

Spin Glass Automata (SGA): An Evolutionary Local Search Automata for Solving Optimization Problems

M. V. Jahan^{1,*}, S. M. S. Kenari²

¹Software Engineering Department, Mashhad Branch, Islamic Azad University, Mashhad, Iran

²Software Engineering Department, Nowshahr Branch, Islamic Azad University, Mazandaran, Iran

Abstract Nowadays, new optimization problems become so complicated that popular classic methods are unable to solve them in a reasonable time. By using heuristic and evolutionary algorithms, these problems can be solved more quickly. Cellular Automata (CA) and Spin Glasses (SG) are examples of such algorithms. A CA is a self-organized machine with a simple structure and complicated behavior. Due to local interactions between its cells, it has a high speed; however, it is unable to solve optimization problems. On the other hand, SGs, due to inter-spin magnetic interaction as well as following thermodynamic rules, continually have the tendency toward lower energy, and this way it can solve optimization problems. The interactions between spins are usually limited and consequently, aforementioned methods cannot solve optimization problems with an expected accuracy and this is a great impact. In this paper, with inspiration from behavior of CA and SG, a new machine named Spin Glass Automata (SGA) is introduced which has behavioral dynamic of SG and CA. It has also the ability of parallel processing while it is rapid enough. It follows rules of a CA too. With these properties at hand, the machine can be used in a vast area of optimization problems. Behavioral dynamic tests of the machine include phase transition, entropy and correlation as well as the comparison to other heuristic algorithms such as (GA, TS, SA and NN) shows the ability of this machine in solving practical optimization problems. It has an appropriate execution speed and accuracy. To test full potential of this machine, the NP problem such as optimal portfolio selection has been solved and compared with the five famous stock market of the world

Keywords Spin Glass Model, Portfolio Selection, Cellular Automata, Phase Transition and Parallel Processing

1. Introduction

Heuristic algorithms are the strongest classes of optimization algorithms which are receiving more attention and application increasingly. Due to a great extension in practical problem domains and increasing of the need for rapid information processing, parallel processing and rapid problem solver algorithms becomes more important. Most of parallel processing enable algorithms are inspired with natural processing. The main idea of these algorithms is local processing and global conclusion which namely are parallel genetic algorithms, parallel colony ants, parallel simulated annealing[1], cellular automata[2] and spin glasses[3].

CA is a self-organized machine that is composed of an array of cells which each of them can own several finite states and can be updated through specific rules, discretely and simultaneously. While having the ability of parallel and local processing, this machine has a great speed in processing, too[2]. Behavioral dynamic of these machines, generally

ends to specific finite states and this ability is denoted as self-organizing[4]. Aforementioned properties caused these machines to be used in encryption[5], image processing[5] and physical behavior simulating[6,7]. The great inabilities of the automata are, being unknown rules and inability in solving engineering problems specially optimization problems[6].

SG is a network of spins in which spins interact magnetically and consistently change their values to achieve a lower level of energy. When the system is in minimum energy (or temperature) state, there is no significant variation in spin values and the system has reached ground state (the state with minimum energy)[8]. This model has many characteristics among which are non-exponential growth of ground states with an increase in the number of spin bonds, effectiveness of environmental factors such as temperature on network behavior and ability to achieve the optimum state at variant temperatures[9]. Also, in this model, each spin is a portion of solution while for example in genetic algorithms a chromosome is, of problem solutions itself[10]. Since the model is physically well specified, via modeling behavior of some engineering problems through it, one can solve these problems which are otherwise unsolvable conveniently [11,12].

* Corresponding author:

VafaeiJahan@mshdiau.ac.ir (M. V. Jahan)

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The differences between SG and CA are that SGs follows thermodynamic rules, they are based on temperature (it is temperature that define the final state of SGs), Spins have interactions in the network and the system has a energy function (or Hamiltonian function)[8]; while CA is based on time (it is time that define the final state of automata), cells have no interaction between themselves, only interact according to defined rule and have no defined energy function [2].

In this paper, a new machine named spin glass automata (SGA) is introduced, which not only has the properties of CA but also follows thermodynamic rules of SG. In other words, this automata has the property of self-organizing, in specific temperature, has local behavior and cells are interacted with each other through defined rules. Furthermore, this machine has the ability of parallel and simultaneous processing capability that can solve optimization problems more efficiently.

