



Aalto University
School of Electrical
Engineering

Stochastic (partial) differential equations and Gaussian processes

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- 1 Basic ideas
- 2 Stochastic differential equations and Gaussian processes
- 3 Stochastic partial differential equations and Gaussian processes
- 4 Conclusion

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Kernel vs. SPDE representations of GPs

GP model $\mathbf{x} \in \mathbb{R}^d, t \in \mathbb{R}$	Equivalent S(P)DE model
Spatial $k(\mathbf{x}, \mathbf{x}')$	SPDE model (\mathcal{L} is an operator) $\mathcal{L} f(\mathbf{x}) = w(\mathbf{x})$
Temporal $k(t, t')$	State-space/SDE model $\frac{d\mathbf{f}(t)}{dt} = \mathbf{A} \mathbf{f}(t) + \mathbf{L} w(t)$
Spatio-temporal $k(\mathbf{x}, t; \mathbf{x}', t')$	Stochastic evolution equation $\frac{\partial}{\partial t} \mathbf{f}(\mathbf{x}, t) = \mathcal{A}_x \mathbf{f}(\mathbf{x}, t) + \mathbf{L} w(\mathbf{x}, t)$

Why use S(P)DE solvers for GPs?

- The $O(n^3)$ computational complexity is a challenge.
- What do we get:
 - $O(n)$ state-space methods for SDEs/SPDEs.
 - Sparse approximations developed for SPDEs.
 - Reduced rank Fourier/basis function approximations.
 - Path to non-Gaussian processes.
- Downsides:
 - We often need to approximate.
 - Mathematics can become messy.

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Ornstein-Uhlenbeck process

- The mean and covariance functions:

$$m(x) = 0$$

$$k(x, x') = \sigma^2 \exp(-\lambda|x - x'|)$$

- This has a *path representation* as a stochastic differential equation (SDE):

$$\frac{df(t)}{dt} = -\lambda f(t) + w(t).$$

where $w(t)$ is a white noise process with x relabeled as t .

- Ornstein-Uhlenbeck process is a Markov process.
- What does this actually mean \implies white board.

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Ornstein-Uhlenbeck process (cont.)

- Consider a Gaussian process regression problem

$$f(x) \sim \text{GP}(0, \sigma^2 \exp(-\lambda|x - x'|))$$
$$y_k = f(x_k) + \varepsilon_k$$

- This is equivalent to the state-space model

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that is, with $f_k = f(t_k)$ we have a Gauss-Markov model

$$f_{k+1} \sim p(f_{k+1} | f_k)$$
$$y_k \sim p(y_k | f_k)$$

- Solvable in $O(n)$ time using Kalman filter/smoothing.

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State Space Form of Linear Time-Invariant SDEs

- Consider a N th order LTI SDE of the form

$$\frac{d^N f}{dt^N} + a_{N-1} \frac{d^{N-1} f}{dt^{N-1}} + \dots + a_0 f = w(t).$$

- If we define $\mathbf{f} = (f, \dots, d^{N-1} f / dt^{N-1})$, we get a state space model:

$$\frac{d\mathbf{f}}{dt} = \underbrace{\begin{pmatrix} 0 & 1 & & \\ & \ddots & \ddots & \\ & & 0 & 1 \\ -a_0 & -a_1 & \dots & -a_{N-1} \end{pmatrix}}_{\mathbf{A}} \mathbf{f} + \underbrace{\begin{pmatrix} 0 \\ \vdots \\ 0 \\ 1 \end{pmatrix}}_{\mathbf{L}} w(t)$$
$$f(t) = \underbrace{\begin{pmatrix} 1 & 0 & \dots & 0 \end{pmatrix}}_{\mathbf{H}} \mathbf{f}.$$

- The vector process $\mathbf{f}(t)$ is Markovian although $f(t)$ isn't.

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Spectra of Linear Time-Invariant SDEs

- By taking the Fourier transform of the LTI SDE, we can derive the **spectral density** which has the form:

$$S(\omega) = \frac{\text{(constant)}}{\text{(polynomial in } \omega^2)}$$

- We can also do this conversion to the **other direction**:

- With certain parameter values, the Matérn has the form:

$$S(\omega) \propto (\lambda^2 + \omega^2)^{-(\rho+1)}$$

- Many non-rational spectral densities can be approximated:

$$S(\omega) = \sigma^2 \sqrt{\frac{\pi}{\kappa}} \exp\left(-\frac{\omega^2}{4\kappa}\right) \approx \frac{\text{(const)}}{N! / 0! (4\kappa)^N + \dots + \omega^{2N}}$$

- For the conversion of a rational spectral density to a Markovian (state-space) model, we can use the **spectral factorization**.

