

The random walkers toolbox for analyzing single-particle tracking data – Supplement

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I. LIST OF ROUTINES IN THE TOOLBOX

Matlab routines and data files are provided on GitHub ([link](#)). For consistency, all Matlab routines should be moved into a subfolder `routines`, data files should be moved into a subfolder `routines/data`). Routines have been tested extensively with Matlab R2018b on MacOSX, R2020b on Linux, and with the comparable open-source clone Octave 6.2.0 on Linux. Full functionality was seen for both Matlab versions. For Octave, the function `wfbm` is not available, i.e. the creation of fractional Brownian motion tracks will require an auxiliary function. Also, in Octave the subroutines `plotter` and `oplot` in the program \rightarrow `driver.m` (with which all figures have been prepared) will have to be placed before the rest of the code (instead at the end), and graphics handles for changing fonts may have to be amended. Generating histograms will require a replacement of the command `histcounts` to the more basic command `histc`. Besides these points, all analysis routines seemed to work also in Octave.

For working with the evaluation routines, the program \rightarrow `driver.m` may serve as an initial template for designing the data-analysis workflow. This master program not only provides examples of how to call the individual evaluation routines but also how to create and store simulated trajectories with different properties by distinct data production routines (these parts are commented by a `%` sign and will have to be uncommented before use). Altogether, the following routines define the toolbox:

make_rndwalk.m	<code>pos = make_rndwalk(N,alpha,dx)</code>
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N	length of trajectories N
alpha	twofold Hurst coefficient
dx	mean step length (per dimension)
pos	$N \times 2$ -array of positions

This routine creates a two-dimensional FBM trajectory of length N with mean step size dx and Hurst coefficient $H = \alpha/2$.

make_switchwalk.m	<code>[sx,sy] = make_switchwalk(dt,N,M,xx,yy,k1,k2,fact)</code>
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dt	frame time Δt
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of FBM trajectories
yy	array of y-coordinates of FBM trajectories
k1	switch rate to low mobility
k2	switch rate to high mobility
fact	ratio of diffusion coefficients at high & low mobility ($f_K = K_{\text{high}}/K_{\text{low}} \geq 1$)

Based on an ensemble of FBM trajectories, this routine creates intermittent FBM trajectories that switch between a high- and a low-mobility state (ratio $f_K = K_{\text{high}}/K_{\text{low}} \geq 1$) with rates k_1 and k_2 , while maintaining the Hurst coefficient $H = \alpha/2$.

make_blurwalk.m `pos = make_blurwalk(N,alpha,dx,np)`

N	length of trajectories N
alpha	twofold Hurst coefficient
dx	mean step length (per dimension)
np	number of photons per substep (reasonable range: $n_p \in [10, 10^3]$)
pos	$N \times 2$ -array of positions

This routine creates a two-dimensional FBM trajectory of length N with mean step size dx and Hurst coefficient $H = \alpha/2$ including static and dynamic localization errors (tuned by the specified number of photons n_p).

ta_msd.m `[tau,msdt] = ta_msd(dt,xx,yy,dim,dis)`

dt	frame time Δt
xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
dim	dimension(s) for TA-MSD calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
tau	array of lag times
msd	array of TA-MSD values

This routine calculates the TA-MSD of a single trajectory in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

eata_msd.m `[tau,msdte] = eata_msd(dt,N,M,xx,yy,dim,dis)`

dt	frame time Δt
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of trajectories
yy	array of y-coordinates of trajectories
dim	dimension(s) for EA-TA-MSD calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
tau	array of lag times
msdte	array of EA-TA-MSD values

This routine calculates the EA-TA-MSD of an ensemble of M trajectories in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

ea_msd.m `[tau,msde] = ea_msd(dt,N,M,xx,yy,dim,dis)`

dt	frame time Δt
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of trajectories
yy	array of y-coordinates of trajectories
dim	dimension(s) for EA-MSD calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
tau	array of lag times
msde	array of EA-MSD values

This routine calculates the EA-MSD of an ensemble of M trajectories in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

eb.m `[tau,E] = eb(dt,N,M,xx,yy,dim,dis)`

dt	frame time Δt
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of trajectories
yy	array of y-coordinates of trajectories
dim	dimension(s) for EB calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spacing on linear/logarithmic scale
tau	array of lag times
E	array of ergodicity breaking parameters

This routine calculates the ergodicity breaking parameter of an ensemble of M trajectories in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

get_increm.m `[dx,dy] = get_increm(dn,xx,yy,chi)`

dn	lag in units of frame time, i.e. $\delta t / \Delta t$
xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
chi	true/false: normalized step increments (y/n)
dx	steps taken within dn frames in x
dy	steps taken within dn frames in y

This routine calculates the steps δx and δy taken within a period $\delta t = n\Delta t$ in a single trajectory, normalization is optional.

ta_quad.m `[tau,quat] = ta_quad(dt,xx,yy,dim,dis)`

dt	frame time Δt
xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
dim	dimension(s) for TA 4th-moment calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
tau	array of lag times
quat	array of TA 4th-moment values

