}(\tilde{r},\omega) \tilde{r} d\tilde{r} d\omega = \int_0^{2\pi} P(\omega) d\omega = 1.$$

Lemmas 6.2 and 4.1 leads to the following corollaries.

COROLLARY 6.3. For any given $0 < \epsilon < \frac{1}{2}, 0 < \delta < 1$,

(6.5)
$$\lim_{\tau \to 0} \int_0^{2\pi} \left\{ \int_0^{1-\delta} P^{\epsilon}_{\tau}(\tilde{r},\omega) \ \tilde{r}d\tilde{r} + \int_{1+\delta}^{\infty} P^{\epsilon}_{\tau}(\tilde{r},\omega) \ \tilde{r}d\tilde{r} \right\} d\omega = 0.$$

16

Proof. From Lemma 4.1, we have for any $\tau > 0$ and $0 < \epsilon < \frac{1}{2}$,

(6.6)
$$\int_0^{2\pi} \int_0^\infty P_\tau^\epsilon(\tilde{r},\omega) \tilde{r} d\tilde{r} d\omega = 1.$$

For the given $0 < \delta < 1$, dividing the integral range $(0, \infty)$ for \tilde{r} into three disjoint regions namely $(0, 1 - \delta)$, $[1 - \delta, 1 + \delta]$ and $(1 + \delta, \infty)$ and letting $\tau \to 0$, we have

$$\lim_{\tau \to 0} \int_0^{2\pi} \left\{ \int_0^{1-\delta} P_\tau^\epsilon(\tilde{r},\omega) \ \tilde{r} d\tilde{r} + \int_{1-\delta}^{1+\delta} P_\tau^\epsilon(\tilde{r},\omega) \ \tilde{r} d\tilde{r} + \int_{1+\delta}^\infty P_\tau^\epsilon(\tilde{r},\omega) \ \tilde{r} d\tilde{r} \right\} d\omega = 1.$$

Pursuant to Lemma 6.2, the limit

(6.7)
$$\lim_{\tau \to 0} \int_0^{2\pi} \int_{1-\delta}^{1+\delta} P^{\epsilon}_{\tau}(\tilde{r},\omega) \ \tilde{r} d\tilde{r} d\omega$$

exists and equals 1. The result then follows.

COROLLARY 6.4. For any given $0 < \epsilon < \frac{1}{2}$, $0 < \delta < 1$, $\omega_0 \in [0, 2\pi)$ and $0 < \Delta < 2\pi$,

(6.8)
$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \left\{ \int_0^{1-\delta} P_{\tau}^{\epsilon}(\tilde{r}, \omega) \ \tilde{r} d\tilde{r} + \int_{1+\delta}^{\infty} P_{\tau}^{\epsilon}(\tilde{r}, \omega) \ \tilde{r} d\tilde{r} \right\} d\omega = 0.$$

Proof. Let $M = \lfloor \frac{2\pi}{\Delta} \rfloor$. Define $\omega_{i+1} \equiv \omega_i + \Delta \mod 2\pi$ for $0 \leq i \leq M - 1$. Then, we have from Corollary 6.3,

(6.9)
$$\lim_{\tau \to 0} \left[\sum_{i=0}^{M-1} \int_{\omega_i}^{\omega_{i+1}} \mathcal{Q}(\omega) d\omega + \int_{\omega_{i+1}}^{\omega_0+2\pi} \mathcal{Q}(\omega) d\omega \right] = 0,$$

where

(6.10)
$$\mathcal{Q}(\omega) = \int_0^{1-\delta} P_{\tau}^{\epsilon}(\tilde{r},\omega) \ \tilde{r}d\tilde{r} + \int_{1+\delta}^{\infty} P_{\tau}^{\epsilon}(\tilde{r},\omega) \ \tilde{r}d\tilde{r}.$$

Since $P^{\epsilon}_{\tau}(\tilde{r},\omega)\tilde{r} \geq 0$, it follows that $\mathcal{Q}(\omega)$ and both integrals in Equation 6.9 are *non-negative* and hence each integral converges to zero *independently* giving us the desired result.

From Theorem 5.1 and Corollaries 6.1 and 6.4, the subsequent results follow almost immediately.

