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Efficient Ant Colony Optimization (EACO) Algorithm for Deterministic Optimization

Urmila M Diwekar*and Berhane H Gebreslassie

Center for Uncertain Systems: Tools for Optimization & Management (CUSTOM), Vishwamitra Research Institute, Crystal Lake, USA

Abstract

In this paper, an efficient ant colony optimization (EACO) algorithm is proposed based on efficient sampling method for solving combinatorial, continuous and mixed-variable optimization problems. In EACO algorithm, Hammersley Sequence Sampling (HSS) is introduced to initialize the solution archive and to generate multidimensional random numbers. The capabilities of the proposed algorithm are illustrated through 9 benchmark problems. The results of the benchmark problems from EACO algorithm and the conventional ACO algorithm are compared. More than 99% of the results from the EACO show efficiency improvement and the computational efficiency improvement range from 3% to 71%. Thus, this new algorithm can be a useful tool for large-scale and wide range of optimization problems. Moreover, the performance of the EACO is also tested using the five variants of ant algorithms for combinatorial problems.

Introduction

A wide range of optimization problems, which include a large number of continuous and/or discrete design variables, fall into the category of linear programming (LP), nonlinear programming (NLP), integer programming (IP), mixed integer linear programming (MILP), and mixed integer nonlinear programming (MINLP). Gradient based methods such as Branch and Bound (BB), Generalized Bender's Decomposition (GBD), and Outer Approximation (OA), are generally used for solving IP, MILP, and MINLP problems. However, these methods have limitations whenever, the optimization problems do not satisfy convexity conditions, the problems have large combinatorial explosion, or the search domain is discontinuous [1]. Metaheuristic optimization strategies such as simulated annealing (SA) [2], genetic algorithm (GA) [3] and ant colony optimization (ACO) [4] can provide a viable alternative to the gradient based programming techniques. Review on metaheuristic algorithms can be viewed at Kahraman et al. [5]. One of the metaheuristic optimization techniques that attract researchers in recent years is an ACO algorithm. ACO algorithm was first introduced by Dorigo [4] and it has been receiving increased attention because of a series of successful applications in different disciplines such as routing, scheduling, machine learning, assignment, and design problems, and different branches of engineering [6-10]. Even though, ACO algorithm was practiced with great success for combinatorial optimization problems, in recent years, the research focus has been on extending the combinatorial ACO algorithm to continuous and mixed variable nonlinear programming problems. In this work, a new variant of ACO algorithm to solve deterministic optimization problems is proposed.

An ant colony is a population of simple, independent, asynchronous agents that cooperate to find a good solution to an optimization problem. In the case of real ants, the problem is to find good quality of food in a close vicinity of the nest, while in the case of artificial ants, it is to find a good solution to a given optimization problem [6]. A single ant is able to find a solution to its problem, but only cooperation among many individual ants through stigmergy (indirect communication) enables them to find good or global optimal solutions. Therefore, the ACO algorithm is inspired by the ants' foraging behavior. Natural ants deposit pheromone on the ground in order to mark some favorable path that should be followed by other members of the colony. ACO

algorithm exploits a similar mechanism for solving optimization problems [4]. It was originally introduced to solve combinatorial optimization problems, in which decision variables are characterized by a finite set of components. However, in recent years, its adaptation to solve continuous [11-14] and mixed variable [15-17] programming problems has received an increasing attention.

In this paper, a new strategy that improves the computational performance of ACO algorithm by increasing the initial solution archive diversity and the uniformity of the random operations using a quasi-random sampling technique referred as Hammersley Sequence Sampling (HSS) [20]. This sampling mechanism has been shown to exhibit better uniformity property over the multivariate parameter space than the crude Monte Carlo sampling (MCS) and the variance reduction techniques such as Latin hypercube sampling (LHS) [14-17]. The ACO benefits from the multivariate uniformity property of the HSS. Therefore, the main advantage of the proposed algorithm is the high computational efficiency attained from the EACO algorithm compared to the conventional ACO algorithm.

The rest of this article is organized as follows: The Ant colony optimization algorithms for combinatorial, continuous and mixed variables are presented in the next section. In Section 3, Oracle penalty method to solve constrained optimization problems is introduced. The "Sampling Techniques" are introduced in Section 4. The efficient ant colony optimization algorithm proposed in this work is presented in Section 5. Following the benchmark problems presentation in Section 6, results and discussions of the benchmark problems are presented in Section 7. Finally, the concluding remarks are presented in the Section 8.

*Corresponding author: Urmila M Diwekar, Center for Uncertain Systems: Tools for Optimization & Management (CUSTOM), Vishwamitra Research Institute, Crystal Lake, USA, Tel:(630)-886-3047, Email: Urmila@vri-custom.org

Ant Colony Optimization

The ACO is a metaheuristic class of optimization algorithm inspired by the foraging behavior of real ants [6]. Natural ants randomly search food by exploring the area around their nest. If an ant locates a food source, while returning back to the nest, it lay down a chemical pheromone trail that marks its path. This pheromone trail will indirectly communicate with other members of the ant colony to follow the path. Over time, the pheromone will start to evaporate and therefore reduce the attraction of the path. The routes that are used frequently will have higher concentration of the pheromone trial and remain attractive. Thus, the shorter the route between the nest and food source imply short cycle time for the ants and these routes will have higher concentration of pheromone than the longer routes. Consequently, more ants are attracted by the shorter paths in the future. Finally, the shortest path will be discovered by the ant colony [6,8].

In ACO algorithms, artificial ants are stochastic candidate solution construction procedures that exploit a pheromone model and possibly available heuristic information of the mathematical model. The artificial pheromone trails (numeric values) are the sole means of communication among the artificial ants. Pheromone decay, a mechanism analogous to the evaporation of the pheromone trial of the real ant colony allows the artificial ants to forget the past history and focus on new promising search directions. Like the natural ants, by updating the pheromone values according to the information learned in the preceding iteration, the algorithmic procedure leads to very good and hopefully, a global optimal solution.