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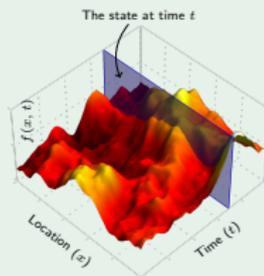
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State-space methods for Gaussian processes

- Approximation:

$$S(\omega) \approx \frac{b_0 + b_1 \omega^2 + \dots + b_M \omega^{2M}}{a_0 + a_1 \omega^2 + \dots + a_N \omega^{2N}}$$



- Results in a linear stochastic differential equation (SDE)

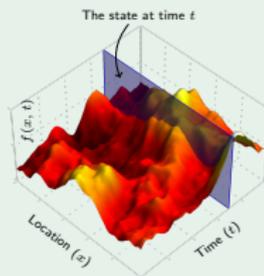
$$df(t) = \mathbf{A} f(t) dt + \mathbf{L} d\mathbf{W}$$

- More generally stochastic evolution equations.
- $O(n)$ GP regression with Kalman filters and smoothers.
- Parallel block-sparse precision methods $\rightarrow O(\log n)$.

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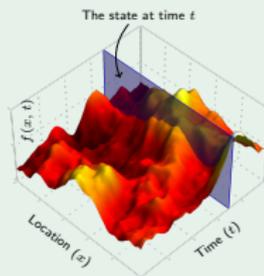
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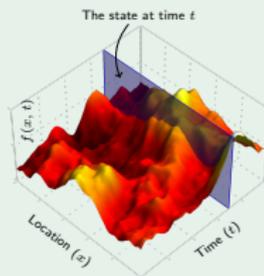
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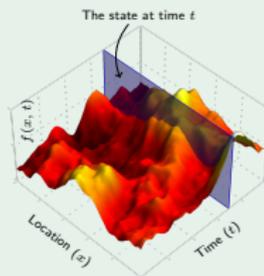
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State-space methods – temporal example

Example (Matérn class 1d)

The Matérn class of covariance functions is

$$k(t, t') = \sigma^2 \frac{2^{1-\nu}}{\Gamma(\nu)} \left(\frac{\sqrt{2\nu}}{\ell} |t - t'| \right)^\nu K_\nu \left(\frac{\sqrt{2\nu}}{\ell} |t - t'| \right).$$

When, e.g., $\nu = 3/2$, we have

$$d\mathbf{f}(t) = \begin{pmatrix} 0 & 1 \\ -\lambda^2 & -2\lambda \end{pmatrix} \mathbf{f}(t) dt + \begin{pmatrix} 0 \\ q^{1/2} \end{pmatrix} dW(t),$$
$$f(t) = (1 \quad 0) \mathbf{f}(t).$$

State-space methods – spatio-temporal example

Example (2D Matérn covariance function)

- Consider a space-time Matérn covariance function

$$k(x, t; x', t') = \sigma^2 \frac{2^{1-\nu}}{\Gamma(\nu)} \left(\sqrt{2\nu} \frac{\rho}{l} \right)^\nu K_\nu \left(\sqrt{2\nu} \frac{\rho}{l} \right).$$

where we have $\rho = \sqrt{(t - t')^2 + (x - x')^2}$, $\nu = 1$ and $d = 2$.

- We get the following representation:

$$df(x, t) = \begin{pmatrix} 0 & 1 \\ \frac{\partial^2}{\partial x^2} - \lambda^2 & -2\sqrt{\lambda^2 - \frac{\partial^2}{\partial x^2}} \end{pmatrix} \mathbf{f}(x, t) dt + \begin{pmatrix} 0 \\ 1 \end{pmatrix} dW(x, t).$$

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Basic idea of SPDE inference on GPs [1/2]

- Consider e.g. the stochastic partial differential equation:

$$\frac{\partial^2 f(x, y)}{\partial x^2} + \frac{\partial^2 f(x, y)}{\partial y^2} - \lambda^2 f(x, y) = w(x, y)$$

- Fourier transforming gives the spectral density:

$$S(\omega_x, \omega_y) \propto (\lambda^2 + \omega_x^2 + \omega_y^2)^{-2}.$$

- Inverse Fourier transform gives the covariance function:

$$k(x, y; x', y') = \frac{\sqrt{(x - x')^2 + (y - y')^2}}{2\lambda} K_1(\lambda \sqrt{(x - x')^2 + (y - y')^2})$$

- But this is just the Matérn covariance function.