To study dynamic of this machine, testing upon random data can be an option but due to lack of a performance metric, it will be uncertain. Therefore the problem of optimal portfolio selection with initial constraints that is a non-deterministic polynomial problem (NP-Complete) is solved by this machine as a case study. Results of this experiment with actual data are compared and evaluated against accurate results from five well-known stock market of the world. The machine shows desirable speed and accuracy in solving the problem. In addition to aforementioned experiments, the dynamic of SGA is investigated through phase transition, correlation and entropy analysis. Finally, the performance of the automata is evaluated, in solving optimal portfolio selection problem, against that of four famous heuristic algorithms namely as genetic algorithm(GA), simulated annealing(SA), tabu search(TS) and neural network(NN).

Therefore, first in section 2, the related activities to spin glasses are described as well as the combination of cellular automata and spin glasses. In sections 3 and 4, a nominal presentation of cellular automata and spin glasses are presented. In section 5, SGA and the problem of optimal portfolio selection are introduced. In section 6, experimental results such as the phase transition, entropy analysis and correlation test, are introduced to evaluate trustiness of prescribed method based on the fundamental data from five stock market of the world. Furthermore, execution speed and accuracy as well as Pareto front of the machine is compared to that of four well-known heuristic algorithms (SA, GA, TS and NN).

2. Related Activities

CA is a simple and rapid machine and its operation is done locally with the ability of parallel processing. This way, the hardware implementation of the machine against other similar models, becomes simpler[2]. CA was used in encryption, compression and image processing[4]. It has also

applications in chemical processes simulation, traffic and numerical calculus[13].

SG is used in solving a lot of optimization problems. As an example, Hartmann et al. in[11,12] introduced several engineering problems such as Maximum Flow and Matching. Nishimori in[14,15], has investigated application of SGs for data transfer in noisy environment. In[16], Horiguchi et al. introduced routing in adaptive and distributed computer networks as another application area. Gabor et al. in[17] are the first ones who used spin glass to solve optimal portfolio selection problem. Finally, Akbarzadeh and Vafaei Jahan are pioneers who introduced a rapid method in[3] which could find optimal state with an appropriate accuracy for spin glasses with a limited bonds (limited connection) and that are using migration and elite operators (similar to genetic algorithm operators).

SG and CA, has been combined in several ways in the past. As an example in[18], the Ising model is simulated using CA in a way that introduced cellular automata demonstrates Ising model like behaviors. In[19], an Ising cellular automata (ICA) is introduced which has thermodynamic properties of Ising model. Also in[20], the Q2R rules are introduced which caused, cellular automata to be used as an Ising model simulator but in the absence of thermal effects.

3. Cellular Automata

CA is a simple machine which is able to show complex behavior. Mathematical model of the machine is a four tuple $\{A, S, N, f\}$, where Z is the cellular network of automata. S is the set of possible values each cell can own. N is the adjacent set of each cell and finally f is transition rule or rule briefly. At each time, the value of cell S can be reached through applying rule f on the adjacent set N [2]. If all cells of the CA include a same rule, it is called homogenous CA and is called heterogeneous otherwise[2,6]. If boundary cells (first and last cells) are neighbors, the cellular automata is called CA with circular boundary and non-circular otherwise [2,6].

4. Spin Glass

Ising model is a network of spins which has either value of +1 or -1. These spins are interacted with each other according to a magnetic property among themselves and continually change themselves to reach lower energy level. Once the network is in the lowest energy level, no additional change will be seen in spin values and the network reaches ground state. Spin Glass Ising model is the same as Ising model except the value of each spin is between $[-1, +1]$ rather than -1 and +1. This model is briefly called Spin Glass, too[8].

Theoretically, SG, is a network of N spins that its structure can be either limited connection (connection with adjacent spins e.g. with 2D or 3D structure) or full connection¹ (connection between each spin to all other spins). Spins have

either ferromagnetic or anti-ferromagnetic property. According to before mentioned properties, network energy is calculated using the following formula[8,12]:

$$E(\{x_i\}) = \left[-\frac{1}{2} \sum_{i=1}^N \sum_{j=1}^m x_i J_{ij} x_j \right] + \left[-\sum_{i=1}^N h_i x_i \right] \quad (1)$$

where $E(\{x_i\})$ is the energy of all spins, The sum i, j runs over all pairs of nearest neighbors, m is the number of nearest neighbors of each spin i that can be $m=4$ in Van Neumann cellular automata (CA) or $m=8$ in Moore CA [16] or $m=N$ for full connection. J_{ij} denotes the strength of the bond connecting spins i and j . $J_{ij} > 0$ describes a ferromagnetic interaction, while $J_{ij} < 0$ describes an anti-ferromagnetic interaction. The quantity h_i is the external field acting on spin i and describes the energy due to the spin's orientation. Also, the factor $\frac{1}{2}$ corrects for double counting of the interaction between every two neighboring spins. Here the task is to find a spin configuration x_i that minimizes the energy of the spin glass, given $\{J_{ij}\}, \{h_i\}$. A scheme for a 2D spin glass is depicted in Figure 1.

