

JIDT: An information-theoretic toolkit for studying the dynamics of complex systems



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European Conference on Artificial Life (ECAL)

York, UK

July 20, 2015

Java Information Dynamics Toolkit (JIDT)



On google code → github (<https://lazier.me/jidt/>)

JIDT provides a standalone, open-source ([GPL v3 licensed](#)) implementation of information-theoretic measures of information processing in complex systems, i.e. information storage, transfer and modification.

JIDT includes implementations:

- Principally for transfer entropy, mutual information, their conditional variants, active information storage etc;
- For both discrete and continuous-valued data;
- Using various types of estimators (e.g. Kraskov-Stögbauer-Grassberger, linear-Gaussian, etc.).

Java Information Dynamics Toolkit (JIDT)



JIDT is written in Java but directly usable in Matlab/Octave, Python, R, Julia, Clojure, etc.

JIDT requires almost zero installation.

JIDT is distributed with:

- A paper describing its design and usage;
 - J.T. Lizier, *Frontiers in Robotics and AI* 1:11, 2014; (arXiv:1408.3270)
- A full tutorial and exercises
- Full Javadocs;
- A suite of demonstrations, including in each of the languages listed above.

Code credits: JL, Ipek Özdemir, Pedro Martínez Mediano

JIDT tutorial – Objectives



Participants will:

- Understand measures of information dynamics;
- Be able to obtain and install JIDT distribution;
- Understand and run sample scripts in their chosen environment;
- Be able to generate new code or modify sample scripts for new analysis;
- Know how and where to seek support information (wiki, Javadocs, mailing list, twitter).

- 1 Introduction
- 2 Information dynamics
 - Information theory
 - The information dynamics framework
- 3 Estimation techniques
- 4 Overview of JIDT
- 5 Demos
- 6 Exercise
- 7 Wrap-up

Hi, I'm Joe ..



- My background.
- Why I started this information-theoretic toolkit project.

Tell me about yourselves ...



- Where are you from? Unis, other organisations?

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- How familiar are you with information theory – what measures do you know? Are you using it for analysis yet?

Tell me about yourselves ...



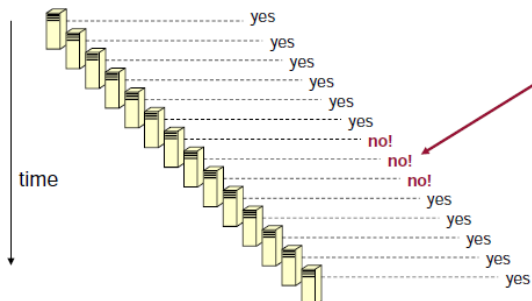
- Where are you from? Unis, other organisations?
- Which languages/environments are you using, and how much coding experience do you have?
- How familiar are you with information theory – what measures do you know? Are you using it for analysis yet?
- What types of information-theoretic analysis do you have planned (perhaps with JIDT)?

Entropy

(Shannon) entropy is a measure of **uncertainty**.

Intuitively: if we want to know the value of a variable, there is uncertainty in what that value is before we inspect it (measured in **bits**).

- e.g. Is the web server cam.ac.uk running?



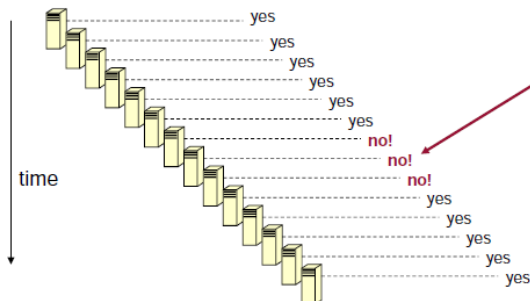
1. There is more uncertainty in more rare events

2. There is more average uncertainty when there is a balance in event probabilities

$$H(X) = - \sum_{x \in \mathcal{X}} P(x) \log P(x)$$

Sample entropy calculation

- e.g. Is the web server cam.ac.uk running?



1. There is more uncertainty in more rare events

2. There is more average uncertainty when there is a balance in event probabilities

$$H(X) = - \sum_{x \in \mathcal{X}} P(x) \log P(x)$$

Here we have $p(\text{yes}) = 11/14$, $p(\text{no}) = 3/14$ – what is H ?

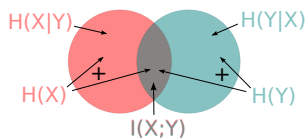
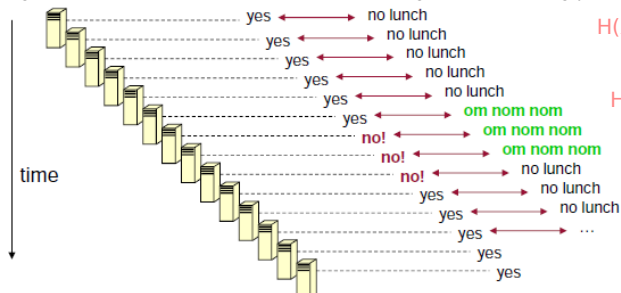
$x = \text{yes no}$	$p(x)$	$-\log_2 p(x)$	$-p(x) \log_2 p(x)$
yes	0.786	0.348	0.273
no	0.214	2.22	0.476
			$H = 0.750$ bits

Conditional entropy

Uncertainty in one variable X **in the context of** the known measurement of another variable Y .

Intuitively: how much uncertainty is there in X after we know the value of Y ?

e.g. How uncertain are we about the web server is running if we know the IT guy is at lunch?



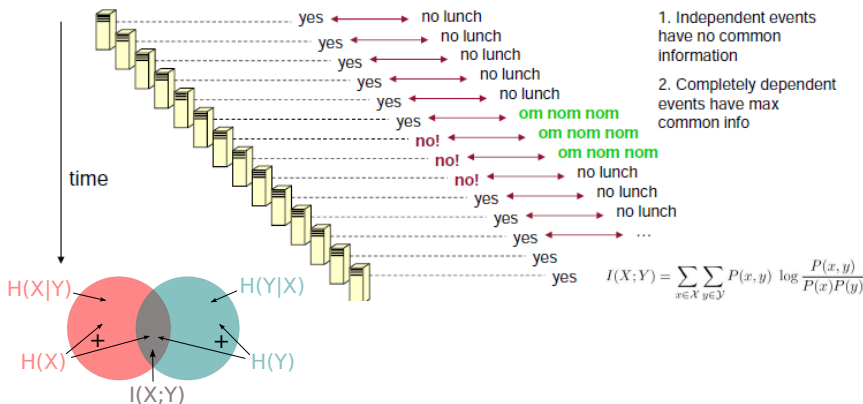
$$H(X|Y) = H(X, Y) - H(Y) = - \sum_{x \in X} \sum_{y \in Y} p(x, y) \log p(x|y)$$

Mutual information

Information is a measure of uncertainty **reduction**.

Intuitively: common information is the amount that knowing the value of one variable tells us about another.

- e.g. How much common info b/w if IT guy is at lunch and the web server running?



Information-theoretic quantities

Shannon entropy

$$\begin{aligned} H(X) &= - \sum_x p(x) \log_2 p(x) \\ &= \langle -\log_2 p(x) \rangle \end{aligned}$$

Conditional entropy

$$H(X|Y) = - \sum_{x,y} p(x,y) \log_2 p(x|y)$$

Mutual information (MI)

$$\begin{aligned} I(X; Y) &= H(X) + H(Y) - H(X, Y) \\ &= \sum_{x,y} p(x,y) \log_2 \frac{p(x|y)}{p(x)} \\ &= \left\langle \log_2 \frac{p(x|y)}{p(x)} \right\rangle \end{aligned}$$

Conditional MI

$$\begin{aligned} I(X; Y|Z) &= H(X|Z) + H(Y|Z) - H(X, Y|Z) \\ &= \left\langle \log_2 \frac{p(x|y,z)}{p(x|z)} \right\rangle \end{aligned}$$

Local measures



We can write **local** (or point-wise) information-theoretic measures for specific observations/configurations $\{x, y, z\}$:

$$h(x) = -\log_2 p(x), \quad i(x; y) = \log_2 \frac{p(x|y)}{p(x)}$$

$$h(x|y) = -\log_2 p(x|y), \quad i(x; y|z) = \log_2 \frac{p(x|y, z)}{p(x|z)}$$

- We have $H(X) = \langle h(x) \rangle$ and $I(X; Y) = \langle i(x; y) \rangle$, etc.
- If X, Y, Z are time-series, local values measure **dynamics** over time.