ACO for Discrete optimization problems

In a discrete optimization problem (shown in Eqn.1), the search domain of the problem is partitioned into a finite set of components, and the discrete optimization algorithm attempts to find the optimal combination or permutation of a finite set of elements from large and finite set of the search domain [17].

$$\min_{x} \quad f(y)$$
$$s.t. \quad h(y) = 0 \tag{1}$$
$$g(y) \le 0$$
$$y \in Y$$

The combinatorial search space Y is given as a set of discrete variables y_i where $i = 1, ..., NDIM$, with possible set of options $v_i^j \in NOP_i = \{v_i^1, v_i^2, ..., v_i^{NOP_i}\}$. If a solution vector $y \in Y$ takes a value for each decision variable that satisfies all the constraints, it is a feasible solution of the discrete optimization problem. A solution $y^* \in Y$ is called a global optimal if and only if; $f(y^*) \le f(y) \quad \forall y \in Y$. Therefore, solving a discrete optimization problem involves finding at least one $y^* \in Y$.

To solve a discrete optimization problem, artificial ants construct a solution by moving from one solution state to another in a sequential manner. Real ants walk to locate the shortest path by choosing a direction based on local pheromone concentrations and a stochastic decision policy. Likewise, the artificial ants construct solutions by moving through each decision variable guided by the value of the artificial pheromone trial and by making stochastic decisions at each state.

The important differences between the real and artificial ants are

- Artificial ants move sequentially through a finite set of decision variable set of options. Real ants move in a direction of the food by selecting from different passible paths, which are not finite.

- Real ants deposit and react to the concentration of the pheromone while walking. However, in artificial ants, sometimes the pheromone update is done only by some of the artificial ants, and often the pheromone updated is only after a complete solution is constructed.

- Artificial ants can introduce additional mechanisms to improve the solution such as adding local search strategies that does not exist in real ants.

The decision policy in choosing the solution component considers a trade-off between the pheromone intensity on a particular edge and the desirability of that edge with respect to the edge contribution on the objective function [4]. Taking these two properties of an edge into account, ACO algorithms effectively utilize the pheromone intensity which is based on the information learnt from the prior solution construction and the edge desirability. The decision policy is given by the transition probability function as shown in Eqn. 2.

$$\text{prob}_{ij}(it) = \frac{\tau_{ij}^{\alpha}(it)\eta_{ij}^{\beta}}{\sum_{il}\tau_{ij}^{\alpha}(it)\eta_{ij}^{\beta}} \tag{2}$$

where $prob_{ij}(it)$ is the probability of choosing edge ij as a solution component when an ant is at node i at iteration it, $\tau_{ij}(it)$ is the pheromone value associated with edge ij at iteration it, $\eta_{ij}(it)$ is the desirability of edge ij. α and β are parameters that control the relative importance of pheromone intensity and desirability, respectively. If $\alpha >> \beta$, the algorithm will make decisions based on mainly on the learned information, represented by the pheromone, and if $\alpha << \beta$, the algorithm will act as a greedy heuristic selecting mainly the cheapest edges, disregarding the learned information [6].

Evaluation of the pheromone value $\tau_{ij}(it)$ at the end of each of the iteration is the core of ACO algorithm. After each ant has constructed a solution (i.e., at each iteration) the pheromone value on each edge is updated. The goal of the pheromone update is to increase the pheromone values associated with good or promising solutions, and to decrease those that are associated with bad ones. The pheromone updating rule consists of two operations:

- The pheromone evaporation operation that reduces the current level of pheromone

- The pheromone additive operation: depends on the quality of the solutions generated at the iteration, pheromone is added to the edge of good solutions. The updating rule is given as follows [6] as shown in Eqn. 3.

$$\tau_{ij}(it + 1) = \rho\tau_{ij}(it) + \Delta\tau_{ij}(it) \tag{3}$$

Where ρ is the pheromone evaporation factor representing the pheromone decay ($0 \le \rho \le 1$) and $\Delta\tau_{ij}(it)$ is the pheromone addition for edge ij. The decay of the pheromone levels enables the colony to 'forget' poor edges and increases the probability of good edges being selected (i.e. the assumption behind this is that as the process continues in time, the algorithm learns to add pheromone only to good edges, implying that more recent information is better than older information). For $\rho \to 1$, only small amounts of pheromone are decayed between iterations and the convergence rate is slower. This is characterized by the high probability of finding the global optimal solution at the expense computational efficiency. Whereas for $\rho \to 0$ more pheromone is decayed resulting in a faster convergence. This trend leads to getting stuck at the local optimal solutions.

The pheromone addition operation is increasing the pheromone values associated with good or promising solutions and it is the main feature that dictates how an ACO algorithm utilizes its learned information. Typically, pheromone is only added to edges that have been selected, and the amount of pheromone added is proportional to the quality of the solution. In this way, solutions of higher quality receive higher amounts of pheromone [17]. The pheromone concentration added at iteration *it* is determined from Eqn. 4.

$$\Delta \tau_{ij}(it) = \sum_{k}^{nAnts} \Delta \tau_{ij}^{k}(it) \tag{4}$$

Where the pheromone concentration associated with ant k as function of the quality of the objective function is determined as shown below.

$$\Delta_{ij}(it) = \begin{cases} \overline{f(y)} & \textit{if } ij \in edge\ visited\ by\ ant\ k \\ & Otherwise \end{cases} \tag{5}$$

ACO algorithm to solve discrete optimization problems has five most popular variants which are the ant system (*AS*) [4], elitist ant system (*EAS*) [4], rank based ant system (*RAS*), ant colony system (*ACS*) and the *Max-Min* ant system (*MMAS*). For details of the ant algorithm variants, the book by Dorigo and Stutzle [6] can be viewed. These variants are derived from the AS algorithm and they follow similar solution construction procedure and pheromone evaporation procedure. The main differences among these variants are the way the pheromone update and pheromone trial management are performed.

ACO for Continuous domain optimization problem

Continuous variable optimization problems could be tackled with a discrete optimization algorithm only if the continuous ranges of the decision variables are discretized into finite sets [17]. The continuous ACO algorithm stores a set of K solutions in a solution archive which represents as the pheromone model of the combinatorial ACO algorithm. This solution archive is used to generate a probability distribution of the promising solutions over the search space. In this algorithm, the solution archive is first initialized and the algorithm iteratively finds the optimal solution by generating new solutions of size equal to the number of ants (nAnts). The new nAnts size solutions are then added to the K size solutions from the previous iteration. To propagate the updated K size solutions to the next iteration, the K + nAnts solutions are first sorted according to the quality of the objective function of each solution. Then, the solution archive is updated by keeping the best K solutions of the combined solutions and removing the worst nAnts size solutions. The details of the algorithm formulation can be found in Socha [17], but for the completion of this work the main features of the formulation are presented below.