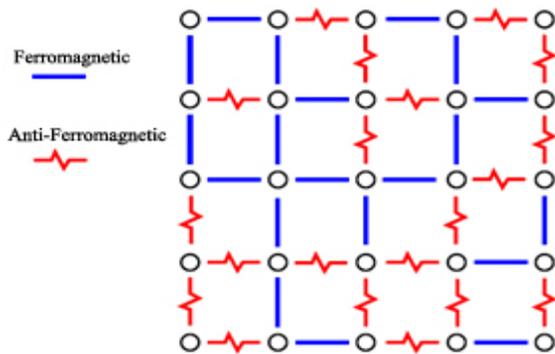


Figure 1. A two-dimensional spin glass with bond disorder. Spins are placed on the sites of a regular grid. They interact with their neighbors, the interaction is random, either ferromagnetic (straight lines) or anti-ferromagnetic (jagged lines)[11].

5. Spin Glass Automata

In this paper, a 2D spin glass CA is defined as an ordered 6-tuple $\{Z, S, J, h, f, H\}$ such that Z is a 2D network where each spin (cell) can interact with adjacent ones based on Van Neumann or Moore models (short range spin glass); S is the value of each spin can take between $[0,1]$; Each spin has an interaction (bond) with its nearest neighbor; A strength of the bond J_{ij} exists between each pair neighbor spins; The quantity h_i is the external field acting on spin i and describes the energy due to the spin's orientation; The rule f is the interaction rule between nearest neighbors which can be any defined rules for the Automata (It is the same as CA rules). This rule defines the type of interactivity; and H is Hamiltonian function of interaction or energy function for the whole automata. It usually constitutes of interactions between composer parts of the system and is calculated

based on the problem type[8], and finally, it is calculated similar to Hamiltonian rule of SG.

In SGA, the subsequent state of a spin is related to current states of the spin and its nearest neighbors and is specified based on rule f . Since SGA is moving through losing energy H , the amount of spins' interaction energy J and external field energy h_i , together prevent spins to take any value S , i.e. only change which can cause any decrease in SGA's energy will be accepted.

When the SGA starts, by using rule f and interactions with its neighbors, each spin tries to decrease its energy (changes its value) according to function H . If the change in value results in local energy decrease, the change will be accepted

and with probability of $e^{-\frac{\Delta H}{kT}}$ otherwise; where k is Boltzmann's constant and T is the environment temperature. Experiments show that in spite of local energy reduction, global decrease is also observed in the machine. Therefore the machine is useful in optimization problem solving.

Using SG properties, the aforementioned SGA can follow thermodynamic rules and goes through equilibrium in any temperatures. Also through using CA capabilities, it can use any rule of it and produce finite states. The advantage of SGA against two other prior models is that it can show stable state based on any interaction rules between spins. If SGA temperature be high, it can produce all possible states. If temperature reduces, the automata goes through equilibrium with environment via losing energy.

5.1. SGA Dynamics

Structural and behavioral dynamic studies of the machine, requires investigations such as thermodynamic behavior of it in various temperatures, entropy changes, the amount of energy decrease to minimum at equilibrium temperature and also phase transition temperature which are all dealt with in detail in the next section. To evaluate dynamic of the proposed machine, a case study is inevitable. To this end, the non-polynomial problem of optimal portfolio selection is chosen. The selection shows not only ability of the machine to solve engineering problems, but also its behavioral dynamic based on real data.

In this paper, a heterogeneous 2D SGA with non-circular boundary condition is used which spin values S include all real numbers in $[0,1]$ and the state of each spin changes because of interactions with adjacent spins. The rule f is an average rule which is calculated via summation of the nearest neighbors values divided by their total count. The Hamilton function H , is the energy function of formula (1) which its calculation method is described in the next section.