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Basic idea of SPDE inference on GPs [2/2]

- More generally, SPDE for some linear operator \mathcal{L} :

$$\mathcal{L} f(\mathbf{x}) = w(\mathbf{x})$$

- Now f is a GP with precision and covariance operators:

$$\mathcal{K}^{-1} = \mathcal{L}^* \mathcal{L}$$

$$\mathcal{K} = (\mathcal{L}^* \mathcal{L})^{-1}$$

- Idea: approximate \mathcal{L} or \mathcal{L}^{-1} using PDE/ODE methods:
 - Finite-differences/FEM methods lead to sparse precision approximations.
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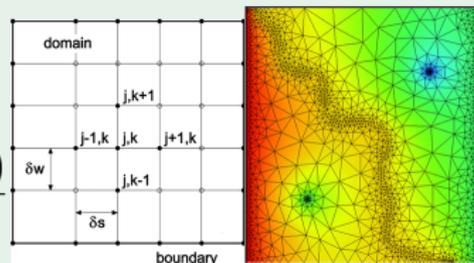
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$$\frac{\partial f(x)}{\partial x} \approx \frac{f(x+h) - f(x)}{h}$$
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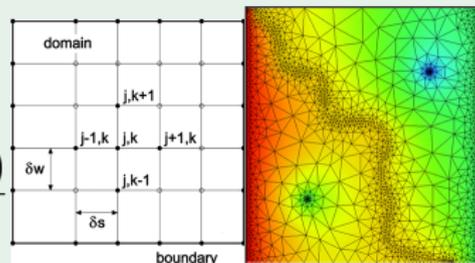
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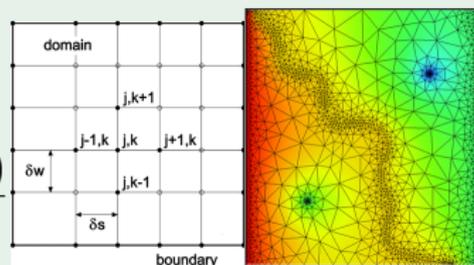
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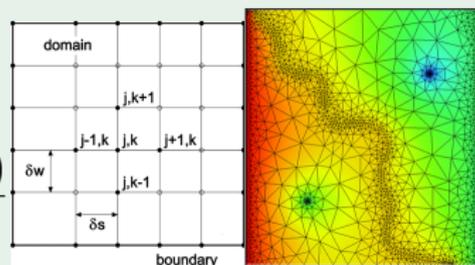
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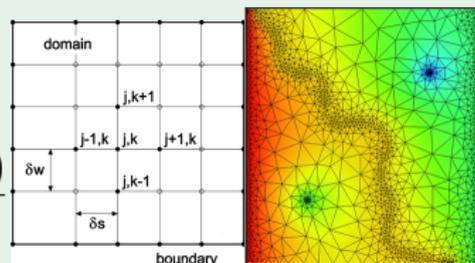
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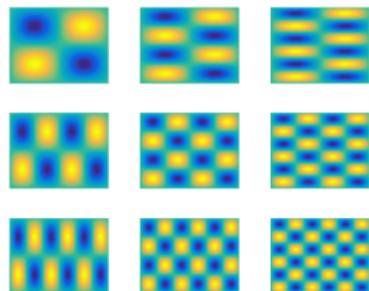
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Classical and random Fourier methods – reduced rank approximations and FFT

- Approximation:

$$f(\mathbf{x}) \approx \sum_{\mathbf{k} \in \mathbb{N}^d} \alpha_{\mathbf{k}} \exp(2\pi i \mathbf{k}^T \mathbf{x})$$

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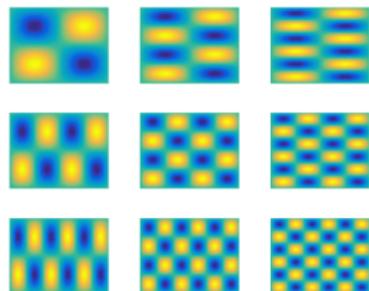
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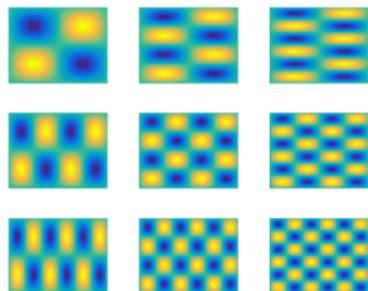
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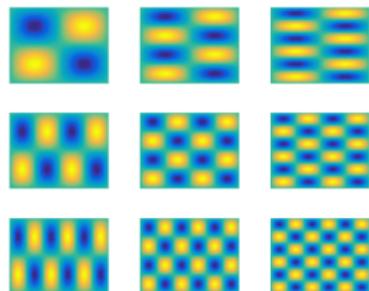
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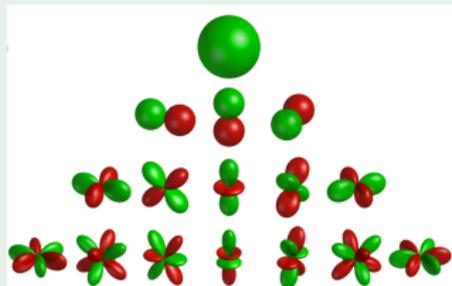
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- Approximation:

$$f(\mathbf{x}) \approx \sum_i c_i \phi_i(\mathbf{x})$$

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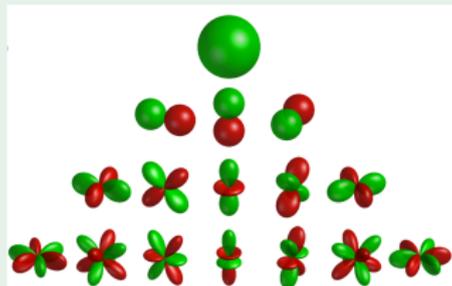
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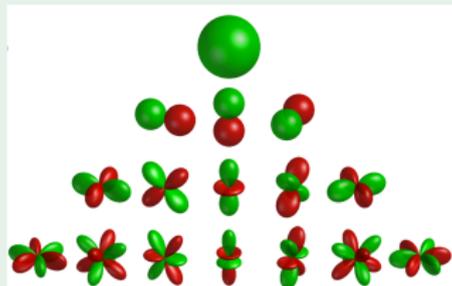
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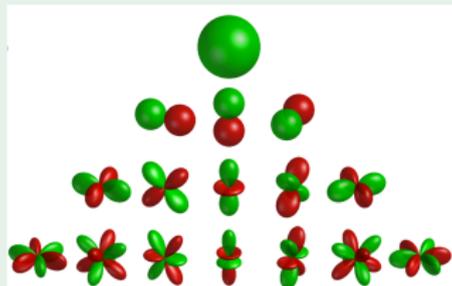
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Contents

- 1 Basic ideas
- 2 Stochastic differential equations and Gaussian processes
- 3 Stochastic partial differential equations and Gaussian processes
- 4 Conclusion

Back to SPDE representations of GPs

GP model $\mathbf{x} \in \mathbb{R}^d, t \in \mathbb{R}$	Equivalent S(P)DE model
Spatial $k(\mathbf{x}, \mathbf{x}')$	SPDE model (\mathcal{L} is an operator) $\mathcal{L} f(\mathbf{x}) = w(\mathbf{x})$
Temporal $k(t, t')$	State-space/SDE model $\frac{d\mathbf{f}(t)}{dt} = \mathbf{A} \mathbf{f}(t) + \mathbf{L} w(t)$
Spatio-temporal $k(\mathbf{x}, t; \mathbf{x}', t')$	Stochastic evolution equation $\frac{\partial}{\partial t} \mathbf{f}(\mathbf{x}, t) = \mathcal{A}_x \mathbf{f}(\mathbf{x}, t) + \mathbf{L} w(\mathbf{x}, t)$

What then?

- Exchange and map approximations between the fields:
 - Inducing points \leftrightarrow point-collocation; spectral methods \leftrightarrow Galerkin methods; finite-differences \leftrightarrow GMRFs;
- Non-Gaussian processes: Student's-t processes, non-linear Itô processes, jump processes, hybrid point/Gaussian processes.
- Hierarchical (deep) SPDE models: we stack SPDEs on top of each other – the SPDE just becomes non-linear.
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Useful references

- **N. Wiener (1950)**. *Extrapolation, Interpolation and Smoothing of Stationary Time Series with Engineering Applications*. John Wiley & Sons, Inc.
- **R. L. Stratonovich (1963)**. *Topics in the Theory of Random Noise*. Gordon and Breach.
- **J. Hartikainen and S. Särkkä (2010)**. *Kalman Filtering and Smoothing Solutions to Temporal Gaussian Process Regression Models*. Proc. MLSP.
- **S. Särkkä, A. Solin, and J. Hartikainen (2013)**. *Spatio-Temporal Learning via Infinite-Dimensional Bayesian Filtering and Smoothing*. IEEE Sig.Proc.Mag., 30(5):51–61.
- **S. Särkkä (2013)**. *Bayesian Filtering and Smoothing*. Cambridge University Press.
- **S. Särkkä and A. Solin (2017, to appear)** *Applied Stochastic Differential Equations*. Cambridge University Press.