What can we do with these measures in ALife/CI?



- Measure the diversity in agent strategies (Miramontes, 1995; Prokopenko et al., 2005).
- Measure long-range correlations as we approach a phase-transition (Ribeiro et al., 2008).
- Feature selection for machine learning (Wang et al., 2014).
- Quantify the information held in a response about a stimulus, and indeed about specific stimuli (DeWeese and Meister, 1999).
- Measure the common information in the behaviour of two agents (Sperati et al., 2008).
- Guide self-organisation using these measures (Prokopenko et al., 2006).
- ...

→ Information theory is useful for answering specific questions about information content, shared information, and where and to what extent information about some variable is mirrored.

Information dynamics



We *talk* about computation as:

- Memory
- Signalling
- Processing

Distributed computation is any process involving these features:

- Time evolution of cellular automata
- Information processing in the brain
- Gene regulatory networks computing cell behaviours
- Flocks computing their collective heading
- Ant colonies computing the most efficient routes to food
- The universe is computing its own future!

Information dynamics

We *talk* about computation as:

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Idea: quantify computation via:

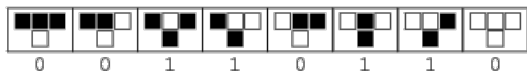
- Information **storage**
- Information **transfer**
- Information **modification**

Distributed computation is any process involving these features:

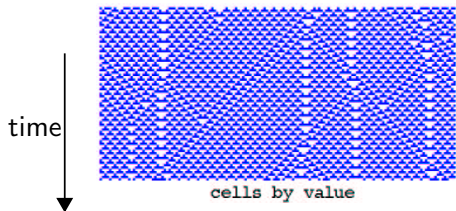
- Time evolution of cellular automata
- Information processing in the brain
- Gene regulatory networks computing cell behaviours
- Flocks computing their collective heading
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- The universe is computing its own future!

General idea: by quantifying intrinsic computation in the language it is normally described in, we can understand how nature computes and why it is complex.

Motivating example: cellular automata



CAs: simple dynamical systems;
known causal structure and rules.

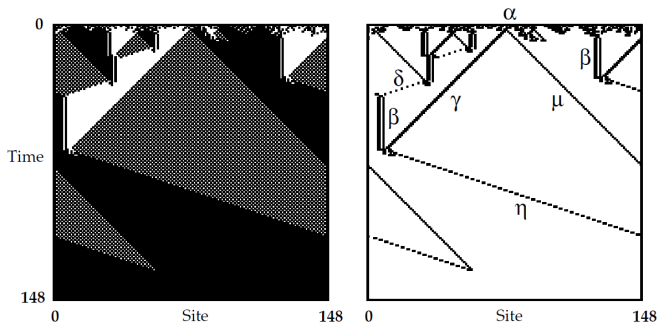


(Wuensche, 1999)

Emergent structure:

- Domain, blinkers
- Particles
 - Gliders, domain walls
- Collisions

Motivating example: cellular automata



Mitchell et al. (1994, 1996) used GAs to evolve CAs to solve specific computational tasks.

In attempting the density classification task (above), the CA uses:

- domains and blinkers β to **store** information;
- gliders γ, η to **transfer** information;
- glider collisions e.g. $\gamma + \beta \rightarrow \eta$ to **modify/process** information.

Information dynamics

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Information dynamics

- Information **storage**
- Information **transfer**
- Information **modification**

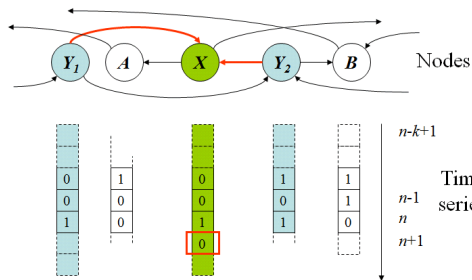
Key properties of the **information dynamics** approach:

- A focus on individual operations of computation rather than overall complexity;
- Alignment with descriptions of dynamics in specific domains;
- A focus on the **local scale** of info **dynamics** in space-time;
- Information-theoretic basis directly measures computational quantities:
 - Captures non-linearities;
 - Is applicable to, and comparable between, any type of time-series.

Information dynamics



Key question: how is the next state of a variable in a complex system **computed**?



Q: Where does the information in x_{n+1} come from, and how can we measure it?

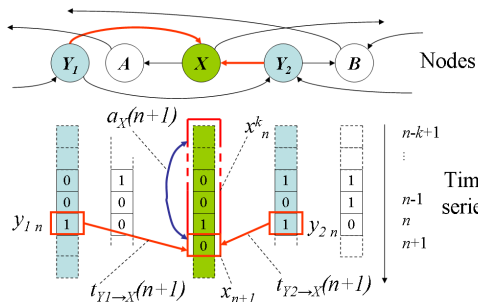
Q: How much was stored, how much was transferred, can we partition them or do they overlap?

Complex system as a multivariate **time-series** of states

Information dynamics



Studies computation of the next state of a target variable in terms of information **storage**, **transfer** and **modification**: (Lizier et al., 2008, 2010, 2012)



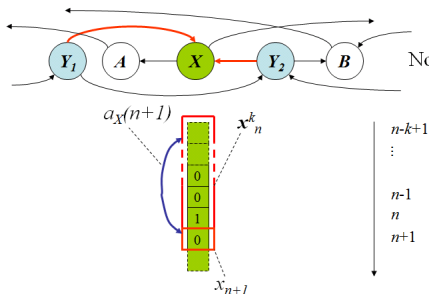
The measures examine:

- **State** updates of a target variable;
- **Dynamics** of the measures in space and time.

Active information storage (Lizier et al., 2012)



How much information about the next observation X_{n+1} of process X can be found in its past **state** $\mathbf{X}_n^{(k)} = \{X_{n-k+1} \dots X_{n-1}, X_n\}$?



Nodes **Active information storage:**

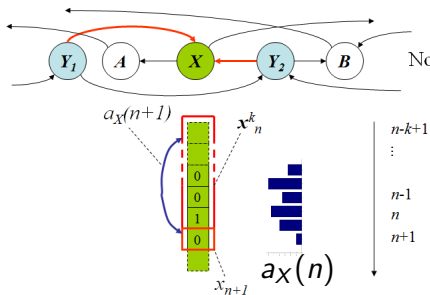
$$A_X = I(X_{n+1}; \mathbf{X}_n^{(k)}) = \left\langle \log_2 \frac{p(X_{n+1} | \mathbf{X}_n^{(k)})}{p(X_{n+1})} \right\rangle$$

Average information from past **state** that is in use in predicting the next value.

Active information storage (Lizier et al., 2012)



How much information about the next observation X_{n+1} of process X can be found in its past **state** $\mathbf{x}_n^{(k)} = \{X_{n-k+1} \dots X_{n-1}, X_n\}$?



$$A_X = \langle a_X(n) \rangle$$

Active information storage:

$$A_X = I(X_{n+1}; \mathbf{x}_n^{(k)}) = \left\langle \log_2 \frac{p(x_{n+1} | \mathbf{x}_n^{(k)})}{p(x_{n+1})} \right\rangle$$

Average information from past **state** that is in use in predicting the next value.

Local active information storage:

$$a_X(n) = \log_2 \frac{p(x_{n+1} | \mathbf{x}_n^{(k)})}{p(x_{n+1})}$$

Information from a **specific** past **state** that is in use in predicting the **specific** next value.

Information transfer

How much information about the **state transition** $\mathbf{X}_n^{(k)} \rightarrow X_{n+1}$ of X can be found in the past **state** $\mathbf{Y}_n^{(l)}$ of a source process Y ?