5.2. Portfolio Selection Problem and SGA

Portfolio selection problem [20] is stated by Markowitz model as if N stocks are assumed and for each stock i , μ_i is the mean return of asset i , and σ_{ij} is the covariance between returns of assets i and j . The decision variable x_i represents the fraction of capital to be invested in asset i . following multi-objective problem has to be optimized:

$$\min \sum_{i=1}^N \sum_{j=1}^N x_i \sigma_{ij} x_j \quad (2)$$

$$\max \sum_{i=1}^N \mu_i x_i \quad (3)$$

$$\text{Subject to } \sum_{i=1}^N x_i = 1 \quad (4)$$

$$0 \leq x_i \leq 1, i = 1..N \quad (5)$$

The formula (2) and (3) are two objective functions which must be satisfied based on the constraints (4) and (5). In formula (3), μ_i is achieved from mean return of each asset in defined intervals, in such a way that if W_{bi} be value of stock i in beginning of the interval and W_{ei} be that of interval end, then

$\mu_i = \sum_{t=1}^n \frac{W_{ei}(t) - W_{bi}(t)}{W_{bi}(t)}$ where n will be the number intervals which stock i is considered in.

In this paper, for a simpler solving of the multi-purpose problem, it converts to a multi-modal problem:

$$\text{Minimize } \lambda \left[\sum_{i=1}^N \sum_{j=1}^N x_i \sigma_{ij} x_j \right] + (1 - \lambda) \left[- \sum_{i=1}^N \mu_i x_i \right] \quad (6)$$

$$\text{Subject to } \sum_{i=1}^N x_i = 1 \quad (7)$$

$$0 \leq x_i \leq 1, i = 1..N \quad (8)$$

In the formula (6), $\lambda \in [0, 1]$ is called risk aversion parameter. It shows the amount of risk effects in optimization function. If $\lambda = 0$ then Eq. (5) represents maximum portfolio mean return (without considering the variance) and the optimal solution will be formed only by the asset with the greatest mean return. The case with $\lambda = 1$ represents minimizing the total variance associated with the portfolio (regardless of the mean returns) and the optimal solution will typically consist of several assets. Any value of λ inside the interval (0,1) represents a tradeoff between mean return and variance, generating a solution between the two extremes, $\lambda = 0$ and 1.

To solve optimal portfolio selection problem, as is also investigated in [3], it is assumed that each stock is a spin which has a value between 0 and 1. This automata owns an energy function like formula (1) while the problem goal is to select an optimal portfolio and to minimize formula (6). Thus, through the following mapping formula (6) can be converted to the formula (1):

$$J_{ij} = -2\lambda\sigma_{ij} \quad (9)$$

$$h_i = (1 - \lambda)\mu_i \quad (10)$$

This way, the interaction between two spins (stocks) can be calculated via formula (9) and the amount of internal energy of each spin can be computed through formulation of (10). Therefore, first the aforementioned mapping in formulations of (9) and (10) is performed; Stocks are dispatched in SGA randomly and then each stock value initializes by a random number such that condition (7) be satisfied. In

thermodynamic view point, the system temperature is raised to the extent that which any possible state can producible at and λ takes a number greater than 0 and lower than 1 (to be equal the effects of risk and mean return, value of 0.5 is selected). The approach, each spin interacts with its nearest neighbors, is such that the amount of each spin is summed with its nearest ones and is averaged while limitations of (4) and (5) are satisfied all over the process. Now, when the machine starts at a random state, through neighbor spins interactions and consequently spins value changes, it tries to find the lowest amount of energy at the current temperature. Since the lowest amount of energy is also the value of cost function (6), states of spins will show optimal portfolio at temperature near to zero.

6. Experimental Results

To evaluate performance of proposed machine, Experiments on the benchmark data were originally performed in [21]. These benchmark data are presented in text file format as follows: Number of assets (N); and for each asset i ($i=1, \dots, N$): mean return as well as standard deviation of return; for all possible pairs of assets: i, j , correlation between asset i and asset j . The above data were taken from five major stock exchange markets, during the time period extending from March 1992 to September 1997. These five stock exchange markets include Hang Seng in Hong Kong (31 assets), Deutscher Aktien Index (DAX100) in Germany (85 assets), Financial Times London Stock Exchange (FTSE100) in Britain (89 assets), Standard & Poor's (S&P100) in USA (98 assets), and Nikkei in Japan (225 assets). Probability density function (pdf) of covariance $P(J)$ of each stock market for the given data has been shown in Figure. 2.

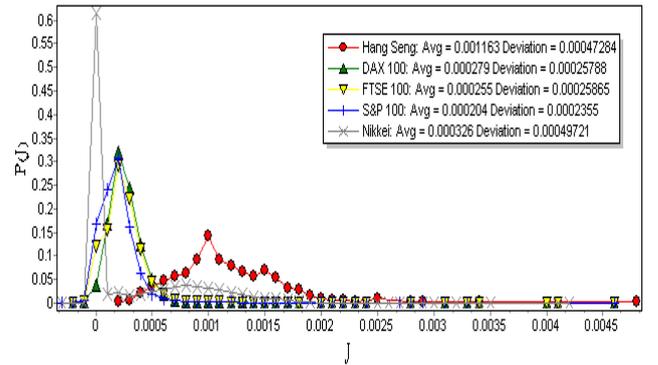


Figure 2. Probability density functions for covariances between assets in 5 major stock markets from 1992 to 1997 from data in[21].