PROPOSITION 6.5. For any given $0 < \epsilon < \frac{1}{2}$, $\omega_0 \in [0, 2\pi)$ and $0 < \Delta < 2\pi$,

(6.11)
$$\lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \left\{ \int_0^\infty P_\tau^\epsilon(\tilde{r}, \omega) \tilde{r} \, d\tilde{r} \right\} d\omega = \int_{\omega_0}^{\omega_0 + \Delta} P(\omega) d\omega.$$

COROLLARY 6.6. For any given $\omega_0 \in [0, 2\pi)$,

(6.12)
$$\lim_{\epsilon \to 0} \lim_{\Delta \to 0} \frac{1}{\Delta} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \left\{ \int_0^\infty P_\tau^\epsilon(\tilde{r}, \omega) \tilde{r} \, d\tilde{r} \right\} d\omega = P(\omega_0).$$

7. Significance of our result and concluding remarks. The integrals

(7.1)
$$\int_{\omega_0}^{\omega_0+\Delta} \int_{1-\delta}^{1+\delta} P^{\epsilon}_{\tau}(\tilde{r},\omega)\tilde{r}d\tilde{r}d\omega, \quad \int_{\omega_0}^{\omega_0+\Delta} P(\omega)d\omega$$

give the interval measures of the density functions P_{τ}^{ϵ} (when polled close to the unit circle $\tilde{r} = 1$) and P respectively. Theorem 5.1 states that at small values of τ , both the interval measures are approximately equal, with the difference between them being o(1). Furthermore the result is also true as $\epsilon \to 0$. Recall that by definition, P_{τ}^{ϵ} is the normalized power spectrum of the wave function $\phi(x, y) = \exp\left(\frac{iS(x,y)}{\tau}\right)$. Hence, we conclude that the power spectrum of $\phi(x, y)$ when polled close to the unit circle $\tilde{r} = 1$ (as $\delta \to 0$ in Theorem 5.1), or when integrated over \tilde{r} (with reference to Proposition 6.5), can potentially serve as a *density estimator* of the orientation of ∇S for small values of τ and ϵ . Our work is essentially an application of the higher-order stationary phase approximation culminating in a new density estimator.

7.1. Advantages of our formulation. One of the foremost advantages of our method is that the orientation gradient density is computed without actually determining the distance transform gradients. Since the stationary points (as seen from the stationary phase approximation) capture gradient information and slot them into the corresponding frequency bins, we can directly work with the distance function—circumventing the need to compute its derivatives. We are *not aware of any previous work* that estimates the orientation gradient density without first computing the gradients of the distance transform.

Recall that we furnished a closed-form expression for the distance transform gradient density function $P(\theta)$ in Equation 3.5. While it initially appears attractive, computing the density function via the closed-form expression is practically cumbersome as we need to first determine the Voronoi region corresponding to each Voronoi center (source point) Y_k and then for each orientation direction θ , compute the ray length $R_k(\theta)$ from Y_k to its Voronoi boundary along θ . These involve unwieldy manipulations of complex data structures. On the other hand, our mathematical result provides an easy mechanism to achieve the same task as it is computationally faster and easier to implement. Given the N sampled values \hat{S} of the distance function S from a point-set of cardinality K, we just need to compute the fast Fourier transform of $\exp\left(\frac{i\hat{S}(x)}{\tau}\right)$ —an $O(N \log N)$ operation—and then subsequently compute the squared magnitude (to obtain the power spectrum)—performed in O(N). Hence the orientation density function can be determined in $O(N \log N)$ independent of the cardinality of the point-set (K). Our algorithm is computationally efficient even when K = O(N).

7.2. Possible Extensions. The present work only deals with special kinds of distance functions, namely those defined from a set of discrete point locations in two dimensions. Other cases include signed distance functions, distance functions defined from a set of curves [18, 20] etc. While our work initially appears to be somewhat restrictive, we should note that as the cardinality and locations in each point-set can be arbitrary, the resulting distance functions can be quite complex. We strongly believe that it is possible to establish a similar Fourier transform-based density estimation result even for *arbitrary*, continuous (and differentiable) functions in 2D. A general result of this nature would subsume our current work on point-set distance functions as well as the other kinds of distance functions mentioned above, as the gradient

magnitude of an arbitrary 2D function can vary across the point locations and need not necessarily be identically equal to one as in the case of distance functions. Here the gradient density function will inherently be two dimensional and defined both along the radial, gradient magnitude direction $\left(r = \sqrt{S_x^2 + S_y^2}\right)$ and the orientation $\left(\theta = \arctan\left(\frac{S_y}{S_x}\right)\right)$. However, at the present time, due to many technical issues, this is merely a conjecture and requires further investigation. Generalization of the present work to three dimensions is also a concrete possibility. These are fruitful avenues for future research and we may explore them in the years to come.

