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run_all (Calls: 1, Time: 66.923 s)

Generated 05-Aug-2024 10:13:54 using performance time.

script in file D:\Aalto\2324\BScThesis\FullRepo\parallelsimulations finitebath\src\run all.m

Copy to new window for comparing multiple runs

Lines where the most time was spent

Line Number	Code	Calls	Total Time	% Time	Time Plot
104	<pre>[vel, el] = diagonal (H);</pre>	25	23.765 s	35.5%	
106	nau = GGE (N, vel);	25	22.556 s	33.7%	
105	E1 = time_evolution (N, hbar,	25	17.355 s	25.9%	
103	H = total_hamiltonian (N,w,mut	25	2.365 s	3.5%	
134	legend([a1(1), a2(1), a3(1)],	1	0.460 s	0.7%	I
All other lines			0.421 s	0.6%	I
Totals			66.923 s	100%	

Children (called functions)

Function Name	Function Type	Calls	Total Time	% Time	Time Plot
diagonal	function	25	23.762 s	35.5%	
GGE	function	25	22.555 s	33.7%	
time_evolution	function	25	17.352 s	25.9%	
total_hamiltonian	function	25	2.282 s	3.4%	I
legend	function	1	0.460 s	0.7%	I
prepareAxes	function	3	0.226 s	0.3%	
xlabel	function	1	0.025 s	0.0%	
title	function	1	0.020 s	0.0%	
hold	function	2	0.018 s	0.0%	
analytical	function	1	0.015 s	0.0%	
ylabel	function	1	0.005 s	0.0%	
Self time (built-ins, overhead, etc.)			0.204 s	0.3%	