As can be observed, the $P(J)$ of the five given stock markets have small mean and variance. And also $P(J)$ are similar together.

6.1. SGA Dynamic: Local Behavior with Parallel Processing Ability

Experiments show that local behavior of SGA results in a considerable increase in convergence speed. In Table (2), run

time of proposed algorithm is presented. As is shown the run time is so much better than global behavior of glass and ever its local behavior. But response accuracy is acceptable while is less than that of global behavior. This shows that trapping in local optimum is possible yet. Among the factors which speed up SGA, is its absolute local while targeted behavior which causes, as an instance for stock market Nikkei, response time to be 6 times more than that of global behavior while the accuracy is appropriate. This local behavior brings machine with parallel processing capability. Experiment results shown in table (1), which is to investigate parallel processing capability of this algorithm and for various stocks, show that parallel processing ability of the process can considerably speeds up it. To this end, in this experiment, it is possible to break down the automata into smaller and discrete automatas and to run each automata on one processor (breaking down method doesn't follow any predefined rule and can be done arbitrarily).

Table 1. Parallel processing ability in SGA, each stock market with defined number of processors, has been run 10 time on average.

SGA	Number of Processors	Average Run Time (ms)
Hang Seng (N=31)	1	490
	2	326
	3	723
DAX 100 (N=85)	1	16189
	2	12543
	3	11321
FTSE (N=89)	1	29071
	2	25654
	3	23541
S&P 100 (N=98)	1	25239
	2	24213
	3	25434
Nikkei (N=225)	1	71578
	2	65212
	3	58212
	4	57992

Simultaneous running of these automata speeds up convergence considerably (since the authors didn't access to multi-processor computers, they used multi-programming instead). In any iteration, each program selects a spin and applies the algorithms on it. For example in a typical iteration 3 spins are selected and optimized by 3 programs. Table (1) shows the average run time of the automata. Since oper-

ating system limitations (e.g. context switch) exist for simultaneous running of several programs, their numbers (processors count) cannot be increased infinitely. As can be seen, the speed of response time is increased due to parallel processing. As an instance, for Hang Seng stock market, average response time for 1 processor is 490 milliseconds and is 326 milliseconds for 2 processors. Average response time for 3 processors is 723 milliseconds which is more than that of 2 processors due to multi programming limitations. For stock market Nikkei which has 225 stocks, 4 processors outperform 2 and 3 processors and make results available sooner.

6.2. SGA Dynamic: Phase Transition

Phase transition typically refers to a transformation of a thermodynamic system from one state to another. This transformation is typically marked by a sudden change in some physical property as a result of a small change in what is called an order parameter (e.g., temperature)[26,27]. Here, the critical temperature (T_c) of SGA is the temperature at

which phase transition occurs. At this temperature, the probability of reaching the ground state of the glass changes significantly, i.e. automata's states become more/less unlikely to be a ground state. Figure. 3, illustrates critical temperature of SGA applied to the portfolio selection problems. As shown in this figure, the probability of reaching ground state ($\frac{E}{E_{min}}$) is equal to 1 in low temperatures, where

E is the amount of SGA energy and E_{min} is the actual minimum of the energy function. As temperature rises, glass energy also changes and the probability of reaching ground state is expected to gradually decrease. However, this drop is delayed until phase transition where a sudden drop occurs (In this paper, phase transition is defined to occur at $\frac{E}{E_{min}} < 0.99$). After the critical temperature, the probability of reaching ground state decreases significantly. Figure 3 illustrates the phase transition for the five stock markets in this paper. The vertical solid red line on $T_c = 1.25 \times 10^{-6}$ highlights the proximity of this transition for most of the benchmarks. After this temperature, system is not likely to be in its optimal ground state. The dashed gray lines indicate the range of critical temperature for different benchmarks.

Table 2. SGA Behavior and its equivalent Local and Global Behavior of Spin Glass from[3] Comparison. The value of cost functions are resulted from 100 times running of the algorithm. Run time, found cost function average and cost function error average against that of actual cost function are be shown.