Appendix A. Proof of Lemma 4.1.

Proof. Define a function H(x, y) by

$$H(x,y) \equiv \begin{cases} 1 : & \text{if } (x,y) \in \Omega^{\epsilon}; \\ 0 : & \text{otherwise.} \end{cases}$$

Let $f(x,y) = H(x,y) \exp\left(\frac{iS(x,y)}{\tau}\right)$. Then,

(A.1)
$$F_{\tau}^{\epsilon}(u,v) = \frac{1}{2\pi\tau l^{\epsilon}} \iint f(x,y) \exp\left(\frac{-i(ux+vy)}{\tau}\right) dxdy.$$

Let $\frac{u}{\tau} = s$, $\frac{v}{\tau} = t$ and $G(s,t) = F_{\tau}^{\epsilon}(s\tau,t\tau)$. Then,

(A.2)
$$\tau l^{\epsilon}G(s,t) = \frac{1}{2\pi} \iint f(x,y) \exp\left(-i(sx+ty)\right) dxdy.$$

Since f is ℓ^1 integrable, by Parseval's theorem [1], we have

(A.3)
$$\iint |f(x,y)|^2 dxdy = \iint |\tau l^{\epsilon}G(s,t)|^2 dsdt = (\tau l^{\epsilon})^2 \iint |F_{\tau}^{\epsilon}(s\tau,t\tau)|^2 dsdt.$$

Letting $u = s\tau$, $v = t\tau$ and observing that

(A.4)
$$\iint |f(x,y)|^2 \, dx \, dy = \iint_{\Omega^{\epsilon}} \left| \exp\left(\frac{iS(x,y)}{\tau}\right) \right|^2 \, dx \, dy = L^{\epsilon}$$

we get

(A.5)
$$(l^{\epsilon})^2 \iint |F^{\epsilon}_{\tau}(u,v)|^2 \, du dv = L^{\epsilon}.$$

Hence

(A.6)
$$\iint |F^{\epsilon}_{\tau}(u,v)|^2 \, du dv = 1$$

which completes the proof. \square

Appendix B. Difficulty with the 5D stationary phase approximation. Since $P^{\epsilon}_{\tau}(\tilde{r}, \omega)$ equals $F^{\epsilon}_{\tau}(\tilde{r}, \omega) \overline{F^{\epsilon}_{\tau}(\tilde{r}, \omega)}$, we have

(B.1)
$$I(\omega) = \sum_{j=1}^{K} \sum_{k=1}^{K} \frac{1}{(2\pi\tau l^{\epsilon})^2} N_{jk}(\omega),$$

where

(B.2)
$$N_{jk}(\omega) = \int_{1-\delta}^{1+\delta} \int_0^{2\pi} \int_{R_k^{(1)}(\theta')}^{R_k^{(2)}(\theta')} \int_0^{2\pi} \int_{R_k^{(1)}(\theta)}^{R_k^{(2)}(\theta)} \exp\left(\frac{i}{\tau}b_{jk}\right) f_1 dr d\theta dr' d\theta' d\tilde{r}.$$

Here,

(B.3)
$$b_{jk}(r,\theta,r',\theta',\tilde{r};\omega) = r \left[1 - \tilde{r}\cos(\theta - \omega)\right] - r' \left[1 - \tilde{r}\cos(\theta' - \omega)\right] - \tilde{r} \left[\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)\right]$$

and

(B.4)
$$f_1(r, r', \tilde{r}) = rr'\tilde{r}.$$

Notice that the phase term of the quantity $C_j \overline{C_k}$, namely

(B.5)
$$-\tilde{r}\left[\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)\right]$$

is absorbed in b_{jk} . Since we are interested only in the limit as $\tau \to 0$, essential contribution to $N_{jk}(\omega)$ comes only from the stationary (critical) point(s) of b_{jk} [23]. The partial derivatives of $b_{jk}(r, \theta, r', \theta', \tilde{r})$ are given by

$$\frac{\partial b_{jk}}{\partial r} = 1 - \tilde{r}\cos(\theta - \omega), \quad \frac{\partial b_{jk}}{\partial \theta} = r\tilde{r}\sin(\theta - \omega),$$
$$\frac{\partial b_{jk}}{\partial r'} = -1 + \tilde{r}\cos(\theta' - \omega), \quad \frac{\partial b_{jk}}{\partial \theta'} = -r'\tilde{r}\sin(\theta' - \omega), \text{ and}$$
$$(B.6)\frac{\partial b_{jk}}{\partial \tilde{r}} = -r\cos(\theta - \omega) + r'\cos(\theta' - \omega) - [\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)].$$