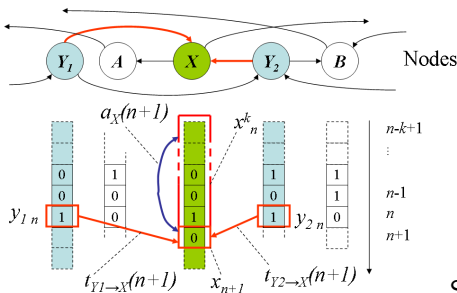
Transfer entropy: (Schreiber, 2000)

$$T_{Y \rightarrow X} = I(\mathbf{Y}_n^{(l)}; X_{n+1} | \mathbf{X}_n^{(k)})$$

$$= \left\langle \log_2 \frac{p(x_{n+1} | \mathbf{x}_n^{(k)}, \mathbf{y}_n^{(l)})}{p(x_{n+1} | \mathbf{x}_n^{(k)})} \right\rangle$$

Average info from source that helps predict next value in context of past.

Storage and transfer are **complementary**:
 $H_X = A_X + T_{Y \rightarrow X} + \text{higher order terms}$



Information transfer

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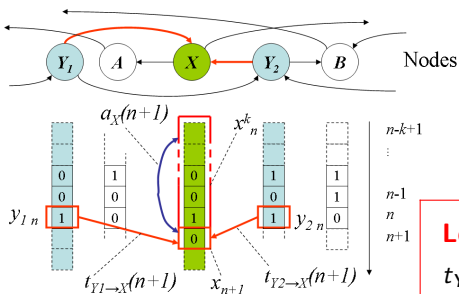
$$= \left\langle \log_2 \frac{p(x_{n+1} | \mathbf{x}_n^{(k)}, \mathbf{y}_n^{(l)})}{p(x_{n+1} | \mathbf{x}_n^{(k)})} \right\rangle$$

Average info from source that helps predict next value in context of past.

Local transfer entropy: (Lizier et al., 2008)

$$t_{Y \rightarrow X}(n) = \log_2 \frac{p(x_{n+1} | \mathbf{x}_n^{(k)}, \mathbf{y}_n^{(l)})}{p(x_{n+1} | \mathbf{x}_n^{(k)})}$$

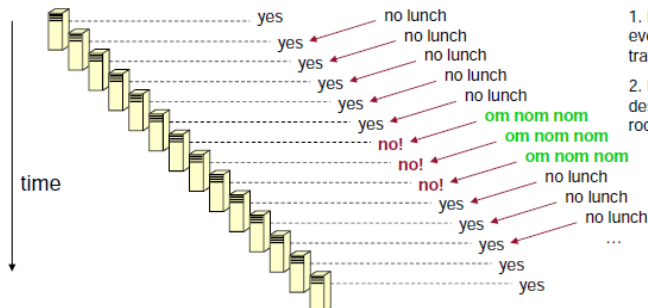
Information from a **specific** observation about the **specific** next value.



Information transfer

Transfer entropy measures directed coupling between time-series.
Intuitively: the amount of information that a source variable tells us about a destination, in the context of the destination's current state.

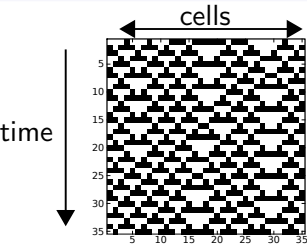
e.g. How much does knowing the IT guy is at lunch tell us about the web server running, given its previous state?



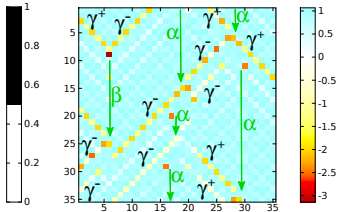
1. Directionally dependent events have max info transfer

2. Internal predictability in destination allows no room for transfer

Information dynamics in CAs

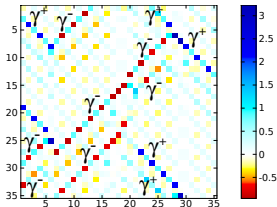


(a) Raw CA

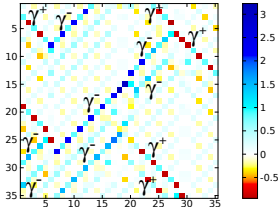


(b) LAIS

Domains and blinkers are the **dominant information storage** entities.



(c) LTE right

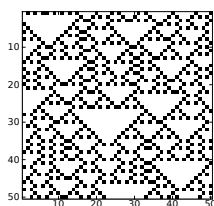


(d) LTE left

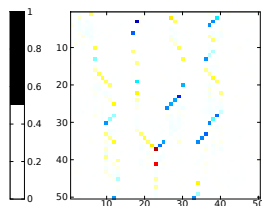
Gliders are the **dominant information transfer** entities.

Other transfer entropy characteristics

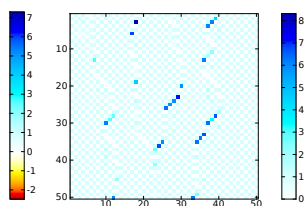
TE can be made **conditional** $T_{Y \rightarrow X|Z} = I(\mathbf{Y}_n^{(l)}; X_{n+1} | \mathbf{X}_n^{(k)}, \mathbf{Z}_n^{(m)})$
 or **multivariate** $T_{Y \rightarrow X|Z} = I(\{\mathbf{Y}_n^{(l)}, \mathbf{Z}_n^{(m)}\}; X_{n+1} | \mathbf{X}_n^{(k)})$ (Lizier et al., 2008, 2010, 2011)



(a) Raw CA



(b) LTE left



(c) LTE left conditional

Lizier
et al.
(2008,
2014)

Computed over delay u as $T_{Y \rightarrow X|Z} = I(\mathbf{Y}_{n-u+1}^{(l)}; X_{n+1} | \mathbf{X}_n^{(k)})$
 (Wibral et al., 2013).

Discrete: plug-in estimator



For discrete variables x and y , to compute $H(X, Y)$

- 1 estimate: $p(x, y) = \frac{\text{count}(X=x, Y=y)}{N}$, where N is our sample size;
- 2 plug-in each estimated PDF to $H(X, Y)$ to get $\hat{H}(X, Y)$

Discrete: plug-in estimator



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Bias: expected offset of estimated value from a finite sample set from the true underlying value of the measure. There are several available bias correction techniques.

Variance: variance in estimated values of the measure (from a finite sample set) around the expected value.

Discrete: plug-in estimator



For discrete variables x and y , to compute $H(X, Y)$

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Bias: expected offset of estimated value from a finite sample set from the true underlying value of the measure. There are several available bias correction techniques.

Variance: variance in estimated values of the measure (from a finite sample set) around the expected value.

A simple way to handle continuous variables is to discretise or bin them.

Continuous variables → Differential entropy



Differential entropy:

$$H_D(X) = - \int_{S_X} f(x) \log f(x) dx$$

for PDF $f(x)$, where S_X is the set where $f(x) > 0$.

Evaluate all measures as sums and differences of $H_D(X)$ terms.

The properties of $H_D(X)$ are slightly odd ... however, the properties of $I_D(X; Y)$ are the same as for discrete variables.

JIDT includes 3 estimation methods for differential entropy based MIs (and conditional MIs) ...

Gaussian model



If a multivariate $\mathbf{X}$ (of d dimensions) is Gaussian distributed (Cover and Thomas, 1991):

$$H(\mathbf{X}) = \frac{1}{2} \ln \left[(2\pi e)^d |\Omega_{\mathbf{X}}| \right]$$

(in *nats*) where $|\Omega_{\mathbf{X}}|$ is the determinant of the $d \times d$ covariance matrix $\Omega_{\mathbf{X}} = \overline{\mathbf{X}\mathbf{X}^T}$.

Any measure is computed as sums and differences of these joint entropies.

Pros: fast ($O(Nd^2)$), parameter free

Cons: subject to the linear-model assumption

Kernel estimation

Estimate PDFs with a *kernel function* Θ , measuring “similarity” between pairs of samples $\{x_n, y_n\}$ and $\{x_{n'}, y_{n'}\}$ using a resolution or *kernel width* r .

