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Submitted on 5 Mar 2015

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Eprints ID: 12657

To link to this article: DOI: 10.1016/j.future.2013.07.010
URL: http://dx.doi.org/10.1016/j.future.2013.07.010


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Exploiting performance counters to predict and improve energy performance of HPC systems

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HIGHLIGHTS
• We present two generic approaches for tracking high performance computing systems’ behaviour.
• We demonstrate that energy performance can be improved without a priori knowledge of application.
• We demonstrate that HPC systems can benefit from more than CPU frequency scaling.
• We introduce an approach to estimating the energy consumption of HPC applications.

ABSTRACT
Hardware monitoring through performance counters is available on almost all modern processors. Although these counters are originally designed for performance tuning, they have also been used for evaluating power consumption. We propose two approaches for modelling and understanding the behaviour of high performance computing (HPC) systems relying on hardware monitoring counters. We evaluate the effectiveness of our system modelling approach considering both optimizing the energy usage of HPC systems and predicting HPC applications’ energy consumption as target objectives. Although hardware monitoring counters are used for modelling the system, other methods – including partial phase recognition and cross platform energy prediction – are used for energy optimization and prediction. Experimental results for energy prediction demonstrate that we can accurately predict the peak energy consumption of an application on a target platform; whereas, results for energy optimization indicate that with no a priori knowledge of workloads sharing the platform we can save up to 24% of the overall HPC system’s energy consumption under benchmarks and real-life workloads.

1. Introduction

The increasing need for performance for the many computational problems in science and engineering has taken high performance computing (HPC) systems to the set of indispensable tools in modern industries and scientific research. HPC systems deliver tremendous raw performance for solving real-life problems. These problems include: turbulence, combustion, genomics, astrophysics, geosciences, molecular dynamics, homeland security, imaging and biomedicine, etc. HPC systems consume a large amount of electrical power, almost all of which is converted in heat requiring cooling. For example, Tianhe-1A consumes 4.04 MW of electricity [1]; a simple calculation at $0.10/KWh yields a powering and cooling cost of about $3.5 million per year, which is significant. Beyond the operating cost, some data-centres are also being limited by the peak power that electric facilities can provide. Consequently, it is necessary to reduce their power consumption; however, this must not result in significant performance degradation. The term significant performance degradation is a relative term; nevertheless, up to 10% performance degradation is often acceptable. Note: throughout this paper, (i) an HPC system is considered as a set of computing and storage nodes excluding network equipments such as routers and switches; whereas, (ii) the term “system” designates a single node of the HPC system.

The importance of power efficiency in HPC systems has attracted enormous attention from both the industrial and the research communities. This is evidenced by the multitude of techniques aiming to understand and reduce the energy consumption of HPC systems. These techniques generally break into hardware and software approaches. Hardware approaches led by manufac-
Embedding hardware event counters of modern microprocessors, or simply hardware performance/monitoring counters, monitor the occurrence of hardware events in a microprocessor with almost no performance penalty. These counters include: the number of cycles, instructions, cache references and hits/misses, main memory writes/backs and references, and branch miss-predictions counts. Although originally designed for performance tuning/adaptation, hardware performance counters have been used in the past for estimating the power usage (or the energy consumption) of both HPC and desktop applications.

Performance tuning mechanisms rely on hardware performance counters along with application's specific metrics (Message Passing Interface – MPI – calls in an MPI program for example) to adapt the system accordingly. This can be accomplished by inserting specific code segments in the source program (often referred to as code instrumentation) or by tracking application’s specific routine calls at runtime.

Adapting the processor’s frequency to meet workloads’ demands is the commonly used hardware adaptation approach for reducing energy consumption. This is eased by the use of Dynamic Voltage and Frequency Scaling (DVFS) [2,3] technology available on modern processors. A fundamental requirement for such adaptation is defining how DVFS control should be performed. In other words, the application designer must decide when and for how long the processor should be kept in a given performance state (P-state). In the performance optimization jargon, a specific part of the program which is defined to run at a given performance state is referred to as area or phase. Although these approaches for adapting HPC systems are efficient, they may fail for two main reasons:

1. HPC systems are often shared (the whole infrastructure is not dedicated to a single workload) by multiple workloads each having its own characteristics, in which case optimizing the energy performance considering some applications is likely to impact the performance of others.
2. In spite of the fact that HPC codes are actively maintained their increasing complexity makes code instrumentation impractical and sometimes require extensive knowledge: a platform provider can dedicate some engineers to “code instrumentation” or ask the programmer to write programs with these constraints in mind, which not only being unacceptable is not likely to happen.

An effective way to overcome these limitations is to optimize energy performance of the HPC system from the infrastructure stand point. This implies understanding the behaviour of the HPC system rather than that of individual applications sharing the platform.