As r, r' and $\tilde{r} > 0$, it is easy to see that for $\nabla b_{jk} = 0$ (stationary), we must have

(B.7)
$$\tilde{r} = 1, \quad \theta = \theta' = \omega, r = r' - [\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)].$$

Let t_0 denote the stationary point. The Hessian matrix \mathcal{W} of b_{jk} at t_0 is given by

$$\mathcal{W}(r,\theta,r',\theta',\tilde{r})|_{t_0} = \begin{bmatrix} 0 & 0 & 0 & 0 & -1 \\ 0 & r_{t_0} & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & -r' & 0 \\ -1 & 0 & 1 & 0 & 0 \end{bmatrix}$$

where $r_{t_0} = r' - [\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)]$. Unfortunately, the determinant of \mathcal{W} at the stationary point t_0 equals 0 as the first and third rows—corresponding to r and r' respectively—are scalar multiples of each other. This impedes us from directly applying the 5D stationary phase approximation [23].

The addition of a 6th integral to the above setup—where the power spectrum is integrated over a small range on the orientation ω (in order to remove cross phase factors)—leads to a 6D stationary phase approximation. This is of no help either, as the Hessian continues to remain degenerate for the same reasons as above.

20

B.1. Avoiding degeneracy by symmetry breaking. As we notice above, the degeneracy in the 5D (and 6D) stationary phase approximation arises because the determinant of the Hessian, namely \mathcal{W} , when evaluated at the stationary point t_0 takes the value zero, as its first and third rows corresponding to r and r' respectively are scalar multiples of each other. Also, observe that the value of either r or r' is not determined at the stationary point and can take on arbitrary values. However, the rows (and columns) of \mathcal{W} corresponding to the other three variables θ, θ' and \tilde{r} are indeed independent of each other and do not cause degeneracy. This strongly suggests that if we *do not* consider both r and r' together and hold back either one of them, say r', the resulting 4D stationary phase approximation will be well-defined. Since the integration range for r' is defined in terms of θ' , we retain both these variables and perform the stationary phase approximation on the other three variables. This manual breaking of symmetry avoids the degeneracy issue.

Appendix C. Proof of Lemma 5.2.

Proof. Recall that the essential contribution to $g_{jk}(r', \theta'; \omega)$ comes only from the stationary points of γ_{jk} as $\tau \to 0$ [23]. The partial derivatives of $\gamma_{jk}(r, \theta, \tilde{r}; r', \theta', \omega)$ are given by

$$\frac{\partial \gamma_{jk}}{\partial r} = 1 - \tilde{r}\cos(\theta - \omega), \quad \frac{\partial \gamma_{jk}}{\partial \theta} = r\tilde{r}\sin(\theta - \omega), \text{ and}$$

$$(C.1)\frac{\partial \gamma_{jk}}{\partial \tilde{r}} = -r\cos(\theta - \omega) + r'\cos(\theta' - \omega) - [\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)].$$

As both r and $\tilde{r} > 0$, for $\nabla \gamma_{jk} = 0$ (stationary), we must have

(C.2)
$$\tilde{r} = 1, \ \theta = \omega, \text{ and}$$

 $r = r' \cos(\theta' - \omega) - [\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)].$

Let t_0 denote a stationary point. Then

$$\gamma_{jk}(t_0) = r' \cos(\theta' - \omega) - [\cos(\omega)(x_j - x_k) + \sin(\omega)(y_j - y_k)] = r_{jk}(r', \theta'; \omega),$$

$$f_2(t_0) = r_{jk}(r', \theta'; \omega)$$

and the Hessian matrix \mathcal{H} of γ_{jk} at the stationary point t_0 is

$$\mathcal{H}(r,\theta,\tilde{r})|_{t_0} = \begin{bmatrix} 0 & 0 & -1 \\ 0 & r_{jk}(r',\theta';\omega) & 0 \\ -1 & 0 & 0 \end{bmatrix}.$$

It can be easily verified that the determinant of \mathcal{H} equals $-r_{jk}(r', \theta')$.