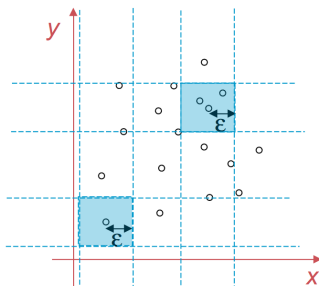
$$\text{E.g.:} \quad \hat{p}_r(x_n, y_n) = \frac{1}{N} \sum_{n'=1}^N \Theta \left(\left| \begin{pmatrix} x_n - x_{n'} \\ y_n - y_{n'} \end{pmatrix} \right| - r \right).$$

By default Θ is the step kernel ($\Theta(x > 0) = 0$, $\Theta(x \leq 0) = 1$), and the norm $|\cdot|$ is the maximum distance.

Pros: model-free (captures non-linearities)

Cons: sensitive to r , is biased, less time-efficient (though can be reduced to $O(N \log N)$).

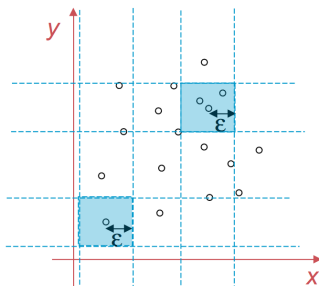
Estimating $p(x, y)$, $p(x)$ and $p(y)$



Kernel estimation

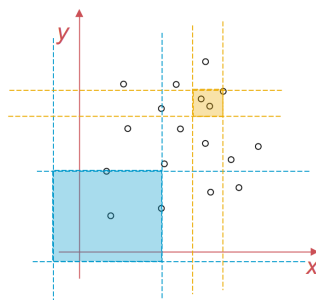
- Fixed width $r = \epsilon$
- MI: “How does knowing x within r help me predict y within r ?”

Estimating $p(x, y)$, $p(x)$ and $p(y)$



Kernel estimation

- Fixed width $r = \epsilon$
- MI: “How does knowing x within r help me predict y within r ?”



Kraskov (KSG) technique (Kraskov et al., 2004)

- Dynamic width r and bias correction
- MI: “How does knowing x within the K nearest neighbours in the joint space help me predict y ?”

- Harnessing Kozachenko-Leonenko entropy estimators;
- Using nearest-neighbour counting, with a fixed number K of neighbours in the full joint space.

KSG estimators (Kraskov et al., 2004)



Improve on box-kernel estimation with lower bias via:

- Harnessing Kozachenko-Leonenko entropy estimators;
- Using nearest-neighbour counting, with a fixed number K of neighbours in the full joint space.

There are two algorithms; algorithm 1 gives:

$$I^{(1)}(X; Y) = \psi(K) - \langle \psi(n_x + 1) + \psi(n_y + 1) \rangle + \psi(N),$$

(in *nats*) where ψ denotes the digamma function.

Extensions to conditional MI are available (Frenzel and Pompe, 2007; Gomez-Herrero et al., 2010; Wibral et al., 2014).

Pros: model-free, bias corrected, best of breed in terms of data efficiency and accuracy, and is effectively parameter free (w.r.t K).

Cons: less time-efficient (though fast nearest neighbour searching reduces this to $O(KN \log N)$).

Why JIDT?



JIDT is unique in the **combination** of features it provides:

- Large array of measures, including all conditional/multivariate forms of the transfer entropy, and complementary measures such as active information storage.
- Wide variety of estimator types and applicability to both discrete and continuous data

Measure-estimator combinations

As of V1.2 distribution:

Measure		Discrete estimator	Continuous estimators			
Name	Notation		Gaussian	Box-Kernel	Kraskov <i>et al.</i> (KSG)	Permutation
Entropy	$H(X)$	✓	✓	✓	*	
Entropy rate	$H_{\mu X}$	✓	Use two multivariate entropy calculators			
Mutual information (MI)	$I(X; Y)$	✓	✓	✓	✓	
Conditional MI	$I(X; Y Z)$	✓	✓		✓	
Multi-information	$I(\mathbf{X})$	✓		✓ ^U	✓ ^U	
Transfer entropy (TE)	$T_{Y \rightarrow X}$	✓	✓	✓	✓	✓ ^U
Conditional TE	$T_{Y \rightarrow X Z}$	✓	✓ ^U		✓ ^U	
Active information storage	A_X	✓	✓ ^U	✓ ^U	✓ ^U	
Predictive information	E_X	✓	✓ ^U	✓ ^U	✓ ^U	
Separable information	S_X	✓				

Why JIDT?



JIDT is unique in the **combination** of features it provides:

- Large array of measures, including all conditional/multivariate forms of the transfer entropy, and complementary measures such as active information storage.
- Wide variety of estimator types and applicability to both discrete and continuous data
- Local measurement for all estimators;
- Statistical significance calculations for MI, TE;
- No dependencies on other installations (except Java);
- Lots of demos and information on website/wiki:
 - <https://code.google.com/p/information-dynamics-toolkit/>

Why implement in Java?



The Java implementation of JIDT gives us several fundamental features:

- Platform agnostic, requiring only a JVM;
- Object-oriented code, with a hierarchical design to interfaces for each measure, allowing dynamic swapping of estimators for the same measure;
- JIDT can be directly called from Matlab/Octave, Python, R, Julia, Clojure, etc, adding efficiency for higher level code;
- Automatic generation of Javadocs.

Installation



- 1 Download the latest full distribution by following the Download link at <https://lazier.me/jidt/> (links to google code or github after JIDT moves there ...)
- 2 Unzip it to your preferred location for the distribution
- 3 To be able to use it, you will need the `infodynamics.jar` on your classpath.

Installation



- 1 Download the latest full distribution by following the Download link at <https://lizier.me/jidt/> (links to google code or github after JIDT moves there ...)
- 2 Unzip it to your preferred location for the distribution
- 3 To be able to use it, you will need the `infodynamics.jar` on your classpath.

That's it!

Installation – caveats



- ① You'll need a JRE installed (Version ≥ 6)
 - ① Should come automatically with Matlab/Octave/Python-JPy installation
- ② You need `ant` if you want to rebuild the project using `build.xml`
- ③ You need `junit` if you want to run the unit tests
- ④ Additional preparation may be required to use JIDT in GNU Octave or Python ...

Check that your environment works



Java:

- 1 Run `demos/java/example1TeBinaryData.sh` or `.bat`

Matlab/Octave:

- 1 For Octave version < 3.8 , first follow steps on the [wiki](#), including installing `octave-java` from `octave-forge`.
- 2 Run `demos/octave/example1TeBinaryData.m`

Python:

- 1 Install `jPy` to connect Python to Java (or `jpype1-py3` for Python 3)
- 2 Run `demos/python/example1TeBinaryData.py`

In case of issues, see the wiki pages on Non-Java environments or the Instructor.

Contents of distribution



- `license-gplv3.txt` - GNU GPL v3 license;
- `infodynamics.jar` library file;
- **Documentation**
- **Source code** in `java/source` folder
- Unit tests in `java/unittests` folder
- `build.xml` ant build script
- **Demonstrations** of the code in `demos` folder.

Documentation



Included in the distribution:

- `readme.txt`;
- `InfoDynamicsToolkit.pdf` – a pre-print of the publication introducing JIDT;
- `tutorial` folder – a full tutorial presentation and sample exercise (also via [JIDT wiki](#))
- `javadocs` folder – documents the methods and various options for each estimator class;
- PDFs describing each demo in the `demos` folder;

Also see:

- The wiki pages on the [JIDT website](#)
- Our email discussion list `jidt-discuss` on [Google groups](#).