Similar to the combinatorial ACO algorithm the solution construction of ants is accomplished in an incremental manner, *i.e* variable by variable. First, an ant chooses probabilistically one of the solutions in the solution archive as a solution construction guide. The probability of choosing solution *j* as guide is given by Eqn. 6:

$$prob_j = \frac{\omega_j}{\sum_{r=1}^{K} \omega_r} \tag{6}$$

where ω_j is the weight associated with solution *j*. The weight is determined using the Gaussian function [17] as follows:

$$\omega_j = \frac{1}{qK\sqrt{2\pi}} e^{\frac{-(j-1)^2}{2(qK)^2}} \tag{7}$$

where, *j* is rank of the solution in the archive, *q* is an algorithmic parameter and *K* is the size of the solution archive. The mean of the

Gaussian function is *1*, so that the best solution has highest weight [17]. The ant treats each problem variable *i* = 1, . . . *NDIM* separately. It takes value x_j^i of variable *i* of the jth solution guide and samples its neighborhood.

The algorithm for continuous domain is designed with the aim of obtaining multimodal one dimensional density function (PDF) from a set of probability density functions. Each PDF is obtained from the search experience and it is used to incrementally build a solution $x \in \mathfrak{R}^n$. Where, x is a vector of the continuous decision variables. To approximate a multimodal PDF [14] proposed a Gaussian kernel which is defined as a weighted sum of several one dimensional Gaussian functions $g_{ji}(x)$ as shown in Eqn. 8.

$$G_i(x) = \sum_{j=1}^{K} \omega_j g_j(x) = \sum_{j=1}^{K} \omega_j \frac{1}{\sigma_j \sqrt{2\pi}} e^{\frac{-(x-\mu_j)^2}{2\sigma_j^2}} \tag{8}$$

where $i \in \{1,2,...,K\}$ and $j \in \{1,2,...,K\}$ identify the optimization problem dimensionality and ranking of the solution in the archive. ω_j is vector of weights associated with the individual Gaussian function. μ_{ij} is the vector of means of the individual solution components. σ_{ij} is vector of the standard deviations. All these vectors have cardinality K.

Each row of the solution archive maintains the solution components $x_j \in \mathfrak{R}^n$ where x_j is vector of solution of row *j* of the solution archive ($x_j = x_{j1}, x_{j2},....,x_{jn}$), the objective function $f(x_j)$ and the weight associated with the solution (ω_j). The objective function values and the corresponding weight are stored in such a way that $f(x_1) < f(x_2),$ . . . , $< f(xj), \ldots , < f(x_K)$ and $\omega_1 > \omega_2, \ldots, > _j, \ldots, > \omega_K$. The solutions in the achieve T are, therefore, used to dynamically generate the PDFs involved in the Gaussian kernels. More specifically, in order to obtain the Gaussian kernel G_i, the three parameters (ω_i, μ_i, and σ_i) has to be determined. Thus, for each G_i, the values of the ith variable of the K solutions in the archive will be elements of vector μ_i, that is, $\mu_i = \mu_{1i}, \mu_{2i},..., \mu_j, ...\mu_K = x_{1i}, x_{2i},..., x_j, ...x_K$. On the other hand, each component of the standard deviation vector $\sigma_i = \sigma_{1i}, \sigma_{2i},...,\sigma_j, ...\sigma_K$ is determined as shown in Eqn. 9.

$$\sigma_j = \xi \sum_{z=1}^{K} \frac{|x_z - x_j|}{K-1} \tag{9}$$

where $j \in \{1,2,...,K\}$, $\xi > 0$ is a parameter that has similar effect to that of the pheromone evaporation parameter in combinatorial ACO algorithm. The higher the value of ξ leads to the lower convergence speed of the algorithm.

The solution archive update is performed following the three steps below.

- The newly generated solutions of size equal to the number of ants (*nAnts)* are added to the old solution archive.

- The combined solutions are sorted according to the quality of the objective functions.

- The first K best solutions are stored and the worst *nAnts* size solutions are removed.

ACO for Mixed-integer nonlinear programming problem

For optimization problem that includes continuous and discrete variables, the ACO algorithm uses a continuous-relaxation approach for *ordering* variables and similar to the combinatorial approach for *categorical* variables [17].

Ordering discrete variables: the algorithm works with indexes. In

similar fashion to that of the algorithm for continuous variables, the ACO algorithm generates values of the indexes for a new solution as real numbers. However, before the objective function is evaluated, a heuristic rule is applied for the continuous values that represent the discrete variables by round off the continuous values to the closest index, and the value at the new index is then used to evaluate the objective function.

Categorical discrete variables: in conventional ACO algorithm for combinatorial problems, solutions are constructed by choosing the solution component values from the finite search domain following the transition probability test that depend on the static pheromone value. However, because the solution archive replaces the pheromone model of the ACO algorithm, the transition probabilistic rule is modified to handle the solution archive.

Similar to the case of continuous variables, each ant constructs the discrete part of the solution incrementally. For each $i = 1, NT$ discrete variables, each ant chooses probabilistically one solution component sc_i from the finite set of solution component options available $v_i^j \in NOP_i = \{v_i^1, v_i^2, ..., v_i^{NOP_i}\}$. The probability of choosing the jth value is given by Eqn. 10.

$$prob_i^j = \frac{\omega_j}{\sum\limits_{r=1}^{\kappa} \omega_r} \qquad (10)$$

where ω_j is the weight associated with the jth available option. It is calculated based on the weights (ω) and some additional parameters as shown below:

$$\omega_j = \frac{\omega_r}{u_i^r} + \frac{q}{\eta} \qquad (11)$$

where the weight ω_{jr} is calculated according to Eqn. 7, jr is the index of the highest quality solution in the solution archive that uses value v_i^j for the ith variable. u_i^r is the number of solutions in the archive that use value v_i^j for the ith variable. Therefore, the more popular value v_i^j in the archive, the lower is its final weight. The second component is a fixed value (i.e., it does not depend on the value v_i^j selected): η is the number of values of v_i^j from the sc_i available that are unused by the solutions in the archive, and q is the same algorithm parameter that was used in Eqn. 7. Some of the available categorical values v_i^j may be unused for a given i^{th} decision variable in all the solutions of the archive. Hence, their initial weight is zero. The second component is added in order to enhance exploration and prevent premature convergence.

