Pi4U tutorial

What is Pi4U?

Pi4U is an extensible framework for non-intrusive Bayesian Uncertainty Quantification and Propagation of complex and computationally demanding physical models, that can exploit massively parallel computer architectures

What is implemented in Pi4U?

- TMCMC (for exact Bayesian inference)
- CMA-ES (for optimization)
- Subset Simulation (for rare event sampling)
- ABC-SubSim (for approximate Bayesian inference)
- A-PNDL (for parallel numerical differentiation)

Next steps in Pi4U?

- APIs for python, MATLAB, R
- Manifold TMCMC
- Hierarchical Bayesian (HB)
- Surrogate modelling
- EM algorithm for HB

TODO:

- Fix the composite prior and document it properly
- Document the Gaussian prior
- What priors do we support? Best option: a keyword for the prior for each parameter, then give prior parameters
- Give the users the possibility to make their own prior
- Add a check for the dimensionality vs number of parameters set by the user
- Provide an example with a fitfun in the script form
- Delete sn-TMCMC and everything else which is obsolete
- Delete with/without exp acceptance (leave only without)
- CMA-style proposal: make clear that it's 0 or 1
- Document interactive dumping more clearly
- Add CMA boundaries to the initials file
- Discuss if we should have MAP (as an option in the parameter file)

INSTALLATION on Mac:

Pi4U is based on TORC, a task-parallel library created by P. Hadjidoukas [1]. The steps to install TORC on a Mac are:

- 1. brew install mpich qsl R # skip if you already have those package
- 2. R # opens R. Inside R type the following:
 - a. install.packages("sp") # for plotting data
 - b. quit() # quits R
- 3. cd <path to your new pi4u installation>
- 5. cd pi4u/torc lite
- 6. ./configure CC=mpicc --prefix=<install folder> # default prefix:
 /usr/local
- 7. make && make install

INSTALLATION on a cluster

Instead of 1. load the corresponding modules

INSTALLATION on Linux (Ubuntu)

Follow the steps for Mac, just instead of step 1. Run:

sudo apt-get install libmpich-dev libgsl-dev r-base For the step "install.packages("sp")" in R and the step "make install" (in the case the default --prefix value was used in the ./configure step), you may need to use sudo.

EXAMPLE to run:

Sample negative Rosenbrock function (a = 1, b = 100) using TMCMC [2]:

- 1. cd pi4u.git/engines
- 2. make
- 3. mpirun -np 1 ./engine tmcmc
- 4. cd postprocessing tools
- 5. **set** fname **in** kdeplotter.R **to** "../curgen_db_008.txt" **which contains** samples of the last generation
- 6. Plot results: R --no-save < kdeplotter.R
- 7. cd ..
- 8. open curgen db 008.png #you should see Fig. 1

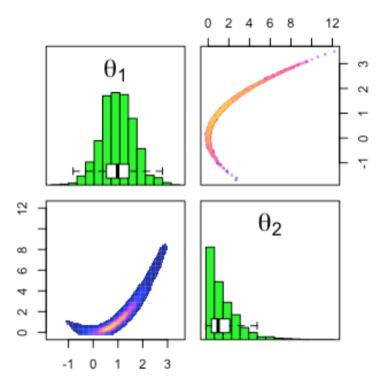


Fig. 1. Posterior PDF for the Rosenbrock function. Diagonals: marginal distributions of the corresponding parameters. Above diagonal: samples colored by the log-likelihood value.

Below diagonal: smoothed histogram of the samples.

PARAMETER file: tmcmc.par

The parameters in this file control the behavior of TMCMC. Be careful to change the prior parameters when changing dimension N of the problem, e.g. for gaussian prior the covariance matrix should be of dimension NxN (in matlab format) and the mean vector of dimension N.

SETTING UP your example

- 1. Write your fitness function:
 - a. for an easy-log-likelihood: modify engines/fitfun.c in such a way that it takes a vector of parameters and returns the log-likelihood value (see engines/fitfun.c for examples). The fitfun should have the following signature: double fitfun (double *x, int N, void *output, int *info), where x is the current sample, N is the dimensionality of the problem (denoted Nth in the tmcmc.par), output is used if the function should return something more than just a log-likelihood value, info contains information about the position of the current sample in the TMCMC task-graph: info[0] = stage, info[1] = chain, info[2] = task, info[3] can be used for subtask index.
 - b. for a complicated log-likelihood: write a script which will compute your log-likelihood value and call it from the <code>engines/fitfun.c</code>
- 2. Modify the tmcmc.par file accordingly: change the dimensionality of the problem (Nth), the population size (PopSize), the type of prior

Running other engines

CMA-ES:

CMA-ES is a black-box minimization algorithm [3].

- 1. mpirun -np 1 ./engine_cmaes # It uses the same fitfun.c and is compiled
 together with engine tmcmc
- 2. Have in mind that CMA-ES does minimization, so the log-likelihood has a minus sign inside <code>engine_cmaes.c</code>, because we look for the parameters which maximize the log-likelihood
- 3. Check the outcmaesxrecentbest.dat for the output
- 4. Now you can play with the parameters: edit <code>cmaes_initials.par</code> (see the comments inside)

Troubleshooting

- MPI: Make sure your localhost name is the same in your terminal and your /etc/hosts file. If not, you will get an error when trying to run mpich, "gethostbyname failed". To fix that: do cat /etc/hosts and it should print out sth like: // 127.0.0.1 my_hostname, where my_hostname should match the one in your terminal. If not, change one of them: change /etc/hosts: sudo vim /etc/hosts. Change terminal hostname: sudo scutil --set HostName my new host name

References

- [1] P.E. Hadjidoukas, E. Lappas, and V.V. Dimakopoulos. "A runtime library for platform-independent task parallelism", *In 20th Euromicro International Conference on Parallel, Distributed and Network-Based Processing (PDP)*, pages 229–236. IEEE Computer Society, Munich, Germany, 2012.
- [2] J. Ching and Y. Chen. "Transitional Markov chain Monte Carlo method for Bayesian model updating, model class selection, and model averaging." *J. Eng. Mech.*, 133(7):816–832, 2007.
- [3] Hansen, Nikolaus, Sibylle D. Müller, and Petros Koumoutsakos. "Reducing the time complexity of the derandomized evolution strategy with covariance matrix adaptation (CMA-ES)." *Evolutionary computation* 11.1 (2003): 1-18.
- [4] Hadjidoukas P.E., Angelikopoulos P., Kulakova L., Papadimitriou C., Koumoutsakos P. "Exploiting Task-Based Parallelism in Bayesian Uncertainty Quantification." *Lecture Notes in Computer Science* (2015): 9233, 532-544
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- [7] Hadjidoukas P.E, Angelikopoulos P., Voglis C., Papageorgiou D.G., Lagaris I.E., "NDL-v2.0: A New Version of the Numerical Differentiation Library for Parallel Architectures", *Comput. Phys. Commun.*, 185:2217-2219, 2014.