If $r_{jk}(r', \theta'; \omega) \leq 0$, no stationary points exist as r > 0 by definition and hence $g_{jk}(r', \theta'; \omega) = 0$ as $\tau \to 0$ [23]. If $r_{jk}(r', \theta'; \omega) > 0$, the determinant of \mathcal{H} is strictly negative and its signature–the difference between the number of positive and negative eigenvalues–is 1. Then, from the higher-order stationary phase approximation [23], we have

$$g_{jk}(r',\theta';\omega) = (2\pi\tau)^{\frac{3}{2}} \sqrt{r_{jk}(r',\theta';\omega)} \exp\left(\frac{ir_{jk}(r',\theta';\omega)}{\tau} + \frac{i\pi}{4}\right) + \epsilon_1(r',\theta',\tau;\omega)$$

as $\tau \to 0$, where $\epsilon_1(r', \theta', \tau; \omega)$ includes the contributions from the boundary in Equation 5.8. Here we have assumed that the stationary point t_0 does not occur on the

boundary and lies to its interior, i.e, $R_j^{(1)}(\theta) < r_{jk}(r', \theta'; \omega) < R_j^{(2)}(\theta)$, as the measure on the set of $\{\omega, \theta', r'\}$ for which $r_{jk}(r', \theta'; \omega)$ (or t_0) can occur on the boundary is zero.

Let Γ denote the boundary in Equation 5.8. If there does not exist a 2D patch on Γ on which γ_{jk} is constant, then we can conclude that $\epsilon_1(r', \theta', \tau; \omega)$ —which includes the contributions from the boundary Γ involving the stationary points of the second kind where the level curves of γ_{jk} are tangential to Γ —should be at least $O(\tau^2)$ as $\tau \to 0$ [12, 3, 23]. From this, we get

(C.3)
$$\epsilon_1(r', \theta', \tau; \omega) = \tau^{\kappa} \xi_{jk}(r', \theta'; \omega)$$

where $\kappa \geq 2$ and $\xi_{jk}(r', \theta'; \omega)$ is some bounded, continuous function. Since the boundary Γ is made of straight line segments, we can show that this is indeed the case. Below, we take care of this technical issue.

The boundary Γ in Equation 5.8 is the *union* of two disconnected surfaces $\Gamma_1 = \mathcal{A}_1 \times [1 - \delta, 1 + \delta]$ and $\Gamma_2 = \mathcal{A}_2 \times [1 - \delta, 1 + \delta]$ where \mathcal{A}_1 is the boundary along $r(\theta) = R_j^{(1)}(\theta)$ and \mathcal{A}_2 is the boundary along $r(\theta) = R_j^{(2)}(\theta)$. Note that both \mathcal{A}_1 and \mathcal{A}_2 are composed of a finite sequence of straight line segments. Consider the surface Γ_1 . The value of γ_{jk} on the surface Γ_1 at a given θ and \tilde{r} (with r', θ' and ω held fixed) equals

(C.4)
$$\gamma_{jk}^{\Gamma_1}(\theta, \tilde{r}; r', \theta', \omega) = R_j^{(1)}(\theta) [1 - \tilde{r} \cos(\theta - \omega)] + \tilde{r} r_{jk}(r', \theta'; \omega).$$

Following the lines of Theorem 4.2, we observe that for a given \tilde{r} , $\gamma_{jk}^{\Gamma_1}(\theta, \tilde{r}; r', \theta', \omega)$ cannot be constant for a contiguous interval of θ as Equation 4.18 cannot be satisfied over any finite interval. By a similar argument, there can exist at most only a *finite* discrete set of θ for which $R_j^{(1)}(\theta) \cos(\theta - \omega) = r_{jk}(r', \theta'; \omega)$. Let \mathcal{Z} denote this finite set. Then, for a given $\theta \notin \mathcal{Z}$, $\gamma_{jk}^{\Gamma_1}$ varies linearly in \tilde{r} and specifically, its derivative with respect to \tilde{r} does not vanish. From the above observations, we can conclude that there does not exist a 2D patch on Γ_1 on which $\gamma_{jk}^{\Gamma_1}$ is constant. A similar conclusion can be obtained even for the surface Γ_2 . Hence, γ_{jk} cannot be constant on the boundary Γ over a 2D region having a finite non-zero measure.