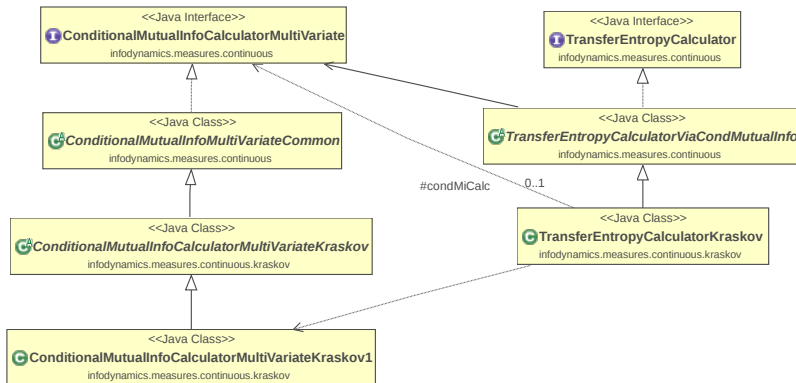
Source code structure



Source code at `java/source` is organised into the following Java *packages* (mapping directly to subdirectories):

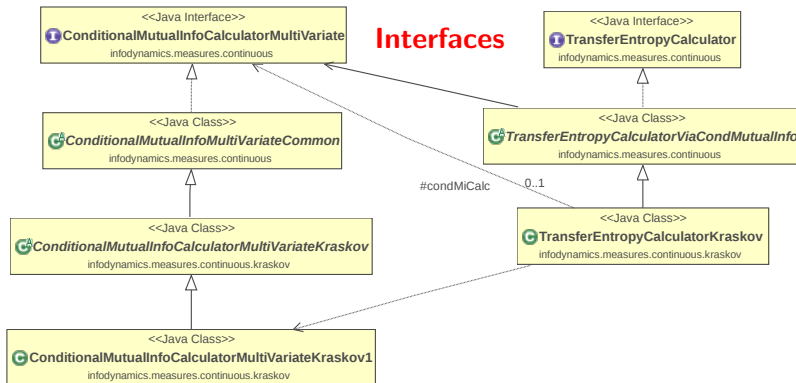
- `infodynamics.measures`
 - `infodynamics.measures.discrete` – for discrete data;
 - `infodynamics.measures.continuous` – for continuous data
 - top level: Java *interfaces* for each measure, then
 - a set of sub-packages (gaussian, kernel, kozachenko, kraskov and symbolic) containing *implementations* of such estimators for these interfaces.
 - `infodynamics.measures.mixed` – experimental discrete-to-continuous MI calculators
- `infodynamics.utils` – utility functions
- `infodynamics.networkinference` – higher-level algorithms

Architecture for calculators on continuous data



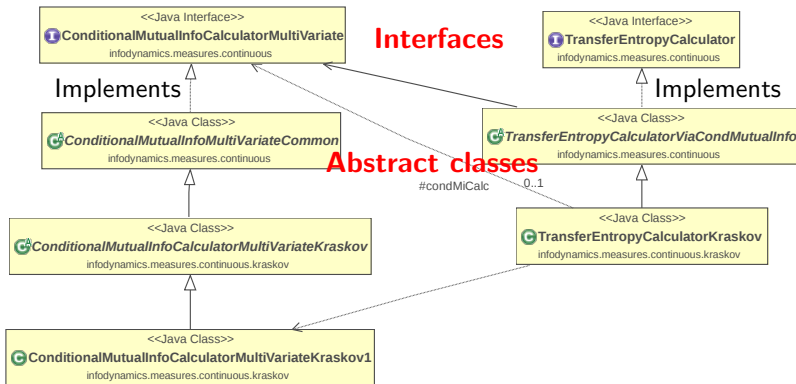
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Architecture for calculators on continuous data



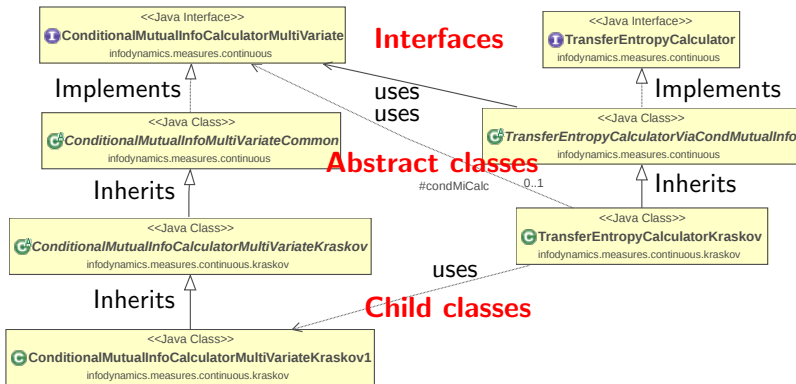
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Architecture for calculators on continuous data



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Architecture for calculators on continuous data



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Demos



JIDT is distributed with the following demos:

- **Auto-analyser GUI** (code generator)
- Simple Java Demos
 - Mirrored in Matlab/Octave, Python, R, Julia, Clojure.
- Recreation of Schreiber's original transfer entropy examples;
- Information dynamics in Cellular Automata;
- Detecting interaction lags;
- Interregional coupling;
- Behaviour of null/surrogate distributions;

All have documentation (PDF and wiki pages) provided to help run them.

Auto Analyser GUI (code generator)



A GUI application to:

- Make simple MI and TE calculations for you;
- Create code for you.

While this auto-analysis is only provided for MI and TE, note that the general coding paradigm for all calculators is the same.

Scripts to start the apps are in the demos/AutoAnalyser folder:

- `runTEAutoAnalyser.sh` (and `.bat`) for TE, and
- `runMIAutoAnalyser.sh` (and `.bat`) for MI.

Run: `runTEAutoAnalyser.sh` (or `.bat`)

Auto Analyser GUI (code generator)



Computing TE and MI could not be easier

JIDT Transfer Entropy Auto-Analyser

Calculation parameters

Calculator Type: **Discrete**

Data file: **hDynamics/demos/data/2coupledBinaryColsUsek2.txt**

Select

Valid data file with 1000 rows and 2 columns

Source column: **0**

Destination column: **1**

Property name	Property value
k_HISTORY	1
base	2

Compute

Generated code

Java Python Matlab

... Awaiting new parameter selection (press compute) ...

JIDT

Auto Analyser GUI (code generator)



Computing TE and MI could not be easier

Calculation parameters

Calculator Type: **Discrete**

Data file: **hDynamics/demos/data/2coupledBinaryColsUsek2.txt**
Select
Valid data file with 1000 rows and 2 columns

Source column: **0**

Destination column: **1**

Property name	Property value
k_HISTORY	1
base	2

Compute

Generated code

Java Python Matlab

... Awaiting new parameter selection (press compute) ...

Just follow the GUI:

- 1 Select estimator
- 2 Select data file
- 3 Identify source/target columns in data
- 4 Fill out properties (use tool tip for descriptions)
- 5 Click "Compute"

Auto Analyser GUI (code generator)



Clicking "compute" then gives you:

- 1 The resulting TE and MI

The screenshot displays the JIDT Transfer Entropy Auto-Analyser GUI. The left panel, titled "Calculation parameters", shows the "Calculator Type" set to "Discrete". The "Data file" is "Dynamics/demos/data/2coupledBinaryColsUseK2.txt", with a "Select" button and a note "Valid data file with 1000 rows and 2 columns". The "Source column" is 0 and the "Destination column" is 1. Below this is a table of properties:

Property name	Property value
k_HISTORY	1
base	2

At the bottom of the left panel is a "Compute" button and the output text "TE_Discrete(col_0 -> col_1) = 0.0003 bits". The right panel, titled "Generated code", shows the generated Java code for the calculation.

```
package infodynamics.demos.autoanalysis;

import infodynamics.utils.ArrayFileReader;
import infodynamics.utils.MatrixUtils;

import infodynamics.measures.discrete.*;

public class GeneratedCalculator {

    public static void main(String[] args) throws Exception {

        // 0. Load/prepare the data:
        String dataFile = "/home/joseph/Dropbox/Work/Investigations/JavaCode/sharedProje
        ArrayFileReader afr = new ArrayFileReader(dataFile);
        int[][] data = afr.getInt2DMatrix();
        int[] source = MatrixUtils.selectColumn(data, 0);
        int[] dest = MatrixUtils.selectColumn(data, 1);

        // 1. Construct the calculator:
        TransferEntropyCalculatorDiscrete calc
            = new TransferEntropyCalculatorDiscrete(2, 1);
        // 2. No other properties to set for discrete calculators.
        // 3. Initialise the calculator for (re-)use:
        calc.initialise();
        // 4. Supply the sample data:
        calc.addObservations(source, dest);
        // 5. Compute the estimate:
        double result = calc.computeAverageLocalOfObservations();
        System.out.printf("TE_Discrete(col_0 -> col_1) = %.4f bits\n", result);
    }
}
```

Auto Analyser GUI (code generator)