Oracle Penalty Method

The problem given in Eqn. 1 is a constrained optimization problem, which can be solved using a penalty method. To solve such problem, first the problem is transformed into unconstrained optimization problem by transforming the original objective function into a penalty function. In most cases, the penalty function is given as a weighted sum of the original objective function and the constraint violations [15,22]. In such a way, the penalty function serves as an objective function. The main advantage of the penalty method is its simplicity. However, the simple penalty methods often perform very poorly on challenging constrained optimization problems, while the more advanced methods need additional parameters thus require an additional tuning of these parameters. An oracle penalty method proposed by Kalagnanam [23] is a simple to implement and general method capable of handling simple and challenging constrained optimization problems. In oracle penalty method, the objective function is first transformed into additional

equality constraint $h_0(x) = f(x) - \Omega$; where Ω is a parameter called oracle. The objective function becomes redundant in the transformed problem definition. Therefore, it is a constant zero function $\widetilde{f}(x)$ and the transformed optimization problem is presented as shown below:

$$\min_{x} \widetilde{f}(x)$$
$$s.t. \quad h_0(x) = f(x) - \Omega \qquad (12)$$
$$h(x) = 0$$
$$g(x) \leq 0$$
$$x \in \Re^n$$

By transforming the objective function into an equality constraint, then minimizing the new constraint $h_0(x)$ and the residual of the original constraints $h(x)$ and $g(x)$ becomes directly comparable. The penalty function balances its penalty weight between the transformed objective function and the violation of the original constraints. The implementation of the oracle penalty function can be found in Appendix A.

Sampling Techniques

Sampling is a statistical procedure that involves selecting a finite number of observations, states, or individuals from a population of interest. A sample is assumed to be representative of the whole population to which it belongs. Instead of evaluating all the members of the population, which would be time consuming and costly, sampling techniques are used to infer some knowledge about the population. Sampling techniques are used in a wide range of science and engineering applications. The most commonly used sampling technique is a probabilistic sampling technique which is based on a pseudo-random number generator called Monte Carlo Sampling (MCS). This sampling technique has probabilistic error bounds and large sample sizes are needed to achieve the desired accuracy. Variance reduction techniques have been applied to circumvent the disadvantages of Monte Carlo sampling [18,19,24].

Monte Carlo Sampling

One of the simplest and most widely used methods for sampling is the Monte Carlo sampling. Monte Carlo method is a numerical method that provides approximate solution to a variety of physical and mathematical problems by random sampling. In crude Monte Carlo approach, a value is drawn at random from the probability distribution for each input, and the corresponding output value is computed. The entire process is repeated n times producing n corresponding output values. These output values constitute a random sample from the probability distribution over the output induced by the probability distributions over the inputs. One advantage of this approach is that the precision of the output distribution may be estimated using standard statistical techniques. On average, the error ϵ of approximation is of the order $O(N^{1/2})$. One important feature of this sampling technique is that the error bound is not dependent on the dimension. However, this error bound is probabilistic, which means that there is never any guarantee that the expected accuracy will be achieved. The pseudorandom number generator produces samples that may be clustered in certain regions of the population and does not produce uniform samples. Therefore, in order to reach high accuracy, larger sample sizes are needed, which adversely affects the computational efficiency [18,19].

Hammersley Sequence Sampling

To improve the efficiency of Monte Carlo simulations and overcome the disadvantages, eg, probabilistic error bounds, variance reduction

techniques were proposed. The sampling approaches for variance reduction that are used more frequently in the chemical engineering applications are importance sampling, Latin Hypercube Sampling (LHS), Descriptive Sampling, and HSS. The HSS belongs to the group of quasi-Monte Carlo methods [19].
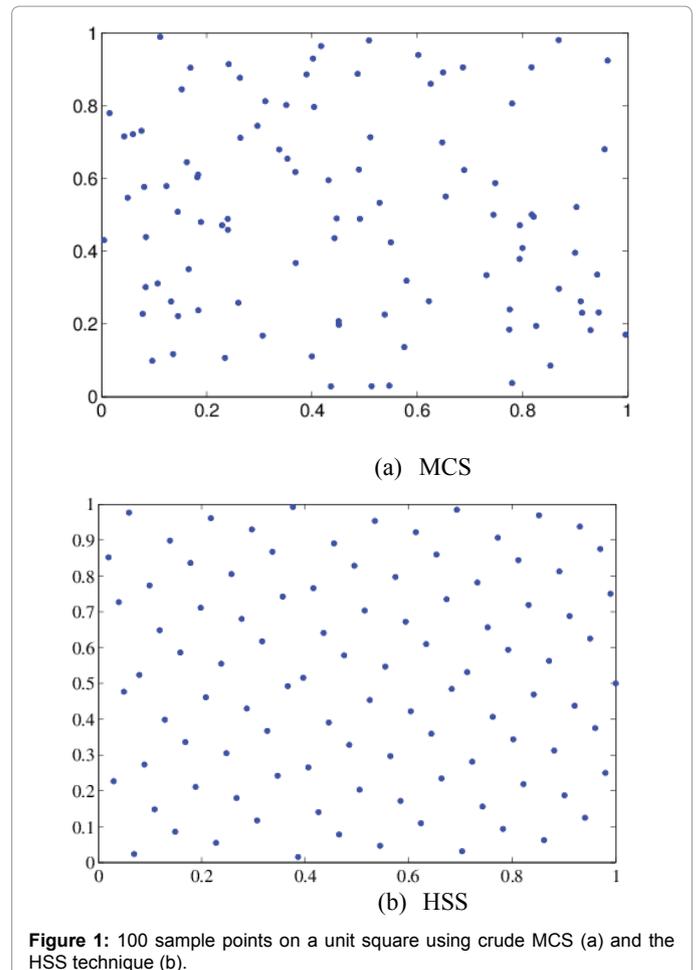
LHS is one form of stratified sampling that can yield more precise estimates of the distribution function, and therefore reduce the number of samples required to improve computational efficiency. This sampling method was used in ACO algorithm to initialize the solution archive for continuous variable optimization problem [25]. LHS was designed to improve the uniformity properties of Monte Carlo methods. It was shown that the error of approximating a distribution by finite sample depends on the equi-distribution properties of the sample used, but the relationship between successive points in a sample, the randomness or the independence is not critical [18,19]. The main drawback of this stratification scheme is that it is uniform in one dimension (1D) but it does not provide the uniformity property in k dimensions.