Appendix D. Proof of Lemma 5.3.

Proof. By construction, the integrals $J_{jk}^{(2)}(\beta,\omega)$ and $J_{jk}^{(3)}(\beta,\omega)$ do not include the stationary point $\theta' = \omega$ and hence $\nabla p \neq 0$ in these integrals. Following the lines of Theorem 4.2, by defining the vector field $\mathbf{u} = \frac{\nabla p}{\|\nabla p\|^2}q$ and then applying the divergence theorem, both $J_{jk}^{(2)}(\beta,\omega)$ and $J_{jk}^{(3)}(\beta,\omega)$ can be shown to be $\tau^{\kappa_2}\zeta^{(2)}(\beta,\omega)$ and $\tau^{\kappa_3}\zeta^{(3)}(\beta,\omega)$ respectively where both κ_2 and $\kappa_3 \geq 0.5$ and $\zeta^{(2)}$ and $\zeta^{(3)}$ are some continuous bounded functions of β and ω . Hence, we can conclude that

(D.1)
$$\left|\lim_{\tau \to 0} \int_0^{2\pi} \frac{\eta_{jk}}{L^{\epsilon}} J_{jk}^{(2)}(\beta, \omega) d\omega\right| \le \lim_{\tau \to 0} \frac{\tau^{\kappa_2}}{L^{\epsilon}} \int_0^{2\pi} \left|\zeta^{(2)}(\beta, \omega)\right| d\omega = 0$$

as $|\eta_{jk} = 1|$ and similarly for $J_{jk}^{(3)}(\beta, \omega)$ for any fixed $\beta > 0$. It follows that the result also holds as $\beta \to 0$ provided the limit for β is considered after the limit for τ , i.e,

(D.2)
$$\lim_{\beta \to 0} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}}{L^{\epsilon}} J_{jk}^{(2)}(\beta, \omega) d\omega = 0, \text{ and}$$
$$\lim_{\beta \to 0} \lim_{\tau \to 0} \int_{\omega_0}^{\omega_0 + \Delta} \frac{\eta_{jk}}{L^{\epsilon}} J_{jk}^{(3)}(\beta, \omega) d\omega = 0.$$

Hence, $I_{jk}^{(1)}(\omega)$ in Equation 5.18 can be approximated by $J_{jk}^{(1)}(\beta,\omega)$ as $\beta \to 0$ and as $\tau \to 0$.

Appendix E. Proof of Lemma 5.4.

Proof. Define

(E.1)
$$\rho_{jk}(\beta,\omega) = \int_{r_k^{(-)}(\beta,\omega)}^{r_k^{(+)}(\beta,\omega)} \sqrt{r'} \sqrt{r' - \alpha_{jk}(\omega)} dr'.$$

We consider two cases, one in which j = k and another in which $j \neq k$.

case(i): If $j \neq k$, then $\alpha_{jk}(\omega)$ varies continuously with ω . Also, notice that $\rho_{jk}(\beta,\omega)$ is *independent* of τ and is also a bounded function of β and ω . The stationary point(s) of α_{jk} —denoted by $\tilde{\omega}$ —satisfy

(E.2)
$$\tan(\tilde{\omega}) = \frac{y_j - y_k}{x_j - x_k},$$

and the second derivative of $\alpha_{jk}(\omega)$ at its stationary point(s) is given by

(E.3)
$$\alpha_{jk}^{\prime\prime}(\tilde{\omega}) = -\alpha_{jk}(\tilde{\omega}).$$

For $\alpha_{jk}^{\prime\prime}(\tilde{\omega}) = 0$, we must have

(E.4)
$$\tan(\tilde{\omega}) = -\frac{x_j - x_k}{y_j - y_k} = \frac{y_j - y_k}{x_j - x_k},$$

where the last equality is obtained using Equation E.2. Rewriting, we get

(E.5)
$$\left(\frac{y_j - y_k}{x_j - x_k}\right)^2 = -1$$

which cannot be true. Since the second derivative cannot vanish at the stationary point $\tilde{\omega}$, from the one-dimensional stationary phase approximation [15], we have