JIDT

Clicking "compute" then gives you:

- 1 The resulting TE and MI, **and**
- 2 Code to generate this calculation in Java, Python and Matlab

JIDT Transfer Entropy Auto-Analyser

Calculation parameters

Calculator Type: **Discrete**

Data file: **Dynamics/demos/data/2coupledBinaryColsUseK2.txt**
 Select
 Valid data file with 1000 rows and 2 columns

Source column: **0**
 Destination column: **1**

Property name	Property value
k_HISTORY	1
base	2

Compute

TE_Discrete(col_0 -> col_1) = 0.0003 bits

Generated code

Java Python Matlab

```
package infodynamics.demos.autoanalysis;

import infodynamics.utils.ArrayFileReader;
import infodynamics.utils.MatrixUtils;

import infodynamics.measures.discrete.*;

public class GeneratedCalculator {

    public static void main(String[] args) throws Exception {

        // 0. Load/prepare the data:
        String dataFile = "/home/joseph/Dropbox/Work/Investigations/JavaCode/sharedProje
        ArrayFileReader afr = new ArrayFileReader(dataFile);
        int[][] data = afr.getInt2DMatrix();
        int[] source = MatrixUtils.selectColumn(data, 0);
        int[] dest = MatrixUtils.selectColumn(data, 1);

        // 1. Construct the calculator:
        TransferEntropyCalculatorDiscrete calc
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        // 2. No other properties to set for discrete calculators.
        // 3. Initialise the calculator for (re-)use:
        calc.initialise();
        // 4. Supply the sample data:
        calc.addObservations(source, dest);
        // 5. Compute the estimate:
        double result = calc.computeAverageLocalOfObservations();
        System.out.printf("TE_Discrete(col_0 -> col_1) = %.4f bits\n", result);
    }
}
```

Auto Analyser GUI (code generator) – discrete



Let's generate a sample TE calculation on **discrete** data:

- ① Select *Discrete* estimator
- ② Select the data file `2coupledBinaryColsUseK2.txt` from the default directory in the *Select* popup.
 - ① Note: the GUI checks validity of the file
 - ② Valid file format is described when you hover on *Data file*
- ③ Leave source and destination columns, and properties, for now
- ④ Click *Compute*

Did everyone get the result 0.0003 bits?

Auto Analyser GUI (code generator) – discrete



Let's generate a sample TE calculation on **discrete** data:

- ① Select *Discrete* estimator
- ② Select the data file `2coupledBinaryColsUseK2.txt` from the default directory in the *Select* popup.
 - ① Note: the GUI checks validity of the file
 - ② Valid file format is described when you hover on *Data file*
- ③ Leave source and destination columns, and properties, for now
- ④ Click *Compute*

Did everyone get the result 0.0003 bits?

What if you change the property `k_HISTORY` to 2?

Hover on the property names to see the description for them

Auto Analyser GUI – discrete – code analysis



Let's examine the code that was generated.

Either:

- ❶ Click on the panel for the language you want to work in, OR
- ❷ Open the file that was automatically generated for you:
 - ❶ `demos/java/infodynamics/demos/autoanalysis/GeneratedCalculator.java`
OR
 - ❷ `demos/AutoAnalyser/GeneratedCalculator.m` OR
 - ❸ `demos/AutoAnalyser/GeneratedCalculator.py`

Auto Analyser GUI – discrete – code analysis



Observe how the classpath is pointed to `infodynamics.jar`:

- **Java**: java command line in
demos/AutoAnalyser/runAutoGenerated.sh/.bat (or in IDE);
- **Matlab/Octave**: `javaaddpath()` statement;
- **Python**: `startJVM()` statement.

Auto Analyser GUI – discrete – code analysis



Observe how the classpath is pointed to `infodynamics.jar`:

- **Java**: java command line in `demos/AutoAnalyser/runAutoGenerated.sh/.bat` (or in IDE);
- **Matlab/Octave**: `javaaddpath()` statement;
- **Python**: `startJVM()` statement.

Note: Discrete data (source and dest) represented as `int[]` arrays:

- 1 with values in the range $0 \dots base - 1$, where e.g. `base=2` for binary.
- 2 for time-series measures, the array is indexed by time.
- 3 for multivariate time-series, we use `int[][]` arrays, indexed first by time then variable number.

Auto Analyser GUI – discrete – Usage Paradigm

```
1 // int[] source and dest defined and loaded earlier
2 TransferEntropyCalculatorDiscrete calc = new TransferEntropyCalculatorDiscrete
   (2, 1);
3 calc.initialise();
4 calc.addObservations(source, dest);
5 double result = calc.computeAverageLocalOfObservations();
```

- ① **Construct** the calculator, providing parameters
 - ① Always check Javadocs for which parameters are required.
 - ② Here the parameters are the number of possible discrete symbols per sample (2, binary), and history length for TE ($k = 1$).
 - ③ Constructor syntax is different for Matlab/Octave/Python.

Auto Analyser GUI – discrete – Usage Paradigm

```
1 // int[] source and dest defined and loaded earlier
2 TransferEntropyCalculatorDiscrete calc = new TransferEntropyCalculatorDiscrete
   (2, 1);
3 calc.initialise();
4 calc.addObservations(source, dest);
5 double result = calc.computeAverageLocalOfObservations();
```

② **Initialise** the calculator prior to:

- ① use, or
- ② re-use (e.g. looping back from line 5 back to line 3 to examine different data).
- ③ This clears PDFs ready for new samples.

Auto Analyser GUI – discrete – Usage Paradigm

```
1 // int[] source and dest defined and loaded earlier
2 TransferEntropyCalculatorDiscrete calc = new TransferEntropyCalculatorDiscrete
   (2, 1);
3 calc.initialise();
4 calc.addObservations(source, dest);
5 double result = calc.computeAverageLocalOfObservations();
```

3 Supply the data to the calculator to construct PDFs:

- ① addObservations() may be called multiple times;
- ② Convert arrays into Java format:
 - From Matlab/Octave using our octaveToJavaIntArray(array), etc., scripts.
 - From Python using JArray(JInt, numDims)(array) for conversion or use numpy arrays directly, etc.

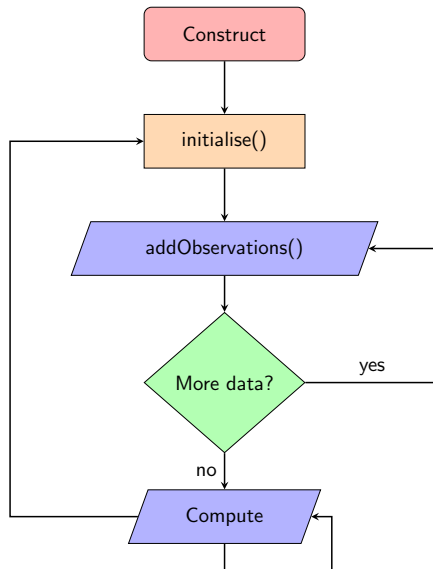
Auto Analyser GUI – discrete – Usage Paradigm

```
1 // int[] source and dest defined and loaded earlier
2 TransferEntropyCalculatorDiscrete calc = new TransferEntropyCalculatorDiscrete
   (2, 1);
3 calc.initialise();
4 calc.addObservations(source, dest);
5 double result = calc.computeAverageLocalOfObservations();
```

④ Compute the measure:

- ① Value is always returned in bits for discrete calculators.
- ② Result here approaches 0 bits for $k = 1$ (as written above), or 1 bit for $k = 2$, as destination is an XOR between the (random) source and it's own value from 2 time steps in the past.
- ③ Other computations include:
 - ① computeLocalOfPreviousObservations() for local values
 - ② computeSignificance() to compute p -values of measures of predictability (see Appendix A5 of paper for description).