In this paper, we introduce two complementary general purpose approaches for modelling and understanding the runtime behaviour of HPC systems. These approaches rely upon hardware performance counters and break into on-line and off-line depending on their use. Information gathered from the off-line approach can be of benefit to its on-line counterpart, which makes them complementary. The off-line approach which we refer to as “DNA-like description” of the system attempts to depict the HPC system as a graph in which each state describes its behaviour (that of the HPC system) over a time interval [4].

The on-line approach which we refer to as “Execution vectors based system behaviour tracking” detects and characterizes the runtime behaviour of the HPC system at runtime [5]. To accomplish this, phase changes are detected in the execution pattern of individual nodes in the HPC system; detected phases are characterized afterwards. Given that HPC workloads generally fall into compute intensive, memory intensive, and communication intensive, we define three types of behaviours including: compute intensive, memory intensive, and communication intensive. Communication intensive behaviour can further be divided into network transmit and receive.

To demonstrate the effectiveness of HPC system’s behaviour tracking approaches just mentioned, we explore several use cases. These use cases show how the energy performance of an application can be predicted using the DNA-like description on one the hand and how energy performance of an HPC system can be improved through behavioural characterization on the other hand. Our work differs from the state of the art in many ways: we address the power/energy consumption issue of HPC systems considering all HPC subsystems (processor, memory, disk and network).

In addition, our system modelling approaches does not rely on any application specific metric, i.e., our approaches do not require any a priori knowledge of applications running on the system.

The major contributions of this work are the following:

1. We present two different (on-line and off-line) approaches enabling a fine-grained control of HPC systems and a better understanding of application’s energy performance. Their strength resides in the fact that they do not need any a priori knowledge of applications/workloads sharing the HPC infrastructure.
2. We present an approach for optimizing/improving energy performance of HPC systems considering HPC subsystems, including the processor, memory, disk and network.
3. We introduce the concepts of cross platform energy prediction and partial phase recognition. Cross platform energy prediction can help choosing the appropriate execution platform for an application; whereas partial phase recognition is an alternative to phase prediction, it has the advantage of not being application or architecture specific.

The remainder of this article is organized as follows. Section 2 gives account of previous work. Our system’s modelling approaches are presented in Section 3. Section 4 presents several use cases of our modelling approaches. Implementation of the two use cases along with experimental results is discussed in Section 5. Finally, Section 6 concludes and gives future directions.

2. Related work

A large body of work investigates the use of hardware performance counters for modelling the power consumption of applications ranging form desktop to HPC applications. In this section, we first detail work using hardware performance counters for estimating the power or energy consumption of individual applications. We next present several work addressing HPC systems’ energy performance and their limitations.

2.1. Power/energy estimation using hardware performance counters

In studies such as [6–12] efforts have been devoted to model or estimate the power usage of individual applications or workloads. These studies monitor the use of system’s component (in particular the processor and memory) during the workload execution via hardware performance counters and correlate them with the power consumed by the system when running that workload to derive a power model. Kadajif et al. [7] propose a model for estimating the energy consumption of the UltraSPARC CPU [13]. Authors estimate the UltraSPARC CPU memory energy consumption considering the following performance counters: Data cache read hits, Data cache read references, Data cache write hits, Data cache...
write references, Instructions cache hits, Instructions cache references, Extended cache misses with writebacks. They claim their energy model to be 2.4% accurate as compared to circuit level simulation. Similarly, Contreras et al. [8] present a first order linear power estimation model that uses hardware performance counters to estimate run-time CPU and memory power consumption of the Intel PXA255 [14]. According to the authors, the proposed model exhibits an average estimation error of 4%. During their analyses, authors considered events listed in the first column of Table 1 for the CPU and those listed in the second column of that very table for the memory.

Energy consumption of the high-performance processor AMD Phenom is estimated in [9]. In their work, authors categorize AMD Phenom performance counters into four buckets: FP Units, Memory, Stalls, and Instruction Retired and consider performance events which express best their power consumption. These performance events include: L2_cache_miss:all, Retire_uops, Retire_mmx_and_fp_instruction:all, and Dispatch_stalls.

More recently, Da Costa et al. [6] have presented a methodology of measurement of the energy consumption of a single process application running on a standard PC. They defined a set of per process and system-wide variables to demonstrate their accuracy in measuring the energy consumption of a given process using multivariate regression.

To summarize, the above researches tell us that performance counters can accurately estimate the power usage of an application, however it is worthwhile to mention that the accuracy of a power/energy model depends on the workload. In other words, a power model designed for estimating the power consumption of a compute-bound workload may not fit well with a memory-bound workload. This is obvious for communication intensive workloads.