Stock	Spin Glass Global Behavior[3]			Spin Glass Local Behavior with Moore Neighbor Model[3]			SGA Behavior with Moore Neighbor Model		
	Run Time (ms)	Cost Function	Cost Function Average Error	Run Time (ms)	Cost Function	Cost Function Average Error	Run Time (ms)	Cost Function	Cost Function Average Error
Hang Seng (N=31)	4626	-0.00336	7.2×10^{-12}	2180	-0.0029	4.6×10^{-4}	490	-0.0031	3.1×10^{-5}
DAX 100 (N=85)	65268	-0.00412	2.6×10^{-11}	17659	-0.0038	3.2×10^{-4}	16189	-0.004	1.3×10^{-5}
FTSE (N=89)	75631	-0.00335	2.0×10^{-11}	31981	-0.0029	4.5×10^{-4}	29071	-0.0028	5.5×10^{-4}
S&P (N=98)	81231	-0.00363	1.76×10^{-11}	32039	-0.0031	5.3×10^{-4}	25239	-0.0035	1.1×10^{-5}
Nikkei (N=225)	418140	-0.00145	2.76×10^{-10}	81321	-0.0012	2.5×10^{-4}	71578	-0.0012	1.15×10^{-4}

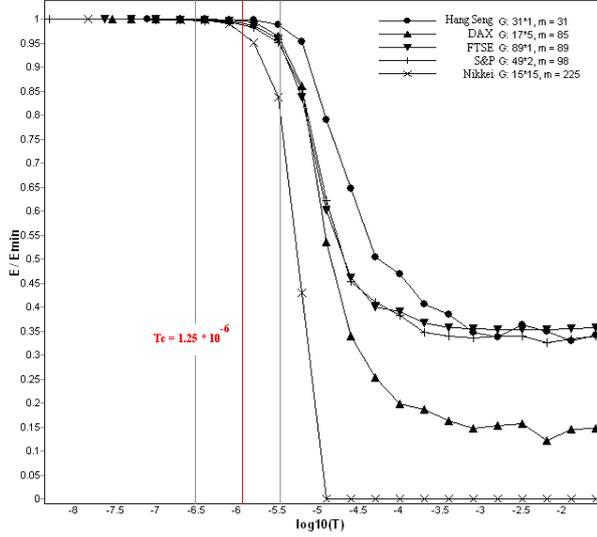


Figure 3. SGA Phase Transition Temperature for the 5 stock markets mentioned in[21]. Phase Transition Temperatures are almost equal for all stocks.

6.3. SGA Dynamic: Correlation Test

The correlation test is used to verify the correctness of our previous test as follows. Two randomly distributed SGAs are independently considered on same stock market data. Both Automatas are used to find the ground state separately. The rate of correlation is then calculated for the two glasses (in lieu of amount of each corresponding cell). The correlation level at each temperature is determined as follows:

$$\rho = \frac{\sigma_{ij}}{\sigma_i \cdot \sigma_j} \quad (11)$$

Where σ_{ij} is the covariance of spins between two corresponding spins of each glass, and σ_i and σ_j are the standard deviation of cells of SGA i and j , respectively.

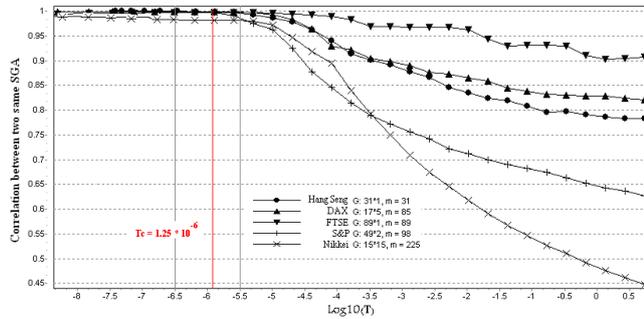


Figure 4. Correlation test of the two SGAs for same stock market.

As illustrated in Figure. 4, the correlation initially remains the same even with gradual increase of temperature. After critical temperature, however, the correlation suddenly drops and each SGA acts separately from the other. It should be noted that critical temperature in this test is the same as the calculated temperature in the first test, shown in Figure 3.

6.4. SGA Dynamic: Entropy Analysis

According to third thermodynamic rule, in near zero

temperature, system has its minimum entropy. In another experiment it is shown that the rule is valid for SGA while is invalid for CA[22,23]. In equilibrium temperature, the probability by which system is placed in state i depends on system energy on that state and is calculated through the following equation[24]:

$$p_i = \frac{e^{-E_i}}{\sum_i e^{-E_i}} \quad (12)$$

Equation (12) is the probability of being in state i divided by summation over all states. Entropy is calculated using equation (13)[24]:

$$H = -\sum_i p_i \log p_i \quad (13)$$

As can be observed in Figure. 5, SGA entropy is decreased when temperature decreases. It is worth noting that entropy is almost constant on phase transition above temperatures but at phase transition temperature starts decreasing suddenly. In fact calculated phase transition temperatures for optimal portfolio selection problem in Figures 3, 4 and 5 are almost equal.