Totals | 66.923 s | 100%

Function listing

```
Calls
 time
                 line
                   6 clearvars
                   7 close all
                   8 clc
                  10 % Enable long format for higher accuracy in the calculations
                  11 format long
                  12
                  13 % Initialize the random number generator based on the current time
                  14 rng("shuffle");
                  15
                  16 % Define parallelisation type. Accepted values are 'modular', 'GPU',
                  17 % 'multicore'.
                  18 type = 'GPU';
                  19
                  20 % Add the folders of the parallelisation in the path
                  21 addpath(fullfile(pwd, type));
                  22
                  23 % Begin timing
                  24 tic;
                  25 profile on
                  27 % Defining example variables of the problem
                  29 % The total number of two level systems (TLSs) in the bath.
                  30 % The intially excited state, the qubit, is not considered to be
                  31 % part of the bath. Therefore N+1 is the overall number of TLSs
< 0.001
                  32 N = 1500;
                  33
                  34 % Number of independent, random iterations
< 0.001
                  35 Nr = 25;
                  36
                  37 % The frequency of the qubit.
                  38 % Take it normalized to 1 for simpler calculations
                  39 w = 1;
< 0.001
              1
                  40
                  41 % The reduced Planck's constant.
                  42 % Take it normalized to 1 for simpler calculations
              1
                  43 \text{ hbar} = 1;
< 0.001
                  44
                  45 % A flag that indicates the consideration of internal
                  46 % couplings of the TLSs in the bath. Use 0 for no
                  47 % internal coupling, 1 to include internal coupling
< 0.001
                  48 mutual = 1;
                  49
                  50 % Sets the magnitude of the internal coupling strength.
                  51 % Taken to be w/(5*sqrt(2)) in the example case.
                  52 % For weak coupling regime, smaller of the frequency of the qubit,
                  53 % but is it enough small? Physical explanation for the choosen value?
                  54 gamma = w/(5*sqrt(2));
< 0.001
              1
                  55
                  56 % The final time at which the populations are calculated.
< 0.001
                  57 \text{ tmax} = 8000000000;
                  58
                  59 % Construct a N-by-1 column vector with (sorted) uniformly distributed
                  60 % random numbers in [0, 2*hbar*w]. It will be the diagonal elements of
                  61 % the bath Hamiltonian, representing the energy levels hbar*frequencies
                  62 % of the spins of the bath hbar*omega j where j is in [1, N].
                  63 % The energy levels are sorted to reflect the ordered energy spectrum of
                  64 % the physical systems.
                  65 % This is a constant random vector during the iterations.
```

```
< 0.001
                  66 omega_j = sort(2*hbar*w*rand(N,1));
                  67
                  68 % The initial state of the system, bath in the ground state
                  69 % and qubit excited
                  70 rho0 = zeros(N+1);
< 0.001
              1
< 0.001
              1
                  71 rho0(N+1, N+1) = 1;
                  72
                  73 % The array for collecting the results of long time evolution
                  74 te_results = zeros(N, Nr);
< 0.001
              1
                  75
                  76 % The array for collecting the results of the GGE prediction
< 0.001
                  77 gge results = zeros(N+1, Nr);
                  78
                  79
< 0.001
              1
                  80 if strcmp(type, 'multicore')
                         % Initiate the parallel poll. In local environment uncomment the next line...
                  82
                         % parpool
                         % ... and comment the next line.
                  83
                  84
                         initParPool()
                         % Initialize the random number generator with the Multiplicative lagged
                  85
                  86
                         % Fibonacci generator, for multiple workers in parallel
                         s = RandStream.create('mlfg6331_64','NumStreams', Nr,'Seed',...
                  87
                  88
                         'shuffle', 'CellOutput', true);
                  89
                         % Iterrate Nr times
                         parfor idx = 1:Nr
                  90
                  91
                         RandStream.setGlobalStream(s{idx});
                  92
                         H = total hamiltonian (N,w,mutual,gamma, omega j);
                  93
                         [vel, el] = diagonal (H);
                  94
                         E1 = time_evolution (N, hbar, tmax, vel, el, rho0);
                  95
                         nau = GGE (N, vel);
                  96
                  97
                         te results(:, idx) = E1;
                  98
                         gge_results(:, idx) = nau;
                  99
                         end
< 0.001
                 100 else
                 101
                         % Iterrate Nr times
             1
                 102
                         for idx = 1:Nr
< 0.001
             25
                103
                         H = total_hamiltonian (N,w,mutual,gamma, omega_j);
  2.365
 23.765
             25
                 104
                         [vel, el] = diagonal (H);
17.355
             25
                 105
                         E1 = time_evolution (N, hbar, tmax, vel, el, rho0);
             25
                         nau = GGE (N, vel);
 22.556
                 106
                 107
< 0.001
             25
                 108
                         te results(:, idx) = E1;
< 0.001
             25
                 109
                         gge_results(:, idx) = nau;
             25
                 110
< 0.001
                         end
< 0.001
              1
                 111 end
                 112
                 113 % Get the mean of the iterations
                114 te results mean = sum(te results, 2) / Nr;
< 0.001
< 0.001
                 115 gge results mean = sum(gge results, 2) / Nr;
                 116
                 117 % The analytical GGE prediction for the populations
                 118 [nl, omega] = analytical (N, w, gamma);
  0.015
                 119
                 120 % Plotting
                 121 % (i) Numerical long-time evolution
                 122 % (ii) Numerical GGE
                 123 % (iii) Analytical
                 124
                125 a1 = semilogy(omega_j, te_results_mean, 'o', "Color", 'b');
  0.227
              1
                126 hold on
  0.015
              1
                127 a2 = plot(omega_j, gge_results(1:N), 'x', "LineWidth", 1.1, "Color", "g");
  0.007
                 128 a3 = plot(omega, nl, "LineWidth", 1.2, "Color", "r");
  0.004
              1
< 0.001
                 130 out1 = sprintf('Long-time evolution for %d spins with %d iterations', N, Nr);
```

```
131 xlabel("$\omega\\Omega$", 'Interpreter', "latex", 'FontSize',18)
0.025
            1 132 ylabel("$n$", 'Interpreter', "latex", 'FontSize', 18)
0.005
            1 133 title(out1);
0.020
              134 legend([a1(1), a2(1), a3(1)], 'Long-time evolution', 'Numerical GGE', ...
0.460
                       'Analytical GGE', 'location', "northwest")
               135
               136 %ylim([0.5*10^(-5),10^(-1)])
            1 137 hold off
0.004
               138
            1 139 profile viewer
0.092
               140
               141 % Save the image
               142 relativeFolder = 'output';
               143 filename = sprintf('time_evolution_%d_%d.png', N, Nr);
               144 fullFolderPath = fullfile(pwd, relativeFolder);
               145 fullFilePath = fullfile(fullFolderPath, filename);
               147 % Ensure the directory exists
               148 if ~exist(fullFolderPath, 'dir')
               149
                       mkdir(fullFolderPath);
               150 end
               151
               152 % Define characteristics for the image
               153 exportgraphics(gcf, fullFilePath, 'Resolution', 300);
               155 % If multicore in local environment unccoment the following line
               156 % delete(gcp('nocreate'));
               157
               158 % Output display
               159 disp('The simulation for')
               160 disp(out1)
               161 disp(['was completed in:', ' ', num2str(toc), ' seconds'])
               162 disp(['using parallelisation type', ' ', type])
               163 disp(['with', ' ', getenv('SLURM_CPUS_PER_TASK'), ' ', 'CPUs'])
```