Discrete Data – Usage Paradigm



Auto Analyser GUI (code generator) – continuous



Next we generate a sample TE calculation on **continuous** data:

- ❶ Select *Kraskov (KSG)* estimator
- ❷ Select the data file `2coupledRandomCols-1.txt` from the default directory in the *Select* popup.
- ❸ Leave source and destination columns
- ❹ Set the property `k_HISTORY` to 2
 - ❶ You can hover on the properties to see their meaning and valid values
- ❺ Click *Compute*

Auto Analyser GUI (code generator) – continuous



Next we generate a sample TE calculation on **continuous** data:

- ① Select *Kraskov (KSG)* estimator
- ② Select the data file `2coupledRandomCols-1.txt` from the default directory in the *Select* popup.
- ③ Leave source and destination columns
- ④ Set the property `k_HISTORY` to 2
 - ① You can hover on the properties to see their meaning and valid values
- ⑤ Click *Compute*

Did everyone get the result 0.2906 nats?

Auto Analyser GUI – discrete – code analysis



Let's examine the code that was generated.

Can you notice anything different to the discrete case?

Auto Analyser GUI – discrete – code analysis



Let's examine the code that was generated.

Can you notice anything different to the discrete case?

Note: Continuous data (source and dest) represented as `double[]` arrays:

- 1 for time-series measures, the array is indexed by time.
- 2 for multivariate time-series, we use `double[] []` arrays, indexed first by time then variable number.

Auto Analyser GUI – continuous – Usage Paradigm



```
1 // double[] source and dest defined and loaded earlier
2 TransferEntropyCalculator calc = new TransferEntropyCalculatorKraskov();
3 calc.setProperty(TransferEntropyCalculator.K_PROP_NAME, "2");
4 calc.initialise();
5 calc.setObservations(source, dest);
6 double result = calc.computeAverageLocalOfObservations();
```

- ① **Construct** the calculator, *possibly* providing parameters
 - ① Always check Javadocs for which parameters are required.
 - ② For continuous calculators, parameters may always be provided later (see next slide) to allow dynamic instantiation.
 - ③ Constructor syntax is different for Matlab/Octave/Python.

Auto Analyser GUI – continuous – Usage Paradigm

```
1 // double[] source and dest defined and loaded earlier
2 TransferEntropyCalculator calc = new TransferEntropyCalculatorKraskov();
3 calc.setProperty(TransferEntropyCalculator.K_PROP_NAME, "2");
4 calc.initialise();
5 calc.setObservations(source, dest);
6 double result = calc.computeAverageLocalOfObservations();
```

② Set properties for the calculator (new method for continuous):

- ① Formally: Check the Javadocs for available properties for each calculator;
 - ① In Auto Analyser GUI, you can hover on each parameter to read about it.
- ② E.g. here we set the embedding (history) length $k = 2$ for the TE calculation.
- ③ Property names and values are always key-value pairs of String objects;
- ④ Only guaranteed to hold after the next `initialise()` call.
- ⑤ Properties can easily be extracted and set from a file (see Simple Demo 6).

Auto Analyser GUI – continuous – Usage Paradigm



```
1 // double[] source and dest defined and loaded earlier
2 TransferEntropyCalculator calc = new TransferEntropyCalculatorKraskov();
3 calc.setProperty(TransferEntropyCalculator.K_PROP_NAME, "2");
4 calc.initialise();
5 calc.setObservations(source, dest);
6 double result = calc.computeAverageLocalOfObservations();
```

③ Initialise the calculator prior to:

- ① use or re-use, as for Discrete.
- ② This clears PDFs ready for new samples, and finalises any new property settings.
- ③ There may be several overloaded forms taking different properties in directly. E.g. for TE, `calc.initialise(1)` sets history length $k = 1$. We could also call `calc.initialise(k, tau_k, 1, 1_tau, delay)` to specify embedding dimensions and source-target delay. Check Javadocs for options here.

Auto Analyser GUI – continuous – Usage Paradigm

```
1 // double[] source and dest defined and loaded earlier
2 TransferEntropyCalculator calc = new TransferEntropyCalculatorKraskov();
3 calc.setProperty(TransferEntropyCalculator.K_PROP_NAME, "2");
4 calc.initialise();
5 calc.setObservations(source, dest);
6 double result = calc.computeAverageLocalOfObservations();
```

- ④ Supply the data to the calculator to construct PDFs:
 - ① setObservations() may be called once, OR
 - ② call addObservations() multiple times, in between startAddObservations() and finaliseAddObservations() calls.
 - ③ Convert arrays into Java format:
 - From Matlab/Octave using our octaveToJavaDoubleArray(array), etc., scripts.
 - From Python using JArray(JDouble, numDims)(array) for conversion or use numpy arrays directly, etc.

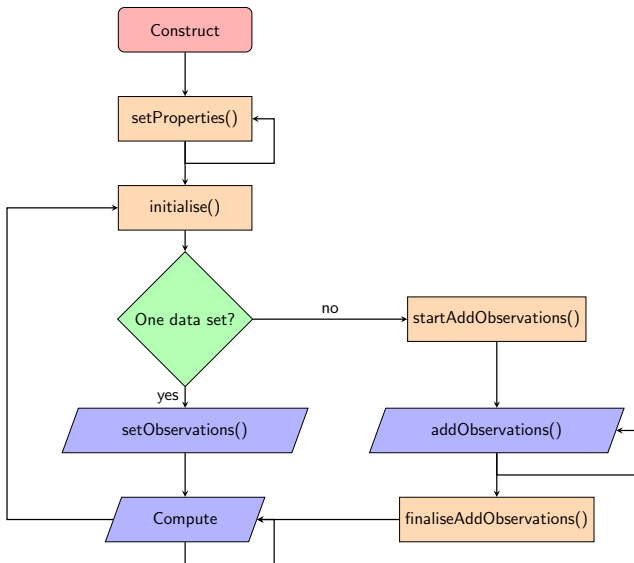
Auto Analyser GUI – continuous – Usage Paradigm

```
1 // double[] source and dest defined and loaded earlier
2 TransferEntropyCalculator calc = new TransferEntropyCalculatorKraskov();
3 calc.setProperty(TransferEntropyCalculator.K_PROP_NAME, "2");
4 calc.initialise();
5 calc.setObservations(source, dest);
6 double result = calc.computeAverageLocalOfObservations();
```

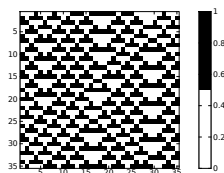
5 Compute the measure:

- ① Value may be in *bits* (kernel) OR in *nats* (KSG, Gaussian) calculators.
- ② Result here is 0.2906 nats since destination is correlated with the (random) source.
- ③ Other computations include:
 - ① `computeLocalOfPreviousObservations()` for local values
 - ② `computeSignificance()` to compute p -values of measures of predictability (see Appendix A5 of paper for description).

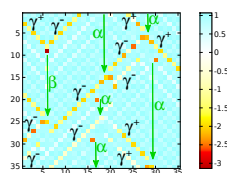
Continuous Data – Usage Paradigm



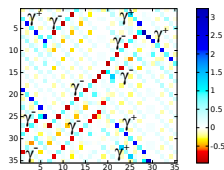
Many other demos – e.g. local dynamics in CAs



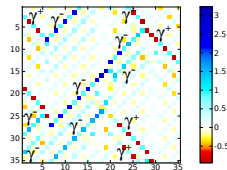
(a) Raw CA



(b) LAIS



(c) LTE right



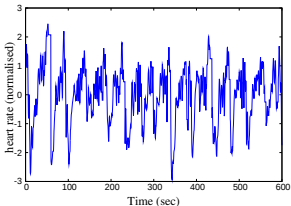
(d) LTE left

See PDF documentation for `demos/octave/CellularAutomata/` to recreate, e.g. run `GsoChapterDemo2013.m`.

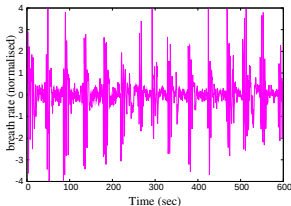
Exercise

- 1 Let's consider the common information between heart rate & breath rate measurements (data from Rigney et al. (1993) in `demos/data/SFI-heartRate_breathVol1_blood0x.txt`).

Open the data file to take a look:



Heart rate



Breath rate

Exercise



- ② Task – on `demos/data/SFI-heartRate_breathVol_blood0x.txt`:
 - ① Compute Mutual Information between the heart and breath rate time-series data in that example (samples 2350 to 3550, inclusive),
 - ② Using a Kraskov (KSG) estimator, algorithm 2, with 4 nearest neighbours.
- ③ How to approach it?

