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run_all (Calls: 1, Time: 87.978 s)

Generated 05-Aug-2024 09:48:48 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 |
|-----------------|-----------------------------------|-------|------------|--------|---------------------------------------------------------------------------------------|
| 105 | E1 = time_evolution (N, hbar, ... | 25 | 56.154 s | 63.8% |  |
| 104 | [vel, e1] = diagonal (H); | 25 | 24.980 s | 28.4% |  |
| 103 | H = total_hamiltonian (N,w,mu...) | 25 | 2.746 s | 3.1% |  |
| 125 | a1 = semilogy(omega_j, te_resu... | 1 | 2.144 s | 2.4% |  |
| 134 | legend([a1(1), a2(1), a3(1)], ... | 1 | 1.211 s | 1.4% |  |
| All other lines | | | 0.743 s | 0.8% |  |
| Totals | | | 87.978 s | 100% | |

Children (called functions)

| Function Name | Function Type | Calls | Total Time | % Time | Time Plot |
|---------------------------------------|---------------|-------|------------|--------|---------------------------------------------------------------------------------------|
| time_evolution | function | 25 | 56.141 s | 63.8% |  |
| diagonal | function | 25 | 24.866 s | 28.3% |  |
| total_hamiltonian | function | 25 | 2.686 s | 3.1% |  |
| prepareAxes | function | 3 | 2.131 s | 2.4% |  |
| legend | function | 1 | 1.173 s | 1.3% |  |
| GGE | function | 25 | 0.462 s | 0.5% |  |
| xlabel | function | 1 | 0.045 s | 0.1% | |
| hold | function | 2 | 0.038 s | 0.0% | |
| analytical | function | 1 | 0.036 s | 0.0% | |
| title | function | 1 | 0.021 s | 0.0% | |
| ylabel | function | 1 | 0.010 s | 0.0% | |
| Self time (built-ins, overhead, etc.) | | | 0.367 s | 0.4% | |

| | | | | | |
|--------|--|--|----------|------|--|
| Totals | | | 87.978 s | 100% | |
|--------|--|--|----------|------|--|

Function listing

| time | Calls | line |
|---------|-------|-----------------------------------------------------------------------------|
| | | 6 clearvars |
| | | 7 close all |
| | | 8 clc |
| | | 9 |
| | | 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 = 'modular'; |
| | | 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 |
| | | 26 |
| | | 27 % Defining example variables of the problem |
| | | 28 |
| | | 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 | 1 | 32 N = 1500; |
| | | 33 |
| | | 34 % Number of independent, random iterations |
| < 0.001 | 1 | 35 Nr = 25; |
| | | 36 |
| | | 37 % The frequency of the qubit. |
| | | 38 % Take it normalized to 1 for simpler calculations |
| < 0.001 | 1 | 39 w = 1; |
| | | 40 |
| | | 41 % The reduced Planck's constant. |
| | | 42 % Take it normalized to 1 for simpler calculations |
| < 0.001 | 1 | 43 hbar = 1; |
| | | 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 | 1 | 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? |
| < 0.001 | 1 | 54 gamma = w/(5*sqrt(2)); |
| | | 55 |
| | | 56 % The final time at which the populations are calculated. |
| < 0.001 | 1 | 57 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. |

```

0.002      1  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
< 0.001    1  70 rho0 = zeros(N+1);
< 0.001    1  71 rho0(N+1, N+1) = 1;
           72
           73 % The array for collecting the results of long time evolution
< 0.001    1  74 te_results = zeros(N, Nr);
           75
           76 % The array for collecting the results of the GGE prediction
< 0.001    1  77 gge_results = zeros(N+1, Nr);
           78
           79
< 0.001    1  80 if strcmp(type, 'multicore')
           81     % Initiate the parallel poll. In local environment uncomment the next line...
           82     % parpool
           83     % ... and comment the next line.
           84     initParPool()
           85     % Initialize the random number generator with the Multiplicative lagged
           86     % Fibonacci generator, for multiple workers in parallel
           87     s = RandStream.create('mlfg6331_64','NumStreams', Nr,'Seed',...
           88     'shuffle', 'CellOutput',true);
           89     % Iterrate Nr times
           90     parfor idx = 1:Nr
           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    1 100 else
           101     % Iterrate Nr times
< 0.001    1 102     for idx = 1:Nr
2.746    25 103         H = total_hamiltonian (N,w,mutual,gamma, omega_j);
24.980    25 104         [vel, el] = diagonal (H);
56.154    25 105         E1 = time_evolution (N, hbar, tmax, vel, el, rho0);
0.467    25 106         nau = GGE (N, vel);
           107
< 0.001    25 108         te_results(:, idx) = E1;
< 0.001    25 109         gge_results(:, idx) = nau;
0.001    25 110     end
< 0.001    1 111 end
           112
           113 % Get the mean of the iterations
0.001    1 114 te_results_mean = sum(te_results, 2) / Nr;
< 0.001    1 115 gge_results_mean = sum(gge_results, 2) / Nr;
           116
           117 % The analytical GGE prediction for the populations
0.036    1 118 [nl, omega] = analytical (N, w, gamma);
           119
           120 % Plotting
           121 % (i) Numerical long-time evolution
           122 % (ii) Numerical GGE
           123 % (iii) Analytical
           124
2.144    1 125 a1 = semilogy(omega_j, te_results_mean, 'o', "Color", 'b');
0.060    1 126 hold on
0.011    1 127 a2 = plot(omega_j, gge_results(1:N), 'x', "LineWidth", 1.1, "Color","g");
0.006    1 128 a3 = plot(omega, nl, "LineWidth", 1.2, "Color", "r");
           129
< 0.001    1 130 out1 = sprintf('Long-time evolution for %d spins with %d iterations', N, Nr);

```

```

0.052 1 131 xlabel("\omega/\Omega$", 'Interpreter',"latex", 'FontSize',18)
0.014 1 132 ylabel("$n$", 'Interpreter',"latex", 'FontSize',18)
0.025 1 133 title(out1);
1.211 1 134 legend([a1(1), a2(1), a3(1)], 'Long-time evolution', 'Numerical GGE', ...
135     'Analytical GGE', 'location', "northwest")
136 %ylim([0.5*10^(-5),10^(-1)])
0.009 1 137 hold off
138
0.051 1 139 profile viewer
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);
146
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);
154
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'])

```