2.2. Energy reduction approaches for HPC systems

In the past few years, HPC systems have witnessed the emergence of energy consumption reduction techniques from the hardware level to the software level. At the hardware level, the majority of Information Technology (IT) equipment vendors works either from bottom up, by using the more efficient components in their equipments, and/or by providing their equipments with technologies that can be leveraged to reduce energy consumption of HPC subsystems – such as processor, network, memory, and I/O – during their operation. For example, the majority of modern processors is provided with Dynamic Resource Sleeping (DRS) which makes components hibernate to save energy and then wakes them on demand. Although major progress has been made, improvements in hardware solutions to energy reduction problem have been slow, due to the high cost of designing equipments with energy-saving technologies and the increasing demand of raw performance. Our work takes advantage of hardware technologies to reduce HPC systems’ energy consumption.

Unlike hardware approaches, software solutions for reducing HPC systems’ energy usage have received extensive attention over time. Rountree et al. [15] use node imbalance to reduce the overall energy consumption of a parallel application in an HPC system. They track successive MPI communication calls to divide the application into tasks composed of a communication portion and a computation portion. A slack occurs when a processor is waiting for data to arrive during the execution of a task. This leaves the possibility to slow the processor with almost no impact on the overall execution time of the application. Rountree et al. developed Adagio which tracks task execution slacks and computes the appropriate frequency at which it should run. Although the first instance of a task is always run at the highest frequency, further instances of the same task are executed at the frequency that was computed after it is first seen. Authors of [16] propose a tool called jitter, which detects slack periods in performance to performance inter-node imbalance and uses DVFS to adjust the CPU frequency accordingly.

Our approach differs from that implemented in Adagio in that our fine-grained data collection offers the possibility to differentiate not only compute-intensive and communication-intensive execution portions (these portions are referred to as phases/regions) but also memory-intensive phases. Memory-intensive phases can be run on a slower core without significant performance penalty [17].

Isci et al. [18] and Choi et al. [19] use on-line techniques to detect applications execution phases, characterize them and set the appropriate CPU frequency accordingly. They rely on hardware monitoring counters to compute runtime statistics such as cache hit/miss ratio, memory access counts, retired instructions counts, etc. which are then used for phase detection and characterization. Policies developed in [18,19] tend to be designed for single task environments. We overcome that limitation by considering each node of the cluster as a black box, which means that we do not focus on any applications, but instead on the platform. The flexibility provided by this assumption enables us to track not applications/workloads execution phases, but node’s execution phases. Our work also differs from previous works in that we use partial phase recognition instead of phase prediction, which is not application specific and does not require multiple executions of the same application. On-line recognition of communication phases in MPI application was investigated by Lim et al. [20]. Once a communication phase is recognized, authors apply CPU DVFS to save energy. They intercept and record the sequence of MPI calls during program execution and consider a segment of program code to be reducible if there are high concentrated MPI calls or if an MPI call is long enough. The CPU is then set to run at the appropriate frequency when the reducible region is recognized again.

Our work differs from those above in two major ways. First, our phase detection approach does not rely on a specific HPC subsystem or MPI communication calls. Second, unlike previous research efforts, our model goes beyond the processor, since it also takes advantage of power saving capabilities available on all HPC subsystems. For example slowing down networks interfaces using Adaptive Link Rate, putting disks in low power-mode, and switching off memory banks are power saving schemes which are not directly linked to the processor.

3. Understanding HPC systems’ behaviour: on-line and off-line models

In this section, we present two effective approaches for tracking HPC systems’ behaviours. The first approach which we refer to as “DNA-like representation” of the HPC system attempts to describe an HPC system as a state graph in which each state represents its behaviour over a time interval; whereas the second approach performs on-line detection and characterization of execution phases of an HPC system at runtime. As mentioned earlier, we assume that an HPC system is a set of computing and storage nodes and will use the term “system” to designate a single node of the HPC system. Network equipments such as routers and switches are not taken into account because of their nearly constant power consumption. In the rest of this paper, unless otherwise expressly stated, we use

<table>
<thead>
<tr>
<th>CPU performance events</th>
<th>Memory performance events</th>
</tr>
</thead>
<tbody>
<tr>
<td>Instructions executed</td>
<td>Instruction fetch misses</td>
</tr>
<tr>
<td>Data dependencies</td>
<td>Data dependencies</td>
</tr>
<tr>
<td>Instruction cache misses</td>
<td></td>
</tr>
<tr>
<td>TLB misses</td>
<td></td>
</tr>
</tbody>
</table>
the term “sensors” to designate performance counters along with network bytes sent/received and disk read/write counts. Sensors related to hardware performance counters provide insight into the processor and memory activities. Likewise, disk and network related sensors provide insight into disk and network activities respectively.

3.1. DNA-like system modelling

The DNA-like system modelling models each node of an HPC system as a graph whose states represent the execution behaviour of the system over fixed length time intervals. We made the assumption that initial and final states of the graph are states or configuration in which the system is idle. A transition from a state $S_1$ to a state $S_2$ of the graph is weighted by the conditional probability that the system goes from $S_1$ to $S_2$. We refer to successive states through which a system goes throughout its life cycle as its “DNA-like” structure. Since not all states of the graph have the same behaviour, the DNA-like structure of a system can be thought of as a succession of behaviours through which the system went over time. From this, we define the terms “letter” and “system description alphabet” as follows: a letter or phase is defined as a behaviour in the DNA-like structure of a system (a state of the graph modelling that system); whereas, the system description alphabet is the set of possible behaviours.