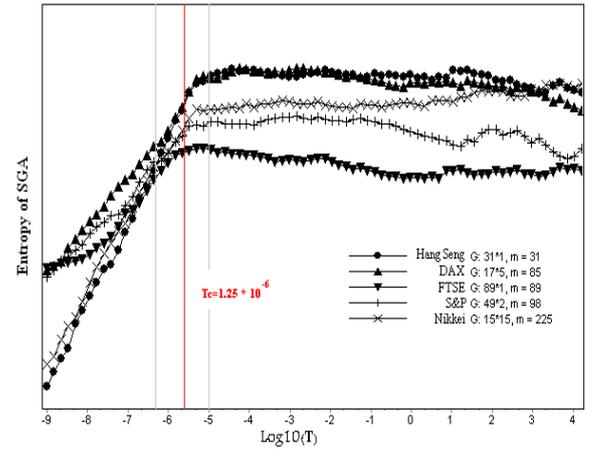


Figure 5. SGA entropy for solving optimal portfolio selection problem.

7. SGA and Other Heuristic Algorithms

SGA performance is evaluated against other heuristic algorithms in another experiment. Table (3) represents SGA behavior against that of four optimization problem solver algorithms. As illustrated in the table, SGA convergence speed is more than that of optimization methods GA, TS, SA and NN which are investigated in[25] (The run time of the proposed method and the current methods in paper[25], due to the differences in computers and similar simulation software, is not the same. Because, the authors implement the GA, TS, SA and NN, according to the author's understanding of the paper[25]; in table (3) the new run time results have been listed. Some differences in the experimental run time results is observed but the conclusion seems generally correct), while its accuracy in finding result is already acceptable. As an instance for Nikkei and Hang Seng stock market, the convergence speed is more considerably.

Table 3. SGA, GA, TS, SA and NN Comparison. Each stock market is solved by any of the methods 100 times; average run times, and average mean return error (AMRE) are listed. Experimental results of GA, TS, SA and NN extracted from[27]. The AMRE result is the exact copy of[27], but the run time is the comparable result (Comparable in the same computer) of the GA, TS, SA and NN personal implemented.

Stock		SGA	GA	TS	SA	NN
Hang Seng (N=31)	AMRE ($\times 10^3$)	1.25	1.1321	1.1237	1.1203	1.2316
	Run Time (s)	4	34	21	20	472
DAX 100 (N=85)	AMRE ($\times 10^3$)	1.2919	2.4457	2.6668	2.3896	1.5776
	Run Time (s)	16	90	59	110	761
FTSE (N=89)	AMRE ($\times 10^3$)	1.1383	0.7310	0.7357	0.9512	1.2513
	Run Time (s)	29	162	60	88	806
S&P (N=98)	AMRE ($\times 10^3$)	1.485	1.3236	1.3130	1.7251	1.7922
	Run Time (s)	25	149	52	71	1319
Nikkei (N=225)	AMRE ($\times 10^3$)	0.9027	1.1415	0.5510	0.5458	1.4737
	Run Time (s)	72	720	132	180	3211

