Detection and localization of peaks in a smooth random field

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Peak inference



Peak Detection

Smoothing

• Suppose we have n subjects and that for each subject we observe i.i.d

$$X_i = \mu_X + \epsilon_i$$

on a finite lattice $\mathcal{L} \subset S \subset \mathbb{R}^D$, where $\mu_X : \mathcal{L} \to \mathbb{R}$.

• Smoothing is typically done in order to increase the signal to noise ratio. I.e. for each i, X_i is smoothed with a C^2 kernel K to give

$$Y_i(v) = \sum_{l \in \mathcal{L}} K(v-l)\mu_X(l) + \sum_{l \in \mathcal{L}} K(v-l)\epsilon_i(l)$$

at each $v \in \mathcal{L}$.

• Let

$$\hat{\mu}_n = \frac{1}{n} \sum_{i=1}^n Y_i = \sum_{l \in \mathcal{L}} K(v-l) \overline{X}_n(l)$$

where $\overline{X}_n = \sum_{i=1}^n X_i$.

Why you should smooth



Lattice smoothing

In applications when smoothing is performed (e.g. in fMRI) you evaluate

$$\hat{\mu}_n(v) = \sum_{l \in \mathcal{L}} K(v-l) \overline{X}_n(l)$$

at every voxel $v \in \mathcal{L}$. An example evaluation $\hat{\mu}_n$ is shown below.



Convolution Random Fields

Definition

Instead we can evaluate the smoothing continuously, meaning that for all $s \in S$,

$$\hat{\mu}_n(s) := \sum_{l \in \mathcal{L}} K(s-l) \overline{X}_n(l).$$

We call this a convolution field.



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Taking slices through a 3D convolution field generated from brain imaging data, you get the following images!



Voxelwise RFT

Let $M_u(\hat{\mu}_n)$ be the number of local maxima of $\hat{\mu}_{Y,n}$ above a threshold u then

$$\mathbb{P}\left(\sup_{s\in S}\hat{\mu}_n(s) > u\right) = \mathbb{P}(M_u(\hat{\mu}_n) \ge 1) \le \mathbb{E}[M_u(\hat{\mu}_n)]$$

because $\hat{\mu}$ exceeds u if and only if there is at least one local maxima above u.

- Many good approximations exist to $\mathbb{E}[M_u(\hat{\mu}_n)]$ at high thresholds u so can control FWER.
- FWER over space is the same as FWER over peaks so we can use this to find significant peaks.
- Using convolution fields to do inference means that RFT works. Doing the same thing on a lattice leads to conservativeness as

$$\sup_{v \in \mathcal{L}} \hat{\mu}_n(v) < \sup_{s \in S} \hat{\mu}_n(s).$$
(1)

FWER for 2D Gaussian random fields

We generate mean-zero random fields by smoothing Gaussian white noise with an isotropic Gaussian kernel and take n = 40. (FWHM is proportional to the smoothing parameter of the kernel.)



Thus RFT provides accurate inference (and is much faster than e.g. bootstrap/permutation resampling methods).

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Peak CRs

Detection and Localization



Peak Confidence Regions

Asymptotic confidence regions

Define $\mu: S \to \mathbb{R}$ s.t. $\mu(s) = \sum_{l \in \mathcal{L}} K(l-s)\mu_X(l)$. Assume that μ has J peaks at locations $\theta_1, \ldots, \theta_J$ within disjoint balls B_1, \ldots, B_J .

Theorem

For each j = 1, ..., J corresponding to a local maximum of μ , let $\hat{\theta}_{j,n} = \operatorname{argmax}_{t \in B_j} \hat{\mu}_n(t)$ (and for the minima let $\hat{\theta}_{j,n} = \operatorname{argmin}_{t \in B_j} \hat{\mu}_n(t)$) and let $\hat{\theta}_n := (\hat{\theta}_{1,n}^T, ..., \hat{\theta}_{J,n}^T)^T$ and $\boldsymbol{\theta} := (\theta_1^T, ..., \theta_J^T)^T$. Then, under regularity assumptions on μ and the noise,

$$\sqrt{n}(\hat{\boldsymbol{\theta}}_{\boldsymbol{n}} - \boldsymbol{\theta}) \stackrel{d}{\longrightarrow} N(0, \boldsymbol{A}\boldsymbol{\Lambda}\boldsymbol{A}^T)$$

as $n \to \infty$. Here $\mathbf{A} \in \mathbb{R}^{DJ \times DJ}$ depends on $\nabla^2 \mu$ and $\mathbf{\Lambda} \in \mathbb{R}^{DJ \times DJ}$ depends on the covariance of ∇Y_1 .

Here ∇ denotes the derivative.