With the description just given, the runtime behaviour of a system can roughly be represented by a sequence of the form $L_1, L_2, \ldots, L_i$; where the $L_i$ are elements of the system description alphabet. It is possible that some states do not appear in the system description alphabet, those are represented by the $X_i$ notation.

A letter is modelled as a column vector of sensors. Details regarding their construction are provided later in this paper. The choice of sensors for representing letters is constrained by the fact that a letter must provide information about the computational behaviour of the workload (is it memory intensive, processor intensive or communication intensive?) and its energy/power consumption.

3.1.1. Letter modelling and representation

The literature corroborates our observation that sensors relevant to power consumption estimation model depend upon workloads/applications being executed (see Section 2 for details). Considering a finite set of sensors is used to estimate the power consumption of a given category of workloads; this suggests that changes in the set of sensors relevant to power consumption estimation over a time interval $T$ reflect changes in the system’s behaviour over that very time interval.

Relying on the above assumption, we propose Algorithm 1 which partitions the runtime of a system into different behaviours according to changes in the set of sensors relevant to power estimation. For the selection of relevant sensors we use this very simple power model $\text{Power} \sim \sum_{i=1}^{p} C_i X_i$, in which $C_i$ and $X_i$ are model coefficients and sensors respectively. We conduct a multivariable linear regression to obtain coefficients $C_i$ and retain sensors $X_i$ exhibiting a 5% level of statistical significance to power consumption estimation given the previous power model. For the sake of simplicity, we limit these sensors to four (i.e., 4 sensors for representing a letter).

Once letters are defined, we use the following formalism for their encoding: Let us assign each sensor to a four-bit aggregation or half byte. Our quadruplet (each letter is a vector of four sensors) is therefore of the form $(X_1, X_2, X_3, X_4)$; where each $X_i$ values is a half byte. Now, deleting commas in between the $X_i$ gives a sixteen-bit aggregation which converted into decimal is an unsigned integer. The unsigned integer obtained from the above transformation is then the final representation of a letter.

Data: $A$: a set of units, where a unit is composed of values of sensors collected at a given time; units are sampled on a per second basis. Note that they are arranged in their order of occurrence in time.

Result: $P = \{t_i\}$ where $t_i$ are points in time at which changes in the behaviour of the system were detected.

Initialization: consider a set $S$ made up of $k$ successive units along the execution time-line. Let us denote by $S_{upper}$ the time at which the last unit of $S$ was sampled, $k$ is chosen such that $k > p + 1$, where $p$ is the number of sensors.

Add the time at which the first unit of $S$ was sampled to $P$; Compute the set $R_0$ of sensors relevant to power consumption estimation using the dataset composed of units in $S$.

while units available do

Add $k$ units to $S$ and update $S_{upper}$ to $S_{upper} + k$

Compute the set $R_i$ of relevant sensors from $S$ if $R_{i-1} \neq R_i$ then

Find the point in time $j \in \{S_{upper} - k, S_{upper}\}$ such that the set of relevant sensors $R$ computed from the set whose last unit was sampled at time $j$ is the same as $R_{i-1}$

Go to Initialization

Algorithm 1: Algorithm to detect application phases.

3.1.2. Example

This example investigates how close to reality is our approach for partitioning a system’s runtime into different computational behaviours (or simply phases). To accomplish this, we successively run two applications IS and EP from the NAS Parallel Benchmark (NPB) suite [21]. These applications are opposite from their computational standpoint in the sense that IS is communication intensive whereas EP is mainly computing. Data collected during their execution are used as input to the algorithm. Fig. 1(b) and (a) (where the curve gives the variation of the power consumption of the system over time and rectangles delimit detected phase) indicates that using a simple power model it is possible to detect phases the system went through. The couple of integers appearing in each rectangle on the figure gives for each phase the corresponding letter coded as an unsigned integer and the amount of time the system spent in that particular phase. For this example, the DNA-like structure of each node is straightforward, for the first node, it is: $\text{Idle}(17694\ 19)\ |\ (17714\ 43)\ \text{Idle}$.

3.2. Execution vectors based system behaviour tracking

It can be seen from Fig. 1(c) and (d) – where the $y$-axis represents the access rate of performance counters and the $x$-axis the execution time-line – that the access pattern of hardware performance counters or sensors in general strongly reflects changes in the behaviour of the application/system. We speak of sensors’ access rate because they are normalized with respect to the number of cycles. From the observation just made, the concept of execution vector which is similar to power vector (PV) [22] seems adequate for phase detection. An execution vector (EV) is defined as a column vector whose entries are system’s metrics, including hardware performance counters, network byte sent/received and disk read/write counts. To remain consistent with previous sections, we shall refer to these system metrics as sensors.