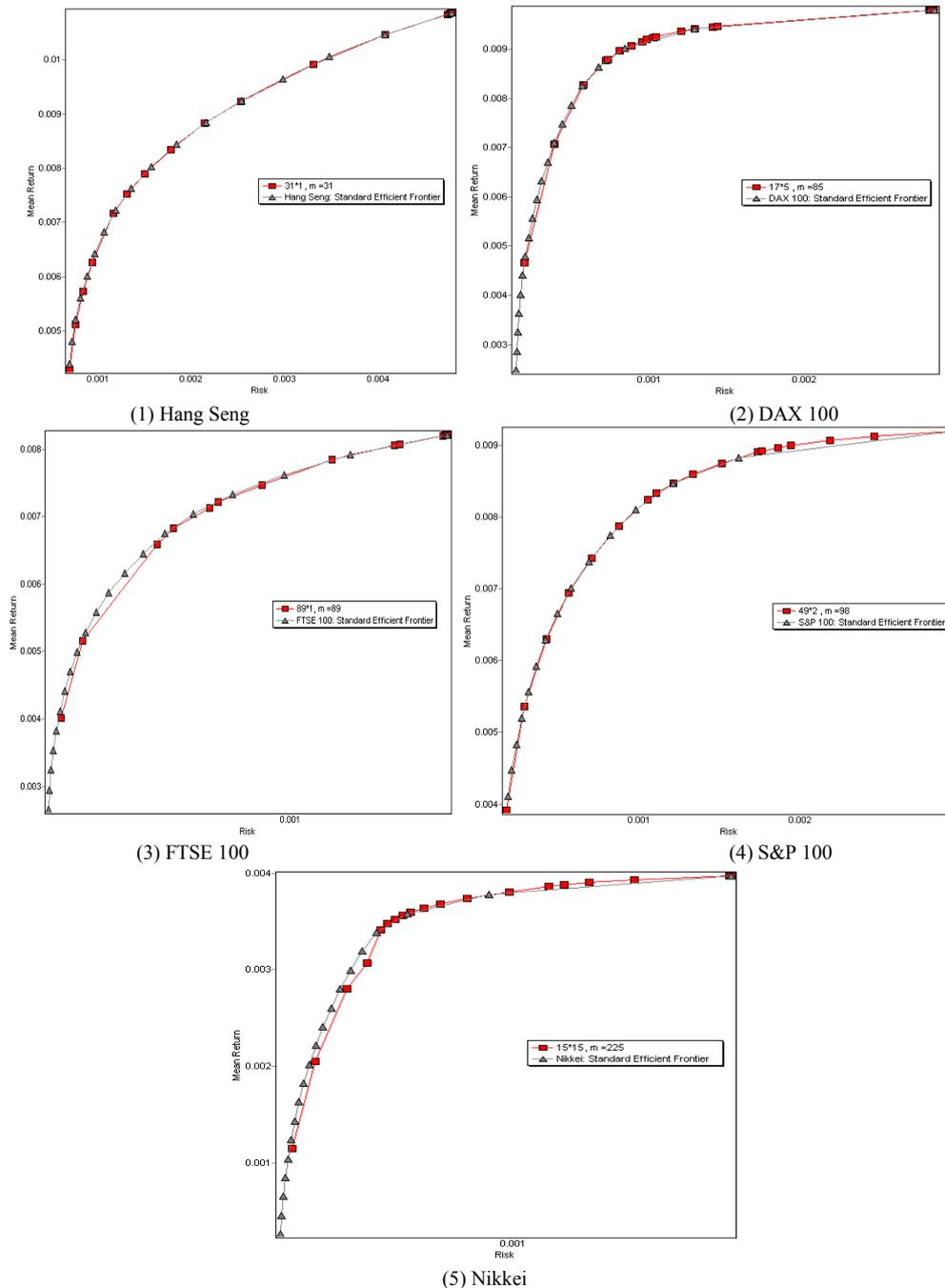


Figure 6. The Pareto front resulted from running SGA against standard optimal front resulted from benchmark data of[21].

8. SGA Pareto Front and Comparison to Optimal Front

In another experiment optimal outcomes front resulted from SGA with various λ s is depicted. In Figure. 6 movement path is toward to Pareto front and for the 5 aforementioned stock market, shows validity of energy decrease and finding result. In the figure, resulted optimal Pareto front for various λ s is depicted in gray lines. To specify Pareto front accurately, for any stock, the λ amount is drawn in range 0.05 to 0.95 with 0.05 differences. For any value of λ , the algorithm is run with 1000 iterations and mean return and its risk is drawn. Comparing base and resulted Pareto front shows operation validity of the proposed machine in finding optimal result with various λ s.

9. Conclusions

In this paper, a new machine called spin glass automata is introduced. The machine has the properties of cellular automata as well as spin glass. In other words, this automata has property of self-organizing in specific temperature; also has local behavior that cells interact according to any defined rule (inspired by cellular automata) and also according to thermodynamic rules (inspired by spin glass). It has parallel and simultaneous processing ability and can actually solve optimization problems.

Also in this paper dynamic of the machine investigated in solving optimal portfolio selection and experimental results compared and evaluated against real data with accurate result from 5 well known stock market of the world. The results show that speed and accuracy of the machine is appropriate in solving the problem and thus can be used as an optimization problem solver algorithm.

In addition to mentioned experiments, SGA dynamic investigated through phase transition, entropy and correlation tests. Phase transition and entropy tests show that the machine follows thermodynamic rules. Correlation test, demonstrates ability of the machine to find global optimums. At the end, performance of the algorithm is compared to that of four famous heuristic algorithms (genetic algorithm, simulated annealing, Tabu search and neural network) in solving optimal portfolio selection problem. Also Pareto front is evaluated to show the ability of the machine to cover finding result for various risks.

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