HSS is an efficient sampling technique developed by Diwekar research group [18,19,24]. It is based on quasi-random number generator. HSS uses Hammersley points to uniformly sample a unit hypercube and inverts these points over the joint cumulative probability distribution to provide a sample set for the variables of interest. HSS technique uses an optimal design scheme for placing n points on a k-dimensional hypercube. This scheme ensures that the sample is more representative of the population showing uniformity properties in multi-dimensions [18,19,24]. A qualitative picture of the uniformity properties of the MCS and HSS techniques on a unit square is presented in Figure 1. In the figure, it is clearly shown that samples generated by the HSS technique achieve better uniformity in the parameter space and hence results better computational efficiency.

Efficient Ant Colony Optimization (EACO) Algorithm

The proposed EACO algorithm improves the conventional ACO algorithm for combinatorial, continuous and mixed variable optimization problems by introducing the HSS technique. The initial solution archive diversity for continuous and mixed-variable optimization problems plays an important role in the performance of ACO algorithm. The uniformity property of the HSS technique is used in this step to avoid initial solution archive clustered in a small region of the potential solution space. Moreover, ACO algorithm is a probabilistic method, several random probability functions are involved in the algorithm procedure. Examples: in combinatorial ACO algorithm, the transition probability that help to choose the next solution component and for continuous and mixed-variable optimization problems, the probability of choosing ant guide from the solution archive. The distribution of the random numbers generated for the acceptance probability of a solution component and ant guide affects the performance of the ACO algorithm. At this stage, the multidimensional uniformity property of HSS is exploited to choose these random numbers.

The EACO algorithm uses as termination criteria maximum number of solution construction steps (*MaxIter*), and the tolerance (ϵ) that is the relative difference between solutions found in two consecutive iterations is less than or equal to a parameter ϵ for a set of consecutive number of iterations, *ICON*. The EACO algorithm is given in Table 1 and Figure 2.

(a) MCS

(b) HSS

Figure 1: 100 sample points on a unit square using crude MCS (a) and the HSS technique (b).

T is solution archive and K is size of T. nAnts is the number of ants. NC, NO, NT and NDIM are the number of continuous, ordinal, categorical and the total number of decision variables, respectively. NOPT is the number of options in categorical variables.

The best features of the EACO algorithm are

- The ability to exploit the multi-dimensional uniformity property of the Hammersley sequence sampling (HSS)

- The high computational efficiency over the conventional ACO algorithm

- The ability to solve combinatorial, continuous and mixed-integer optimization problems

- The ability to solve large scale convex and nonconvex optimization problems

Benchmark problems

The main goal of the benchmark problems is to establish the performance comparison between the proposed EACO algorithm and the conventional ACO algorithm. For combinatorial, continuous and mixed-variable optimization problems, 2, 3 and 4 benchmark problems, respectively, are used for the experimentation. In all of the benchmark problems, we did the experiments for 5, 10, 15, and 20 dimensional variables. 10 experiments for each combination of the benchmark and the dimensions were performed. Moreover, for the combinatorial ant algorithm,

there are five most popular algorithm variants [6]. The performance of each variant is tested using the EACO algorithm. The performance measure used to evaluate the algorithms was based on the number of iterations that the algorithm needed to reach the global optimal of the test functions for the same algorithmic parameter setting and terminating criterion. The benchmark problems are explained below and are given in Table 2.

Start program
• Set K, $nAnts$, NC, NO, NT, $NOPT$, , q, ξ and termination criteria
• Initialize solution archive $T(K, NC + NO)$ using HSS
• Initialize solution archive $T(K, NT)$ randomly from the possible options
• Combine and evaluate the objective function of the K solutions $T(K, NDIM)$
• Rank solutions based on the quality of the objective function ($T = rank(S_1 \dots S_K)$)
• For categorical optimization problems, introduce multidimensional random number generated with HSS ($IterMax \times nAnts \times NT$, $NOPT$)
While Termination criterion is not satisfied
• Generate solutions equivalent to the number of ants ($nAnts$)
For all # nAnts
• Incremental solution construction
For all # NDIM
○ Probabilistically construct continuous decision variables
○ Probabilistically construct ordinal decision variables
○ Probabilistically construct categorical decision variables
End for # NDIM
• Store and evaluate the objective function of the newly generated solutions
End for # nAnts
• Combine, rank and select the best K solutions, $T = Best(rank(S_1 \dots S_K \dots S_{K+nAnts}), K)$
• Update solution
End while
End program

Table 1: EACO Algorithm for continuous, combinatorial and mixed variable problem.

Figure 2: EACO algorithm for continuous, combinatorial and mixed variable problems.

	Function	Formula	Range
Combinatorial Optimization Problems			
1	Travelling Salesman Problem		
2	EX. II	$f_{BC}(y) = \sum_{i=1}^{y_1}(y_1 - 3)^2 + (y_2^i - 3)^2 + (y_3^i - 3)^2$	$[1,5]^{NC}$
Continuous Combinatorial Optimization Problems			
3	Parabolic (PR_{CV})	$f_{PR}(x) = \sum_{i-1}^{NC} x_i^2$	$[-3,3]^{NC}$
4	Ellipsoid (EL_{CV})	$f_{EL}(x) = \sum_{i=1}^{NC} 5^{\frac{i-1}{n-1}} x_i^2$	$[-3,3]^{NC}$
5	Cigar (CG_{CV})	$f_{CG}(x) = x_1^2 + 10^4 \sum_{i=2}^{NC} x_i^2$	$[-3,3]^{NC}$
Mixed-variable Combinatorial Optimization Problems			
6	Parabolic (PR_{MV})	$f_{PR}(x,y) = \sum_{i=1}^{NC} x_i^2 + \sum_{i=1}^{ND} y_i^2$	$[-3,3]^{NM}$
7	Ellipsoid (EL_{MV})	$f_{EL}(x,y) = \sum_{i=1}^{NC} 5^{\frac{i-1}{n-1}} x_i^2 + \sum_{i=1}^{ND} 5^{\frac{i-1}{n-1}} y_i^2$	$[-3,3]^{NM}$
8	Cigar (CG_{MV})	$f_{CG}(x,y) = x_1^2 + 10^4 \sum_{i=2}^{NC} x_i^2 + y_1^2 + 10^4 \sum_{i=2}^{ND} y_i^2$	$[-3,3]^{NM}$
9	EX. III	$f(x,y) = \sum_{i=1}^{NC}(x_i - i/NC)^2 + \sum_{i=1}^{ND} y_i^2 + \prod_{i=1}^{ND} \cos(4\pi y_i)$	$[-3,3]^{NM}$

