

Verifying Parallel Programs with MPI-SPIN

Part 3: Using MPI-Spin

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EuroPVM/MPI 2007
Paris, France
30 September 2007

Overview

1. 4 steps
2. Using `ms`
3. Using `mscc`
4. Executing `pan`
5. Interpreting the output of `pan`
6. Reduction theorems

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4. if counterexample found, play back the trail
 - `./pan -r`
 - sends output to `stdout`

ms syntax

- `ms` (with no arguments)
 - see all command-line options
- `ms [options] foo.prom`
 - generate analyzer source code from `foo.prom`
- `-v`
 - verbose mode
- `-np=<INT>`
 - number of MPI processes (**required**)
- `-DMYMACRO`
 - equivalent to adding `#define MYMACRO` to beginning of model file
- `-DMYTHING=VAL`
 - equivalent to adding `#define MYMACRO VAL` to beginning of model file

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- **-req=<INT>**
 - upper bound on the number of request objects that can be allocated at any one time
 - used only in nonblocking mode (**required**)
 - when upper bound is reached, attempt to post a send or receive request results in **error**

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- `-chansize=<INT>`
 - channel size
 - used only in blocking mode (**required**)
 - for any processes p and q , this is an upper bound on number of messages send from p to q but not yet received
 - when upper bound is reached, sends from p to q block until number of such messages falls below bound
 - hence some executions may not be modeled
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- **-nocancel**
 - optimization for models that do not use `MPI_Cancel`

ms options: deadlock

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- the option has no effect if used with
 - `-req=0`, or
 - `-block -chansize=0`

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- `-n`
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- `-m<INT>`
 - bound search depth to specified number
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- `-i`
 - find a counterexample with the **minimal** number of steps
 - require `mscc` option `-DREACH`

Interpreting the output of pan

- the most important thing
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 - `errors: 0`
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 - an error was found: your property does not hold
- the most important stats
 - number of states explored
 - `74 states, stored`
 - amount of memory used: sum of
 1. `2.622 memory usage (Mbyte)`
 2. `MPI-Spin memory usage (bytes): 125633`

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 - if you restrict to a subset of blocking operations...
 - `MPI_Send`, `MPI_Recv`, `MPI_Sendrecv`,
`MPI_Sendrecv_replace`, `MPI_ANY_TAG`, `MPI_Bcast`,
`MPI_Barrier`, ...
 - **no** `MPI_ANY_SOURCE`

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 - ... then you can conclude
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 - enormous savings in terms of number of states, memory, time
 - even better:
 - any property of the final state of program is independent of interleavings
 - i.e., you can choose any interleaving you want
 - e.g., place an **atomic** block around the body of each process

Reduction with wildcard receives

- **Efficient verification of halting properties for MPI programs with wildcard receives**
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- **Combining Symbolic Execution with Model Checking to Verify Parallel Numerical Programs**
 - S. Siegel, A. Mironova, G. Avrunin, L. Clarke (TOSEM, to appear)
 - if model uses `MPI_ANY_SOURCE`
 - use `atomic` blocks where you like
 - as long as every `MPI_ANY_SOURCE` receive starts a new atomic block

More reduction theorems

- **Verification of Halting Properties for MPI Programs Using Nonblocking Operations**
 - S. Siegel, G. Avrunin (EuroPVM/MPI 2007)
 - extends results above to certain **nonblocking** MPI operations
 - good: `MPI_Isend`, `MPI_Irecv`, `MPI_Wait`
 - bad: `MPI_Test`, `MPI_Waitany`, `MPI_Testany`, `MPI_Waitsome`, `MPI_Testsome`, ...

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- **Semantics Driven Partial-order Reduction of MPI-based Parallel Programs**
 - R. Palmer, G. Gopalakrishnan, R.M. Kirby (PADTAD 2007)
 - **dynamic** partial order reduction

Exercises

1. Consider the C/MPI program `exchange.c` listed on the following slide. Create an MPI-SPIN model of this program and use it to determine whether the program can deadlock.
2. Modify the model to use nonblocking communication to accomplish the exchange. Use MPI-SPIN to determine whether this model can deadlock.
3. Can you find a way to correct `diffusion_par2.c` without introducing a barrier? Use MPI-SPIN to verify your solution.

exchange.c

```
#define UP 0
#define DOWN 1
int main(int argc, char *argv[]) {
    int np, rank, i, sbuf[1], rbuf[1];
    MPI_Status s;
    MPI_Init(&argc, &argv);
    MPI_Comm_size(MPI_COMM_WORLD, &np);
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);
    for (i = 0; i < 3; i++) {
        sbuf[0] = UP;
        MPI_Send(sbuf, 1, MPI_INT, (rank+1)%np, 9, MPI_COMM_WORLD);
        MPI_Recv(rbuf, 1, MPI_INT, (rank+np-1)%np, 9, MPI_COMM_WORLD, &s);
        fprintf(stdout, "Proc %d received %d\n", rank, rbuf[0]);
        sbuf[0] = DOWN;
        MPI_Send(sbuf, 1, MPI_INT, (rank+np-1)%np, 9, MPI_COMM_WORLD);
        MPI_Recv(rbuf, 1, MPI_INT, (rank+1)%np, 9, MPI_COMM_WORLD, &s);
        fprintf(stdout, "Proc %d received %d\n", rank, rbuf[0]);
    }
    MPI_Finalize();
}
```