This routine calculates the TA 4th-moment of a single trajectory in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

ta_gaussianity.m `[tau,g] = ta_gaussianity(dt,xx,yy,dim,dis)`

dt	frame time Δt
xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
dim	dimension(s) for TA-gaussianity calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
tau	array of lag times
g	array of TA-gaussianity values

This routine calculates the TA-gaussianity of a single trajectory in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

eata_gaussianity.m	[tau,g] = eata_gaussianity(dt,N,M,xx,yy,dim,dis)
dt	frame time Δt
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of trajectories
yy	array of y-coordinates of trajectories
dim	dimension(s) for EA-TA-gaussianity calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
tau	array of lag times
g	array of EA-TA-gaussianity values

This routine calculates the EA-TA-gaussianity of an ensemble of M trajectories in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

acf_sqinc.m	[tau,acf] = acf_sqinc(dt,dn,xx,yy,dim,dis)
dt	frame time Δt
dn	frame lag for steps, given by $\delta t/\Delta t$
xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
dim	dimension(s) for correlator calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
tau	array of lag times
acf	autocorrelation of squared increments

This routine calculates the autocorrelation of fluctuations of squared increments of a single trajectory in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

ea_acf_sqinc.m	[tau,acfe] = ea_acf_sqinc(dt,dn,N,M,xx,yy,dim,dis)
dt	frame time Δt
dn	frame lag for steps, given by $\delta t/\Delta t$
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of trajectories
yy	array of y-coordinates of trajectories
dim	dimension(s) for correlator calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
tau	array of lag times
acfe	ensemble-averaged autocorrelation of squared increments

This routine calculates the ensemble-averaged autocorrelation of fluctuations of squared increments in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

lch.m	[tau,Sd] = lch(dt,dn,xx,yy)
dt	frame time Δt
dn	number of positions to be used for local convex hull, typically 3-10
xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
tau	array of lag times
Sd	normalized maximum diameter of local convex hull, S_d

This routine calculates the maximum diameter of the LCH (based on dn points) of a single trajectory as a function of time. Values are normalized to the mean within the trajectory.

vacf.m `[xi,vacf] = vacf(dt,dn,xx,yy,dim,dis)`

dt	frame time Δt
dn	period for velocity, given by $\delta t / \Delta t$
xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
dim	dimension(s) for VACF calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
xi	array of rescaled time $\xi = \tau / \delta t$
vacf	array of VACF values

This routine calculates the VACF of a single trajectory in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

ea_vacf.m `[xi,vacfte] = ea_vacf(dt,dn,N,M,xx,yy,dim,dis)`

dt	frame time Δt
dn	period for velocity, given by $\delta t / \Delta t$
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of trajectories
yy	array of y-coordinates of trajectories
dim	dimension(s) for EA-VACF calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
xi	array of rescaled time $\xi = \tau / \delta t$
vacfte	array of EA-VACF values

This routine calculates the EA-VACF of an ensemble of M trajectories in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of lag times on a linear or logarithmic scale.

vacf_fbm_theo.m `[xi,vacf] = vacf_fbm_theo(alpha)`

alpha	scaling exponent (twofold Hurst coefficient)
xi	array of rescaled time $\xi = \tau / \delta t$
vacf	array of VACF values for FBM

This routine calculates the FBM prediction for the VACF.

psd.m `[f,psdt] = psd(dt,xx,yy,dim,dis)`

dt	frame time Δt
xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
dim	dimension(s) for PSD calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
f	array of frequencies
psdt	array of PSD values

This routine calculates the PSD of a single trajectory in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of frequencies on a linear or logarithmic scale.

ea_vacf.m **[f,psde] = ea_psd(dt,N,M,xx,yy,dim,dis)**

dt	frame time Δt
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of trajectories
yy	array of y-coordinates of trajectories
dim	dimension(s) for EA-PSD calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
f	array of frequencies
psde	array of EA-PSD values

This routine calculates the EA-PSD of an ensemble of M trajectories in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of frequencies on a linear or logarithmic scale.

cov_gamma.m **[fT,gam] = cov_gamma(dt,N,M,xx,yy,dim,dis)**

dt	frame time Δt
N	length of trajectories N
M	trajectory ensemble size M
xx	array of x-coordinates of trajectories
yy	array of y-coordinates of trajectories
dim	dimension(s) for calculation (0=xy, 1=x, 2=y)
dis	'lin'/'log': lag times equi-spaced on linear/logarithmic scale
fT	array of frequencies times total time, fT
gam	array of coeff. of variation values γ

This routine calculates the coefficient of variation γ of PSDs from an ensemble of M trajectories in two dimensions (xy) or along an individual coordinate (x or y), with an equi-distant spacing of dimensionless frequencies fT on a linear or logarithmic scale.

straightness.m **S = straightness(xx,yy,lb,ub)**

xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
lb	position at which straightness calculation starts
ub	position at which straightness calculation ends
S	straightness