(E.6)
$$\lim_{\tau \to 0} \frac{1}{L^{\epsilon}} \int_{\omega_0}^{\omega_0 + \Delta} \exp\left(\frac{-i\alpha_{jk}(\omega)}{\tau}\right) \rho_{jk}(\beta, \omega) d\omega = \lim_{\tau \to 0} O(\tau^{\kappa}) = 0$$

where $\kappa = 0.5$ or 1 depending upon whether the interval $[\omega_0, \omega_0 + \Delta)$ contains the stationary point $(\tilde{\omega})$ or not. Hence, we have $\psi_{jk}(\beta) = 0$ for $j \neq k$.

case(ii): If j = k, then $\alpha_{kk}(\omega) = 0$ and

(E.7)
$$\rho_{kk}(\beta,\omega) = \int_{r_k^{(-)}(\beta,\omega)}^{r_k^{(+)}(\beta,\omega)} r' dr',$$
$$\psi_{kk}(\beta) = \frac{1}{L^{\epsilon}} \int_{\omega_0}^{\omega_0 + \Delta} \rho_{kk}(\beta,\omega) d\omega.$$

From the definitions of $r_k^{(1)}(\beta,\omega)$ and $r_k^{(2)}(\beta,\omega)$ in Equation 5.22, we observe that

(E.8)
$$\begin{split} \lim_{\beta \to 0} r_k^{(1)}(\beta, \omega) \uparrow R_k^{(1)}(\omega), \text{ and} \\ \lim_{\beta \to 0} r_k^{(2)}(\beta, \omega) \downarrow R_k^{(2)}(\omega). \end{split}$$

Since $r_k^{(-)}(\beta,\omega) \to r_k^{(1)}(\beta,\omega)$ and $r_k^{(+)}(\beta,\omega) \to r_k^{(2)}(\beta,\omega)$ as $\beta \to 0$, we have $\lim_{\beta \to 0} r_k^{(-)}(\beta,\omega) = R_k^{(1)}(\omega), \text{ and}$ (E.9) $\lim_{\beta \to 0} r_k^{(+)}(\beta,\omega) = R_k^{(2)}(\omega).$

Since $r_k^{(-)}(\beta,\omega) \ge r_k^{(1)}(\beta,\omega)$ and $r_k^{(+)}(\beta,\omega) \le r_k^{(2)}(\beta,\omega)$ at a fixed β and r' > 0, we see that $\rho_{kk}(\beta,\omega)$ can be bounded from above by a positive *decreasing* function of β , namely,

(E.10)
$$\rho_{kk}(\beta,\omega) \le \int_{r_k^{(1)}(\beta,\omega)}^{r_k^{(2)}(\beta,\omega)} r' dr',$$

and is also independent of τ . As both $r_k^{(1)}(\beta, \omega)$ and $r_k^{(2)}(\beta, \omega)$ are also bounded functions, by the Lebesgue dominated convergence theorem, we get

(E.11)

$$\lim_{\beta \to 0} \lim_{\tau \to 0} \psi_{kk}(\beta) = \frac{1}{L^{\epsilon}} \int_{\omega_0}^{\omega_0 + \Delta} \lim_{\beta \to 0} \rho_{kk}(\beta, \omega) d\omega$$

$$= \frac{1}{L^{\epsilon}} \int_{\omega_0}^{\omega_0 + \Delta} \left\{ \int_{R_k^{(1)}(\omega)}^{R_k^{(2)}(\omega)} r' dr' \right\} d\omega$$

$$= \frac{(1 - 2\epsilon)}{L^{\epsilon}} \int_{\omega_0}^{\omega_0 + \Delta} \frac{R_k^2(\omega)}{2} d\omega.$$

Recall that $L^{\epsilon} = (1 - 2\epsilon)L$. Hence,

(E.12)
$$\sum_{j=1}^{K} \sum_{k=1}^{K} \lim_{\beta \to 0} \lim_{\tau \to 0} \psi_{jk}(\beta) = \frac{1}{L} \sum_{k=1}^{K} \int_{\omega_0}^{\omega_0 + \Delta} \frac{R_k^2(\omega)}{2} d\omega$$
$$= \int_{\omega_0}^{\omega_0 + \Delta} P(\omega) d\omega$$

which completes the proof. \square

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24

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