Table 2: Benchmarks problems.

- Traveling Salesman Problem (TSP): Given the number of cities to be visited and the distances between each pair of cities, find the shortest possible route that lead to visiting each city exactly once and returns back to the origin city. The TSP is an NP-hard combinatorial optimization problem.

- A TSP with known global optimal solution is also retrieved from the TSPLIB benchmark library accessible at http://www.iwr.uni-heidelberg.de/groups/comopt/software/TSPLIB95/. For each variant of the combinatorial ACO optimization algorithm, a 16 Odyssey geographical locations (ulysses16.tsp) with global optimal solution of 6859 is used to perform the computational efficiency comparison between the EACO and the conventional ACO algorithm.

- A pure combinatorial problem, example II of Kim and Diwekar [26]. This problem has one global minimum 0 when all y_1; y_{2i}; and y_{3i} are equal to 3.

- Three test functions are used for the continuous optimization problems. The parabolic function is from Kim and Diwekar example I [26]. This function is a multidimensional parabolic function that has one global optimum at 0 for all decision variables equal to 0. The second and third test functions, which are the ellipsoid and cigar functions, are from Socha [17].

- The three test functions used for the continuous optimization problem are further modified to represent mixed variable optimization problems as shown in Table 2.

- Example III from Kim and Diwekar [26] which is an MINLP problem that has one global minimum -1 is further used as an MINLP benchmark problem.

- Case studies on real world optimization problems solved by the proposed EACO algorithm can be viewed in references [28,29].

Results and Discussions

Benchmark Problems

This section presents a comparative study between the EACO and the conventional ACO algorithm that use MCS for initializing the solution archive and generating the multidimensional random number to optimize the 9 benchmark problems presented in Table 2. The proposed EACO algorithm effectively solves discrete, continuous, and mixed-integer optimization problem benchmarks. The algorithm terminates when it reaches maximum number of solution construction steps, *MaxIter*, or if the tolerance *(ϵ)* that is the relative difference between solutions found in two consecutive iterations is less than or equal to a parameter ϵ for a set of consecutive number of iterations, *ICON*.

In order to make unbiased comparison between the EACO and the conventional ACO algorithms, we use the same parameters to conduct the experiments. The results are based on 10 independent runs on each benchmark problem. Each run differs in the seed to generate the random numbers and the solution archive initialization. The parameters used to tackle the problems are selected after performing a number of experimentations using different combination of the parameters. The experiments to get the parameter settings are performed using the conventional ACO algorithm. The summary of the parameters that we use for the algorithms are presented in Table 3. As shown in Table 3, the parameter setting differs from problem to problem. The archive size *K* and the number of ants *nAnts* for 5 dimensional variables are 50 and 2,

respectively.

x and y are vector of continuous and discrete variables, respectively. NC, ND, and NM are number of continuous, discrete and mixed variable, respectively.

EX. II and III refers to examples from Diwekar [26].

Combinatorial EACO algorithms: The results of the discrete benchmark problems are presented in Tables 4-5. As shown in the tables, for each of the 40 (3×10 + 10) pair of the experimental runs of the benchmark problems, the EACO algorithm shows a better computational efficiency than the conventional ACO algorithm. The iteration improvement ranges from 24.2% of the multistage turbine problem of example II from Kim and Diwekar [26] to 33% of the travelling salesman problem with 10 and 20 numbers of cities to be visited. For the discrete optimization problems, by keeping the diversity of the multidimensional random numbers, the EACO algorithm produces more uniform operation to guide ants in making decisions of accepting or rejecting the current transition probability that lead to the finding of a new solution component. The results in Table 4 and 5 show that the EACO benefits from the uniformity property and diversified multidimensional random number generations.

Comparison of variants of the combinatorial EACO algorithms: To compare the performance of the EACO algorithm with the variants of the combinatorial ACO algorithms, the following algorithm parameter settings are used. Maximum number of tour construction steps (*MaxIter*) of 1000 and absolute error between the global optimal and the best so-far solution ϵ = 1E-6 are used as stopping criteria. Moreover, the parameter settings dependent on the algorithm variant are given as follows. *AS*: α = 1; β = 3; ρ = 0.5, and nAnts = n, that is number of ants equal to the number of the geometrical locations (n). *EAS*: α = 1; β = 5; ρ = 0.5, e = n, and nAnts = n. *RAS*: α = 1; β = 4; ρ = 0.6, w = 10, and nAnts = n. *ACS*: α = 1; β = 3; ρ = 0.9, ξ = 0.1, and nAnts = n. *MMAS* : α = 1; β = 5; ρ = 0.98, p_{bst} = 0.05; and nAnts = n. The pheromone update for *MMAS* is performed using only the iteration-best ant.

The results are based on 25 independent runs of the Ulysses16.tsp problem using the five ant colony algorithm variants. The same parameter settings are used for the same ant colony algorithm variant except the random number generated using the Monte Carlo sampling (ACO) and the multidimensional quasi-uniform HSS sampling methods (EACO). As shown in Figure 3, for ant system (AS), rank based ant

Function	α	β	ρ	nAnts	K	q	ε
TSP	1	4	0.5	nCity	-	-	1E-4
Ex. II [27]	-	-	0.5	20	150	0.001	1E-6
Other test functions	-	-	0.5	5	150	0.001	1E-5

Table 3: Summary of the parameters used for the benchmark problems.