This routine calculates the straightness of a trajectory between the specified time points.

asphericity.m **[An,Ad] = asphericity(xx,yy)**

xx	array of x-coordinates of single trajectory
yy	array of y-coordinates of single trajectory
An	nominator of asphericity A_s
Ad	denominator of asphericity A_s

This routine calculates nominator and denominator of the asphericity A_s of a single two-dimensional trajectory.

vacf_fbm_err.m **[xi,vacf] = vacf_fbm_err(dn,theta,alpha)**

dn	period for velocity, given by $\delta t / \Delta t$
theta	constant for static localization offset
alpha	scaling exponent (twofold Hurst coefficient)
xi	array of rescaled time $\xi = \tau / \delta t$
vacf	array of VACF values for FBM with localization errors

This routine calculates the FBM prediction for the VACF with localization errors.

II. LIST OF DATA FILES

For test purposes, several ensembles of two-dimensional fractional Brownian motion (FBM) based trajectories and an experimental data set are included in subfolder `routines/data`. All numerically obtained data sets, produced with the master program \rightarrow `driver.m`, consist of $M = 100$ trajectories, each with a length of $N = 500$ positions, using a time increment $\Delta t = 0.1$ s, and a (basic) average step size $\Delta x = 0.01 \mu\text{m}$.

data set 1

FBM tracks (created with `make_rndwalk.m`).

1. `FBM_xx_a_0.6.dat` & `FBM_yy_a_0.6.dat` ($\alpha = 2H = 0.6$, subdiffusion)
2. `FBM_xx_a_1.0.dat` & `FBM_yy_a_1.0.dat` ($\alpha = 2H = 1.0$, normal diffusion)
3. `FBM_xx_a_1.4.dat` & `FBM_yy_a_1.4.dat` ($\alpha = 2H = 1.4$, superdiffusion)

data set 2

Intermittent FBM tracks (created with `make_switchwalk.m`).

1. `FBM_switch_xx_a_0.6.dat` & `FBM_switch_yy_a_0.6.dat`
($\alpha = 2H = 0.6$, subdiffusion)
created from data set (1a) with $k_1 = 0.4/\text{s}$, $k_2 = 0.1/\text{s}$, and $f_K = 2$
2. `FBM_switch_xx_a_1.0.dat` & `FBM_switch_yy_a_1.0.dat`
($\alpha = 2H = 1.0$, normal diffusion)
created from data set (1b) with $k_1 = 0.2/\text{s}$, $k_2 = 0.3/\text{s}$, and $f_K = 3$

data set 3

FBM tracks with static localization error (created with `make_blurwalk.m`, $n_p = 50$).

1. `FBM_xx_a_0.6_np_50.dat` & `FBM_yy_a_0.6_np_50.dat` ($\alpha = 2H = 0.6$, subdiffusion)
2. `FBM_xx_a_1.0_np_50.dat` & `FBM_yy_a_1.0_np_50.dat` ($\alpha = 2H = 1.0$, normal diffusion)
3. `FBM_xx_a_1.4_np_50.dat` & `FBM_yy_a_1.4_np_50.dat` ($\alpha = 2H = 1.4$, superdiffusion)

data set 4

FBM tracks with dynamic localization error (created with `make_blurwalk.m`, $n_p = 900$).

1. `FBM_xx_a_0.6_np_900.dat` & `FBM_yy_a_0.6_np_900.dat` ($\alpha = 2H = 0.6$, subdiffusion)
2. `FBM_xx_a_1.0_np_900.dat` & `FBM_yy_a_1.0_np_900.dat` ($\alpha = 2H = 1.0$, normal diffusion)
3. `FBM_xx_a_1.4_np_900.dat` & `FBM_yy_a_1.4_np_900.dat` ($\alpha = 2H = 1.4$, superdiffusion)

data set 5

Ensemble of $M = 100$ experimentally obtained two-dimensional tracks of telomeres in the nucleus of untreated mammalian cells, each having a length $N = 2000$ and a frame time $\Delta t = 0.125$ s (part of the data set that has been analyzed and discussed in detail in (Krapf *et al.*, 2019; Stadler and Weiss, 2017)). Data have been shown to exhibit a transient FBM-like subdiffusion with $\alpha = 2H \approx 0.5$, followed by normal diffusion on larger time scales (Stadler and Weiss, 2017), in line with the motion of monomers in a Rouse polymer (for which such a subdiffusion is expected below the Rouse time). In addition, an excellent agreement of these data with key predictions for the power-spectral density of FBM trajectories has been found (Krapf *et al.*, 2019).

1. `telomers_xx.dat` & `telomers_yy.dat`

References

- Krapf, D., N. Lukat, E. Marinari, R. Metzler, G. Oshanin, C. Selhuber-Unkel, A. Squarcini, L. Stadler, M. Weiss, and X. Xu, 2019, *Phys Rev X* **9**, 011019.
 Stadler, L., and M. Weiss, 2017, *New J. Phys.* **19**, 113048.