Exercise



- ② Task – on `demos/data/SFI-heartRate_breathVol_blood0x.txt`:
 - ① Compute Mutual Information between the heart and breath rate time-series data in that example (samples 2350 to 3550, inclusive),
 - ② Using a Kraskov (KSG) estimator, algorithm 2, with 4 nearest neighbours.
- ③ How to approach it?
 - ① **Easy way**: use the **MI** Auto Analyser GUI to generate code, then you can alter which samples it is using.

Exercise



- ② Task – on `demos/data/SFI-heartRate_breathVol_blood0x.txt`:
 - ① Compute Mutual Information between the heart and breath rate time-series data in that example (samples 2350 to 3550, inclusive),
 - ② Using a Kraskov (KSG) estimator, algorithm 2, with 4 nearest neighbours.
- ③ How to approach it?
 - ① **Easy way**: use the **MI** Auto Analyser GUI to generate code, then you can alter which samples it is using.
 - ② **More challenging**: Start from KSG TE demo on this data set, in your preferred environment, and modify it to compute MI:
 - ① `demos/java/infodynamics/java/schreiberTransferEntropyExamples/-HeartBreathRateKraskovRunner.java`
 - ② `demos/octave/SchreiberTransferEntropyExamples/runHeartBreathRateKraskov.m`
 - ③ `demos/python/SchreiberTransferEntropyExamples/runHeartBreathRateKraskov.py`
 - ④ **Hint 1**: A KSG MI calculator is used in Simple Demo 6.
 - ⑤ **Hint 2**: Remember the usage paradigm on slide 62.

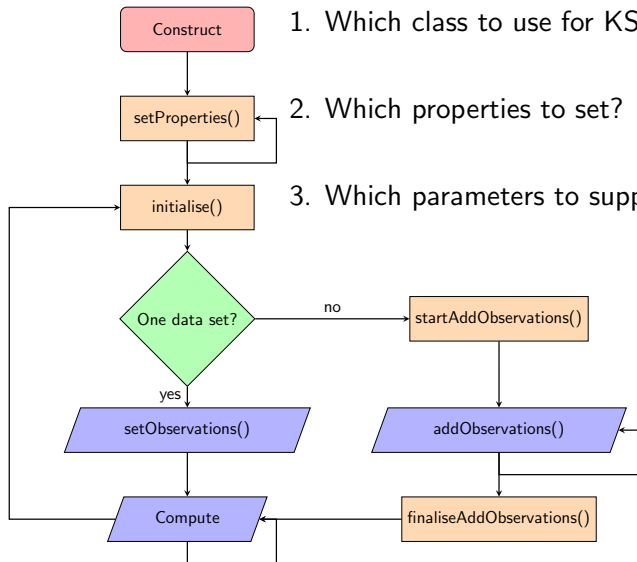
Exercise



- ② Task – on `demos/data/SFI-heartRate_breathVol_blood0x.txt`:
 - ① Compute Mutual Information between the heart and breath rate time-series data in that example (samples 2350 to 3550, inclusive),
 - ② Using a Kraskov (KSG) estimator, algorithm 2, with 4 nearest neighbours.
- ③ How to approach it?
 - ① **Easy way**: use the **MI** Auto Analyser GUI to generate code, then you can alter which samples it is using.
 - ② **More challenging**: Start from KSG TE demo on this data set, in your preferred environment, and modify it to compute MI:
 - ① `demos/java/infodynamics/java/schreiberTransferEntropyExamples/-HeartBreathRateKraskovRunner.java`
 - ② `demos/octave/SchreiberTransferEntropyExamples/runHeartBreathRateKraskov.m`
 - ③ `demos/python/SchreiberTransferEntropyExamples/runHeartBreathRateKraskov.py`
 - ④ **Hint 1**: A KSG MI calculator is used in Simple Demo 6.
 - ⑤ **Hint 2**: Remember the usage paradigm on slide 62.

Answer: **0.134 ± 0.002 nats**; if you use all data: **0.0994 ± 0.0005**.

Exercise – Remember the usage paradigm!



1. Which class to use for KSG MI algorithm 2?

2. Which properties to set? Check Javadocs!

3. Which parameters to supply? Check Javadocs!

4. Supply one set of data

5. Calculate!

Exercise – sample answer

```
1 import infodynamics.measures.continuous.kraskov.  
    MutualInfoCalculatorMultiVariateKraskov2;  
2 // ...  
3 // New KSG MI (algorithm 2) calculator:  
4 miCalc = new  
    MutualInfoCalculatorMultiVariateKraskov2();  
5 miCalc.initialise(1,1); // univariate  
    calculation  
6 miCalc.setProperty("k", "4"); // 4 nearest  
    neighbours  
7 miCalc.setObservations(heart, chestVol);  
8 double miHeartToBreath = miCalc.  
    computeAverageLocalOfObservations();
```

Debrief



How did you find the exercise?

Was it difficult? Which parts?

Did you know where to find the information you needed?

Any questions arising from the exercise?

Exercise – Challenge task 1



Extend to compute $MI(\text{heart}; \text{breath})$, for a variety of **lags** between the two time-series. E.g., investigate lags of $[0, 1, \dots, 14, 15]$.

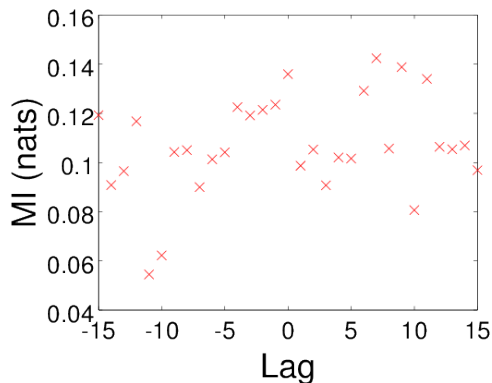
HINT: You could either:

- shift the time-series with respect to each other to affect the lag, or
- (cleaner) check out the available properties for this MI calculator in the Auto Analyser GUI or the Javadocs.
(Challenge: what to do if you want to use negative lags, i.e. a positive lag from breath to heart, with this property?)

Exercise – Challenge task 1



Extend to compute $MI(\text{heart}; \text{breath})$, for a variety of **lags** between the two time-series. E.g., investigate lags of $[0, 1, \dots, 14, 15]$.



What would you interpret from the results? Can you think of some logical further investigations here?

Summary



What I wanted you to take away today:

- Understand measures of information dynamics;
- Be able to obtain and install JIDT distribution;
- Be able to use the GUI Auto Analyser;
- Understand the code it generates in your chosen environment;
- Be able to modify sample scripts for new analysis;
- Know how and where to seek support information (wiki, Javadocs, mailing list, twitter).

Summary



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Did we get there?

Summary



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Did we get there?

Did you get what you came here for?

Summary



What I wanted you to take away today:

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- Know how and where to seek support information (wiki, Javadocs, mailing list, twitter).

Did we get there?

Did you get what you came here for?

Any other questions?

Final messages

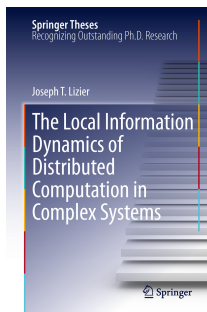


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Final messages



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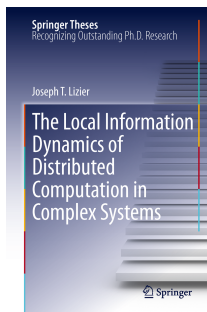


"The local information dynamics of distributed computation in complex systems", J. T. Lizier, Springer, Berlin (2013).

Final messages

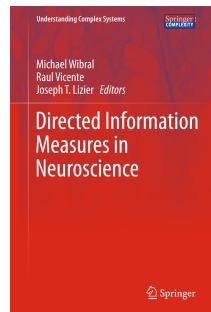


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“The local information dynamics of distributed computation in complex systems”, J. T. Lizier, Springer, Berlin (2013).

“Directed information measures in neuroscience”, edited by M. Wibral, R. Vicente and J. T. Lizier, Springer, Berlin (2014).



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