The sampling rate corresponding to the time interval after which sensors’ values are read depends on the granularity. While a larger sampling rate may hide information regarding the system’s behaviour, a smaller sampling rate may incur a non negligible overhead. In this paper, we use a sampling rate of one second and
We define the resemblance or similarity metric between two EVs as the Manhattan distance between them. The Manhattan distance suits the case since it weighs more heavily differences. A phase change is detected when the Manhattan distance between consecutive EVs exceeds a preset threshold. The threshold is fixed in the sense that it is always the same percentage – we refer to that percentage as the detection threshold – of the maximum distance between consecutive EVs (e.g., if the detection threshold is X%, then the threshold is X% of the maximum distance between consecutive EVs). However, the maximum distance between consecutive EVs is zeroed once a phase change is detected. So, technically, the threshold varies throughout the system's life cycle. The maximum existing distance between consecutive EVs is continuously updated until a phase change is detected where it is zeroed. The idea behind zeroing the maximum existing distance when a phase change occurs is to allow detecting phase changes when changing from a phase where distances between consecutive EVs are big to a phase where they are not and vice versa. For example, given a 10% similarity threshold, two consecutive EVs belong to the same group if the Manhattan distance between them is less than 10% of the maximum existing distance between all consecutive execution vectors. Fig. 2(b) shows the distance between consecutive execution vectors and the variation of the threshold along the execution time-line when the system was running the Advance Research Weather Research and Forecasting (WRF-ARW) model [23].

A graphical view of our phase detection mechanism is provided in Fig. 2(a), where dashed vertical lines indicate the start and end times of WRF-ARW. The left extremity of horizontal solid lines indicates the point (in the execution time-line) at which phase changes are detected and their length indicate the duration or length of corresponding phases. The x-axis represents the execution time-line and the y-axis represents IDs associated to detected phases. Note that IDs of phases are non zero integers ordered by their appearance order. It can be seen from Fig. 2(c) that detected phases are corroborated by the access pattern of sensors. The key motivation behind phase tracking is the use of characteristics of known phases for optimizing similar phases. Depending on the needs, optimizations aim to reduce the execution time of workloads, and/or their energy consumption. An effective phase characterization is therefore needed. To accomplish this, once a phase is detected, we apply principal component analyses (PCA) on the dataset composed of EVs pertaining to that phase. We next select five sensors among those contributing the least to the first principal axis of PCA. These 5 sensors serve as the characteristic of the phase. Since data collected during a phase can be too large for efficient representation and comparison in hardware, we summarize a phase by the EV at the centroid of the group made up of EVs sampled during that phase. We refer to the EV summarizing a phase as the reference vector of that phase.

3.2.1. Partial phase recognition

Phase identification is an interesting property of phase detection mechanisms, since it allows reuse of system reconfiguration information for reoccurring phases. However, a phase cannot be identified with a known phase before its completion. The literature suggests using phase prediction, which predicts the next phase of an application before it gets started; however, it might not work very well when you do not have any a priori information about applications sharing your platform. Therefore, instead of identifying...
a whole phase with a known, we propose to only identify a part of that phase with the known phase and extrapolate the result to the remaining part. We refer to that process as partial phase recognition; further details are provided in the next paragraph.

Partial phase recognition actually consists of identifying an ongoing phase (the phase has started and is not yet finished) $P_i$ with a known phase $P_j$ only considering the already executed part of $P_i$. The already executed part of $P_i$ expressed as a percentage of the length (duration) of $P_j$ is referred to as the recognition threshold $RT$. Thus, with a $RT\%$ recognition threshold, and assuming that the reference vector of $P_j$ is $EV_{P_j}$ and that its length is $m_j$, an ongoing phase $P_i$ is identified with $P_j$ if the Manhattan distance between $EV_{P_j}$ and each EV pertaining to the already executed part of $P_i$ (corresponding in length to $RT\%$ of $m_j$) are within the similarity threshold $ST$. The pseudo algorithm below summarizes partial phase recognition.

- let $P_j$ be a completed phase, $EV_{P_j}$ its reference vector, and $m_j$ its duration
- $P_i$ is partially recognized as $P_j$ if
  - $\forall \upsilon \in$ EV in the already executed part of $P_i$ (that is, EVs sampled between the time stamp corresponding to the start time of $P_i$ and the time stamp start time of $P_j + RT\%$ of $m_j$), the distance between $\upsilon$ and $EV_{P_j}$ is within threshold.

4. Model use cases: system adaptation and energy consumption prediction

Understanding the different behaviours of a high performance computing (HPC) system goes through throughout its life cycle can lead to a multitude of optimization opportunities. In this section, we shall investigate how one can leverage those behaviours for reducing the energy consumption of its infrastructure. Platform as a service becoming more and more attractive users often face the dilemma of choosing between multiple platforms for their applications. This section also investigates how users can determine the least energy consuming platform for their applications. In the following, we first show how our DNA-like modelling approach can be used for predicting the energy consumption of an application on a given platform. We next use our phase tracking methodology for reducing the energy consumption of applications without a priori knowledge.