	HSS		MCS		Improve [%]	
NCITY	Iter	Length	Iter	Length	[%] Iter	[%] Length
10	6	123	9	123	33	0
20	14	164	21	164	33	0
40	31	209	45	207	31	-1

Table 4: Combinatorial: Travelling Salesman Problem for 10, 20 and 40 cities.

			EACO	ACO	Improve
ND	**Runs**	**Glob. Opt.**	Iter	Iter	[%]
11	10	0	16	22	24.2

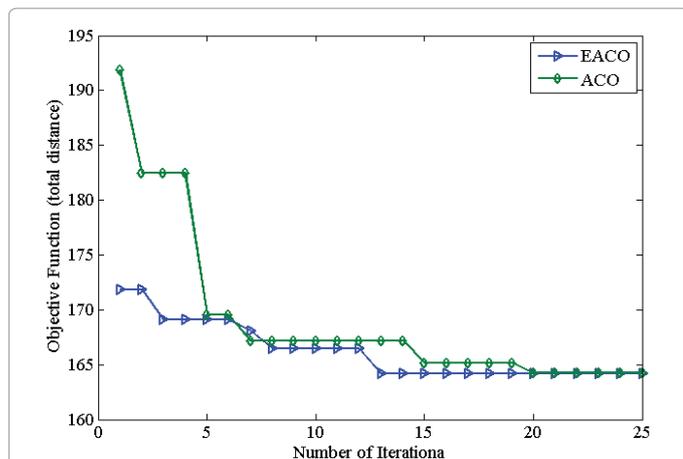**Table 5:** Combinatorial: Example II of Diwekar [26].

Figure 3: Number of runs that find the global optimal solution of the ulysses16. tsp (6859) for each variant the combinatorial ACO and EACO.

system (RAS), and ant colony system (ACS), the EACO form of these algorithms perform better than the conventional ACO algorithm form of these variants. That is, the numbers of runs that find the global optimal solution using the EACO form of AS, RAS and ACCS are higher than that of the conventional ACO algorithm forms of the same variants. In the case of elitist ant system, both methods find the global optimal solutions for the same number of runs (17/25). However, for *Max-Min* ant system, the conventional ACO algorithm find global optimal solution in more runs (23/25) than the EACO algorithm (22/25). This could be attributed to the fact that the small problem size and already *MMAS* is an efficient method of all the ant variants.

Continuous EACO algorithm: The results of the continuous benchmark problems are presented in Tables 6-8. Among the 120 (3 ×4×10) pair of the experimental runs of the benchmark problems, only in one instant (Table 6 of parabolic function with 5 decision variables) the two algorithms have same performance. For 99.2% of the experimental runs of the continuous functions the performance of

NDIM	Glob. Opt	EACO Iter	ACO Iter	Improve [%]
5	0	90	90	0
10	0	151	165	8
15	0	166	187	11
20	0	250	270	7

Table 6: Continuous: Parabolic function for 5, 10, 15 and 20 decision variables.

NDIM	Glob. Opt	EACO Iter	ACO Iter	Improve [%]
5	0	32	37	14
10	0	169	183	8
15	0	198	233	15
20	0	294	341	14

Table 7: Continuous: Ellipsoid function for 5, 10, 15 and 20 decision variables.

NDIM	Glob. Opt	EACO Iter	ACO Iter	Improve [%]
5	0	85	105	19
10	0	304	325	6
15	0	326	426	23.5
20	0	493	536	8

Table 8: Continuous: Cigar function for 5, 10, 15 and 20 decision variables.

the EACO algorithm is better than the conventional ACO algorithm. Because of the multidimensional uniformity property of HSS, the EACO algorithm needed less iteration than using the MCS to find the global optimal solutions. The computational efficiency improvement ranges from 7% for the parabolic function with 20 decision variables (Table 6) to 23.5% of the cigar function with 15 decision variables (Table 8).

MINLP EACO algorithm: The results of the mixed-integer benchmark problems are presented in Tables 9-12. Among the 160 (4 ×4×10) pair of the experimental runs of the benchmark problems, all of the 160 pair experimental runs show that the performance of the EACO algorithm is better than the conventional ACO algorithm. The computational efficiency improvement ranges from 3% for the mixed variable cigar function with 20 decision variables (see Table 11) to 71% of the same function with 5 decision variables. Moreover, the tables also show that for the mixed variable optimization problems, the improvement on the computational efficiency of EACO is more pronounced.

Convergence Trajectory: Furthermore, the convergence paths of the EACO and the conventional ACO from each type of the benchmark problems are presented in Figures 4 - 6. The figures show the trajectories of the objective function value for combinatorial (Figure 4), continuous (Figure 5) and mixed variable (Figure 6) of the benchmark problems as function of the number of iterations to find the optimal solution. As shown in the figures, EACO algorithm found the global optimal solutions with 14, 196 and 326 iterations,

NDIM	Glob. Opt	EACO Iter	ACO Iter	Improve [%]
5	0	77	85	9
10	0	100	108	7
15	0	118	135	13
20	0	166	234	29

Table 9: Mixed variable: Parabolic function for 5, 10, 15 and 20 decision variables.

NDIM	Glob. Opt	EACO Iter	ACO Iter	Improve [%]
5	0	15	43	65
10	0	95	117	19
15	0	156	177	12
20	0	188	266	29

Table 10: Mixed variable: Ellipsoid function for 5, 10, 15 and 20 decision variables.

NDIM	Glob. Opt	EACO Iter	ACO Iter	Improve [%]
5	0	26	89	71
10	0	156	189	17
15	0	245	273	10
20	0	337	348	3

Table 11: Mixed variable: Cigar function for 5, 10, 15 and 20 decision variables.