Accounting for the variance

Like for the MLE, Taylor expanding $\nabla^T \mu_n$ around $\hat{\theta}_{n,j}$ we have,

$$\nabla^T \hat{\mu}_n(\hat{\theta}_{j,n}) = 0 = \nabla^T \hat{\mu}_n(\theta_j) + \nabla^2 \hat{\mu}_n(\theta_{j,n}^*)(\hat{\theta}_{j,n} - \theta_j)$$

and so

$$\hat{\theta}_{j,n} - \theta_j = -\left(\nabla^2 \hat{\mu}_n(\theta_{j,n}^*)\right)^{-1} \nabla^T \hat{\mu}_n(\theta_j)$$

In particular

$$\sqrt{n}(\hat{\theta}_{j,n} - \theta_j) \xrightarrow{d} \mathcal{N}(0, (\nabla^2 \mu(\theta_j))^{-1} \Lambda(\theta_j) (\nabla^2 \mu(\theta_j))^{-1})$$

where $\Lambda(\theta_j) = \operatorname{cov}(Y_1(\theta_j))$. In practice to make an asymptotic confidence region one approximations $(\nabla^2 \mu(\theta_j))^{-1}$ by $\left(\nabla^2 \hat{\mu}_n(\hat{\theta}_{j,n})\right)^{-1}$.

$$\hat{\theta}_{j,n} - \theta_j = -\left(\nabla^2 \hat{\mu}_n(\theta_j) + \frac{1}{2}(\hat{\theta}_{j,n} - \theta_j)^T \nabla^3 \hat{\mu}_n(\tilde{\theta}_{j,n})\right)^{-1} \nabla^T \hat{\mu}_n(\theta_j)$$
$$\approx -\left(\nabla^2 \hat{\mu}_n(\theta_j)\right)^{-1} \nabla^T \hat{\mu}_n(\theta_j)$$

We have

$$\begin{pmatrix} \nabla^T \hat{\mu}_n(\theta_j) \\ \mathbf{vech}(\nabla^2 \hat{\mu}_n(\theta_j)) \end{pmatrix} \sim \mathcal{N} \left(\begin{pmatrix} 0 \\ \mathbf{vech}(\nabla^2 \mu_n(\theta_j)) \end{pmatrix}, \frac{1}{n} \begin{pmatrix} \Lambda & 0 \\ 0 & \Omega \end{pmatrix} \right)$$

and for $1 \le k \le K$ $(K \in \mathbb{N})$ we can approximate this by simulating from the following distribution

$$\begin{pmatrix} A_k \\ B_k \end{pmatrix} \sim \mathcal{N} \left(\begin{pmatrix} 0 \\ \mathbf{vech}(\nabla^2 \hat{\mu}_n(\hat{\theta}_{j,n})) \end{pmatrix}, \frac{1}{n} \begin{pmatrix} \hat{\Lambda} & 0 \\ 0 & \hat{\Omega} \end{pmatrix} \right).$$

and calculating $\delta_{k,n} = (\mathbf{vech}^{-1}(B_{k,n}))^{-1}A_{k,n}$.

Let $\hat{\Sigma}'_j = (\nabla^2 \hat{\mu}_n(\hat{\theta}_j))^{-1} \hat{\Lambda} (\nabla^2 \hat{\mu}_n(\hat{\theta}_j))^{-1}$ and for $0 < \alpha < 1$, choose λ_α such that

$$\frac{1}{K}\sum_{k=1}^{K} \mathbb{1}\Big[n(\hat{\delta}_{k,n}^{T}(\hat{\Sigma}_{j}')^{-1}\delta_{k,n}) > \lambda_{\alpha}\Big] = \frac{\lfloor \alpha K \rfloor}{K}.$$

Given this we define a $(1 - \alpha)$ Monte Carlo confidence region to be

$$\Big\{\theta: n(\hat{\theta}_{j,n}-\theta)^T(\hat{\Sigma}'_j)^{-1}(\hat{\theta}_{j,n}-\theta) < \lambda_\alpha\Big\}.$$

- These regions are asymptotically valid (for the same reason as the asymptotic cases)
- Under stationarity we can prove that these intervals are bigger than the asymptotic ones.

Example simulations



Figure 1: Left: True signal. Right: one realisation.

Simulate n random fields about the signal 5000 times and compute the coverage of the regions.



Application to fMRI



Figure 3: Peaks of the mean of 125 subjects

- Using convolution fields accurately and quickly controls the FWER at the right level. (This is valid under non-stationarity.)
- Here we assumed Gaussianity but the framework works for t-fields and F-fields as well.
- You can derive asymptotic confidence regions and improve upon these if you additionally assume stationarity.
- These results are valid in full generality not just in the convolution framework.
- Software (in MATLAB) to perform inference on random fields is available at the RFTtoolbox (github.com/sjdavenport/RFTtoolbox). (E.g. for LKC estimation, Peak Inference, Peak Height distribution, confidence regions)
- Slides available at sjdavenport.github.io/talks.

- Pre-print on convolution fields soon to be out, in the mean time you can read a lot of the details in my thesis here: https://sjdavenport.github.io/research/papers/thesis.pdf.
- Pre-print on peak confidence regions is available on arxiv (Davenport, Nichols, & Schwarzman, 2022).

Additional simulations



Figure 4: Left: True signal. Right: one realisation.

Additional simulation coverages



Theorem

Suppose that A and B are independent real valued random variables with well defined densities p_A and p_B which are symmetric about $\mathbb{E}[A]$ and $\mathbb{E}[B]$ respectively. Assume that $p_A(x)$ is decreasing for x > 0 and increasing for x < 0, B is positive and that $\mathbb{E}[|B|] < \infty$. Then for all x > 0,

$$\mathbb{P}\bigg(\frac{A}{\mathbb{E}[B]} > x\bigg) \le \mathbb{P}\bigg(\frac{A}{B} > x\bigg).$$

Davenport, S., Nichols, T. E., & Schwarzman, A. (2022). Confidence regions for the location of peaks of a smooth random field. arXiv preprint arXiv:2208.00251.