4.1. Use case 1: cross platform energy prediction

The main motivation behind this use case is that users often have more than one candidate platform for running their jobs, in which case choosing the least energy consuming platform may be beneficial both for them and the platform provider. For simplicity, we assume that the application whose energy consumption is being predicted is the only application running on the platform under consideration. In other words, the DNA-like structure of the system at hand is that of the application.

The prediction model implicitly uses two sets of data, one from a reference platform provided by the DNA-like structure of the application, and one from a target platform which is the platform on which we want to estimate the overall energy consumption of the application. The aforementioned reference platform is the platform on which the application was first run and where its DNA-like structure was built. In most cases, the reference platform will
be the platform on which the application was tested (nearly all the information – including the application’s energy consumption – in regard to the reference platform is known).

Knowing the DNA-like structure of an application, one can easily identify applications with the same computational requirements. We accomplish this by comparing the DNA-like structure of the ongoing application to known DNA-like structures. A match is found when the already executed part of the ongoing application matches with a given percentage of the known DNA-like structure.

Let us denote $E_{tar}$ the energy consumed by the already executed part of the application of which we want to estimate the energy consumption and $E_{ref}$ the energy consumed by the corresponding part of the application whose DNA-like matches with the already executed part of the application at hand (the application that we want to estimate the energy consumption). For example, considering an application that lasts 60 min on its reference platform, let us assume that $E_{tar}$ represents the energy consumed by the same application on the target platform after 10 min run; therefore, $E_{ref}$ represents the proportion of energy consumed by the application on the reference platform during the first 10 min of its execution. Denoted as $E_{rel}$, the relative energy consumption between the two platforms is given by the following equation:

$$E_{rel} = \frac{E_{tar}}{E_{ref}}. \quad (1)$$

With the above relative energy, the estimated energy consumption of the application on the target platform is given by Eq. (2):

$$E_{est} = \int_0^{X} P(t)_{tar} \, dt + E_{rel} \star \int_{X}^{end} P(t)_{ref} \, dt. \quad (2)$$

Where $E_{est}$ is the estimated energy consumption on the target platform; $E_{rel}$ the relative energy consumption between the target platform and the reference platform. In the above equation, $\int_0^{X} P(t)_{tar} \, dt$ represents the energy consumed by the application before a match is found with a known DNA-like structure. Either measured or estimated, $P(t)_{tar}$ is the instantaneous power usage of the application on the target platform. Likewise, $P(t)_{ref}$ is the instantaneous power usage of the application on the reference platform and can be obtained from its DNA-like structure. In cases where the power consumption may radically change after the X% threshold, if the change in power consumption does not imply any change in the set of sensors used to estimate the power consumption, then we still assume it is the same application otherwise we attempt to find another match.

We define the estimation accuracy as the ratio between estimated and measured energy on the target platform, i.e.,

$$\text{Accuracy} = \frac{E_{est,tar}}{E_{tar}} \quad (3)$$

where $E_{est,tar}$ is the estimated energy consumption on the target platform.

Although comparing two DNA-like structures boils down to comparing two strings, the overhead associated with matching the DNA-like structure of a running application with previously seen known applications is proportional to the size of already known applications times the size of the DNA-like structure of the running application. We simplify this with the assumption that our profile database only contains one application which is that of the application of which we want to predict the energy consumption. We also assume that the application follow a very simple pattern which starts with an initialization phase and finishes which a finalization phase. Between the initialization and the finalization phases, there are some iterative computations and optional communications. Finally, assuming that the instantaneous power usage of the application is approximately the same during each of its iterations; meaning that the energy consumed by the application in each iteration throughout its life cycle is nearly the same, Eq. (2) can be simplified to Eq. (4); where $E_{int}$ represents the energy consumed by the application on the target platform during its initialization phase; $E_{ref-end}$ is the energy consumed by the application on the reference platform from the end of the initialization phase to the end of its whole execution; and $E_{ref-exe}$ is the measured energy consumption (resulting from complete execution) of the application on the reference platform.

$$E_{est} = E_{int} + E_{rel} \star (E_{ref-exe} - E_{ref-end}). \quad (4)$$

4.2. Use case 2: optimizing energy performance of HPC systems

The methodology described in Section 3.2 permits online detection and characterization of different runtime behaviours of the system. We use the coupling with partial phase recognition to guide on-the-fly system adaptation considering three HPC subsystems: processor, disk and network interconnect. The power consumption of these HPC subsystems along with that of the memory is about 55% [24] of the total power consumption of a typical HPC system. For the processor, we define three computational levels according to the characteristics of the workload:

- High or CPU-bound: the CPU-bound computational level corresponds to the maximum available CPU frequency, and is used for CPU-bound workloads.
- Medium or memory-bound: corresponds to an average in between the maximum and the minimum available frequencies; it is mainly used for memory-bound workloads.
- Low: the system is in the low computational level when the CPU frequency is set to the minimum available.