NDIM	Glob. Opt	EACO Iter	ACO Iter	Improve [%]
5	-1	60	65	8
10	-1	112	128	13
15	-1	180	197	9
20	-1	232	268	13

Table 12: Mixed variable: Example III of Diwekar UM for 5, 10, 15 and 20 decision variables.
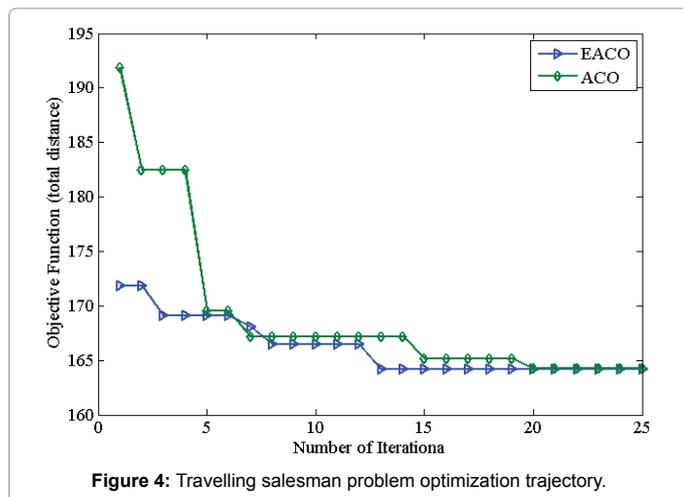
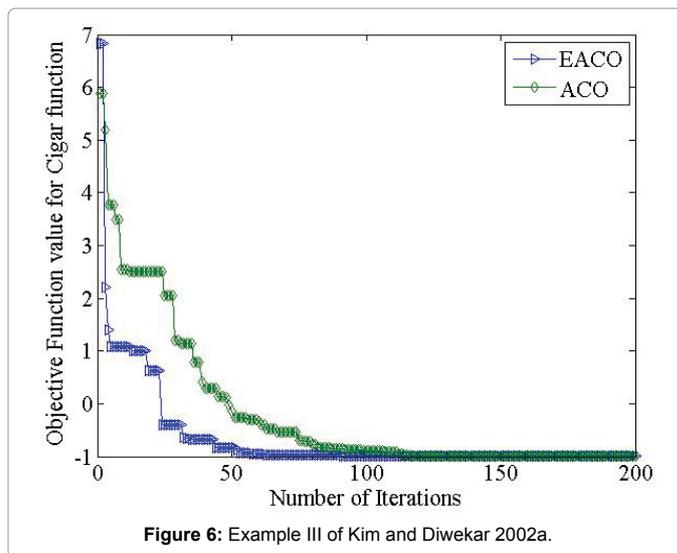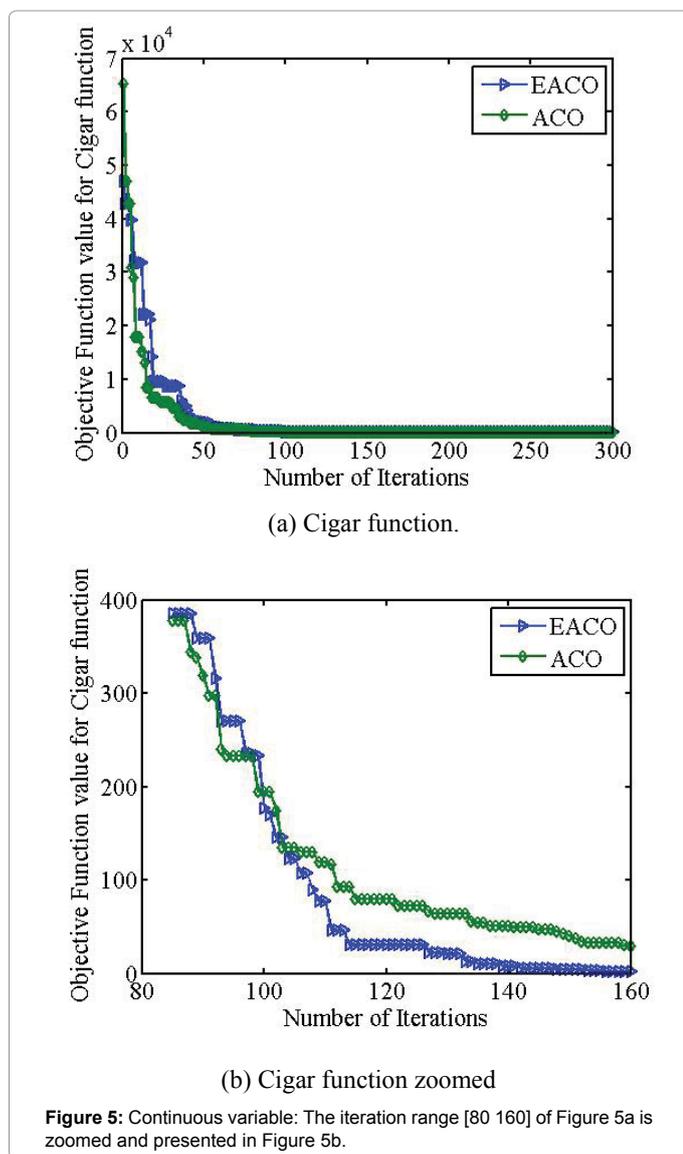**Figure 4:** Travelling salesman problem optimization trajectory.

Figure 6: Example III of Kim and Diwekar 2002a.

(a) Cigar function.

(b) Cigar function zoomed

Figure 5: Continuous variable: The iteration range [80 160] of Figure 5a is zoomed and presented in Figure 5b.

respectively. While the conventional ACO that uses MCS needs hogher number of iteration 21, 226, and 426 to reach the same optimal solutions. All the above observations prove that EACO algorithm benefits from the uniformity property of the HSS, by producing more uniform and diverse samples.

Conclusions

The ACO algorithm is a simple to implement yet a powerful and effective optimization framework for handling discrete, continuous and mixed-variable optimization problems. In this work, we proposed EACO as an alternative to the conventional ACO algorithm and the gradient based algorithms for optimization of large scale problems. ACO algorithm is probabilistic optimization strategy and the performance of the algorithm depends on the sampling strategies implemented in the algorithm. EACO algorithm is developed based on efficient sampling technique that keeps the diversity and the multidimensional uniformity property of samples. The capabilities of the proposed method are illustrated through 9 benchmark problems. The results show that the computational efficiency of the conventional ACO is improved in more than 99% of the experiments and the computational efficiency improvement ranges from 3% to 71%. Moreover, the tests on the five combinatorial ant colony algorithm variants show that the probability of finding the global optimal solution using the EACO algorithm is higher than that of the conventional ant colony algorithms.

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