For the disk, we define two states: active and sleep; where active includes both the disk’s active and standby modes. Finally, for the network interconnect we define two data transfer speeds:

- The communication-intensive speed: corresponds to the highest available transfer rate of the network card.
- Low-communication speed: where the speed of the network interconnect is set to the lowest speed.

As mentioned earlier in this paper, principal component analysis (PCA) is applied to vectors belonging to any newly created phase for selecting five sensors which are used as phase characteristics. These characteristics are translated into system adaptation as detailed in Table 2. Let us comment the first row of that table. Workloads/applications with frequent cache references and misses are likely to be memory bound. In our case, having these sensors (cache_reference and cache_misses) selected from PCA indicates that the workload is not memory bound. If in addition that workload does not issue a high I/O rate (presence of I/O related sensors in the first column), then we assume that it is CPU-bound; consequently, the frequency of the processor can be scaled to its maximum, the disk sent to sleep and the speed of the interconnect scaled down. For the second row of Table 2, the characteristics do not include any I/O related sensor, this implies that the system was running and I/O intensive workload; thus, the processor’s speed can be set to its minimum. Note in passing that changing the disk’s state from sleep to active does not appear in Table 2, this is because the disk automatically enters the active state when it is accessed.

5. Experimentation and validation

5.1. Evaluation for energy consumption prediction

We evaluate our energy estimation model considering two workloads: the first workload (workload_1) iteratively computes the inverse of a $10 \times 10$ matrix and copies a large file from a remote repository; the second workload is GeneHunter a real life program for linkage analysis [25].
We further consider three scenarios: (i) the first scenario estimates the energy consumption of workload\textsubscript{1} on a node running at 2.13 GHz using the same node running at 1.6 GHz as the reference platform; (ii) scenario 2 still estimates the energy usage of workload\textsubscript{1}, but uses a Dell Power Edge server and a Sun Fire V20z as reference and target platforms respectively; (iii) in the third and last scenario we attempt to estimate the energy consumption of GeneHunter. For this specific case, our reference platform is an Intel Xeon E5506 Quad-core with 8 cores and 12 GB of RAM while the target platform is an Intel Xeon X3440 with 4 cores and 16 GB of RAM. For all three scenarios, an empiric partial execution threshold of 20% is used. This means that, a match with an existing DNA-like structure $DS$ is found if the already executed part of our synthetic application matches with 20% of $DS$, i.e., assuming $DS$ lasted 60 s, a match will be found if the already executed part of the synthetic application matches with the DNA-like structure describing the first 12 s of $DS$.

We compute for each scenario the expected energy consumption based on Eq. (4); results are summarized in Fig. 3. We can see from those results that the accuracy is very good. Notice that the accuracy is higher because it is computed considering the average energy consumption. We believe that overestimating the accuracy is higher because it is computed considering the average energy consumption. We also consider three benchmarks (LU, BT and SP) from NPB suite and two real-life applications: Molecular Dynamics Simulation (MDS)\cite{27} and the Advance Research Weather Research and Forecasting (WRF-AWR) model\cite{28,23}. WRF-AWR is a fully compressible conservative-form non-hydrostatic atmospheric model. It uses an explicit time-splitting integration technique to efficiently integrate the Euler equation. The classical Molecular Dynamics solves numerical Newton’s equations of motion for the interaction of the many particles system. We monitored each node’s power usage with one sample per second using a power distribution unit.

5.2.2. Results analysis and discussion

To evaluate our system adaptation policy, we consider 3 basic configurations of the monitored cluster: the first configuration which we refer to as on-demand is the configuration in which the default Linux’s on-demand governor is enabled on all the nodes of the cluster; the second configuration called performance is the configuration in which the Linux’s performance governor is enabled on each node of the cluster; and finally the third configuration which we refer to as managed is the configuration in which our system adaptation policy is applied. We also consider two levels of system adaptation:

- System adaptation level one: It corresponds to the situation in which only processor related optimization is made.
- System adaptation level two: It embraces level one, and additionally considers optimizing the interconnect and the disk.

Results we present here are obtained using an empirical similarity threshold $ST$ of 5%. The same goes for the recognition threshold $RT$ which is set to 10%.

(a) System adaptation level one: Processor’s only optimization: when an ongoing phase is identified with an existing phase, the characteristics of the existing phase are used to adapt the processor’s frequency accordingly. Diagrams of Fig. 4 show the average energy consumption (Fig. 4(a)) and execution time (Fig. 4(b)) of MDS and WRF-AWR under the three system’s configurations. These diagrams indicate that our management policy can save up to 19% of the total energy consumption with less than 4% performance loss.

Fig. 4(b) shows that on-demand and performance governors nearly achieve the same performance. This is because Linux’s on-demand governor do not lower the CPU frequency unless the system load decreases below a given threshold. Traces of CPU load under WRF-AWR for one node of our cluster are shown in Fig. 5, where the y-axis represents the percentage of load. The plot indicates that the CPU load remains above 85%, in which case the
on-demand and performance governors almost have the same behaviour. With processor’s only optimization, our management policy differs from that of the Linux’s on-demand governor in that we do not use the system’s load as system adaptation metric, which enable us to consume less energy.

(b) System adaptation level two: Processor, disk and network optimization: Fig. 6 presents the energy performance considering the processor, along with disk and network. These graphs indicate that considering the processor along with disk and network interconnect improves energy performance up to 24% with the same performance degradation as with processor’s only optimization. In other words, the disk and the network also contribute in improving energy performance.

5.2.3. Threshold selection and energy performance
In this section, we investigate the influence of phase detection parameters $ST$ and $RT$ on application performance and energy consumption. Firstly, we set the detection threshold and vary the partial recognition threshold. Secondly, we set the partial recognition threshold and vary the detection threshold. Fig. 7 — where the y-axis represents either the average energy consumption (Fig. 7(a)) or the average execution time (Fig. 7(b)) and the x-axis the detection threshold — shows the impact of the detection threshold on the execution time and energy consumption of WRF-AWR. It can be seen from that figure that for WRF-AWR, a detection threshold either of 5% or 20% could be a good choice. However, 5% could be preferable since there is a difference in energy consumption of up to 5000 J, for a less than 10 s difference in execution time. Fig. 7 also reveals that the detection threshold might have a significant impact on performance (both in terms of energy consumption and execution time). For these experiments, the partial recognition threshold was fixed to 10%.

The influence of the recognition threshold $RT$ on energy performance is summarized by Fig. 8, where the x-axis represents the recognition threshold and the y-axis either the average energy consumption (Fig. 8(a)) or the average execution time (Fig. 8(b)).
According to Fig. 8, a partial recognition threshold of 15% is convenient both in terms of energy and execution time.

What we can learn from Figs. 7 and 8 is that these values must be chosen depending on the target objective. A small recognition threshold may limit the impact of wrong decisions; however it may also have an influence on right decisions. For, making the right decision earlier allows saving more energy, in reverse, making a wrong decision earlier can result in significant energy waste and/or performance degradation.

5.3. Performance analyses of multiple applications

Earlier herein, we talked about the usefulness to look at a whole system; however, our evaluations has so far focused on only one application at a time. In this section, we evaluate our strategy with multiple applications running at the same time. We consider as test applications the WRF-AWR model and BT from the NAS Parallel Benchmark (NPB) suite. WRF-AWR spans 48 processes and run on 48 cores (in other words 12 nodes) whereas BT spans 48 processes on 12 nodes. The two applications do not share any resources except the network that interconnect them. We measure the power consumption of each of our 24-node cluster using a power distribution unit (PDU). Fig. 9, where the x-axis represents the average percentage of energy improvement or performance loss, indicates that our methodology still function when there are multiple applications sharing the platform; however, we also notice a performance degradation of up to 15% for BT. We are currently investigating the reason of BT performance degradation. Nevertheless, we believe this can be attributed to the fact that some phases might have been considered as memory intensive or communication intensive while they were actually compute intensive. In addition, applications (BT for example) which do not implement load imbalance are likely to experience higher performance degradation than those which implement load imbalance (WRF-AWR for example) [29,30].

6. Conclusion and future works

In this paper, we present two generic approaches for tracking high performance computing systems’ behaviour regardless of the applications being executed. We show through two use cases how they can be used for improving energy performance of a HPC system at runtime on the one hand and on the other hand how they can be used for estimating the energy consumption of an application given a target platform. Experimental results reveal the effectiveness of these two methodologies under real-life workloads and benchmarks. Comparison of our system adaptation policy with baseline unmanaged execution shows that we can save up to 19% of energy with less than 4% performance loss considering the processor only and up to 24% considering the disk, processor and network. As our methodology for improving energy performance does not depend on any application specific metric, we expect it to be
extended to a large number of power-aware HPC systems. Future works include combining the two system's behaviour tracking to enable live feedback on whether a system adaptation will be effective or not. This can help improve the energy performance while reducing performance loss. We also plan on adding memory optimization to processor, disk and network optimization and integrating workload consolidation and migration as core functionalities of the system management policy.

Acknowledgments

This work is supported by the INRIA large scale initiative Hemera focused on “developing large scale parallel and distributed experiments”. Experiments presented in this paper were carried out using the Grid5000 experimental testbed, being developed under the INRIA ALADDIN development action with support from CNRS, RENATER and several Universities as well as other funding bodies (see https://www.grid5000.fr).

References


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