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# ▶ To cite this version:

Andrea Fox, Francesco De Pellegrini, Eitan Altman. Learning Optimal Edge Processing with Offloading and Energy Harvesting. 2023. hal-04022507v1

# HAL Id: hal-04022507 https://hal.science/hal-04022507v1

Preprint submitted on 10 Mar 2023 (v1), last revised 23 Mar 2023 (v2)

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# Learning Optimal Edge Processing with Offloading and Energy Harvesting

Andrea Fox\*, Francesco De Pellegrini\* and E. Altman\*<sup>†</sup>

# ABSTRACT

Modern portable devices can execute increasingly sophisticated AI models on sensed data. The complexity of such processing tasks is data-dependent and has relevant energy cost. This work develops an Age of Information markovian model for a system where multiple battery-operated devices perform data processing and energy harvesting in parallel. Part of their computational burden is offloaded to an edge server which polls devices at given rate. The structural properties of the optimal policy for a single device-server system are derived. They permit to derive a new model-free reinforcement learning method specialized for monotone policies, namely Ordered Q-Learning, providing a fast procedure to learn the optimal policy. The method is oblivious to the devices' battery capacities, the cost and the value of data batch processing and to the dynamics of the energy harvesting process. Finally, the polling strategy of the server is optimized by combining such policy improvement techniques with stochastic approximation methods. Extensive numerical results provide insight into the system properties and demonstrate that the proposed learning algorithms outperform existing baselines.

# **KEYWORDS**

AoI, Energy-Harvesting, Offloading, Reinforcement Learning

#### **ACM Reference Format:**

Andrea Fox\*, Francesco De Pellegrini\* and E. Altman\*<sup>†</sup>. 2023. Learning Optimal Edge Processing with Offloading and Energy Harvesting. In Proceedings of ACM Conference (Conference'17). ACM, New York, NY, USA, 

#### 1 **INTRODUCTION**

New generations of mobile access networks promise low delay and high-speed throughput data connections paired with in-network processing capabilities [1]. They will support mobile computing services able to integrate AI-intensive processing tasks in their workloads such as, e.g., smart city services or virtual and augmented reality applications. In the next future, the vast majority of enterprise IoT projects are expected to have an AI component, up from less

Conference'17, July 2017, Washington, DC, USA

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ACM ISBN 978-x-xxxx-x/YY/MM...\$15.00

https://doi.org/10.1145/nnnnnn.nnnnnn

than 10% in 2018 [2]. To support these applications, edge networks must evolve to solve the key challenges of such scenario [3]. In fact, the energy consumption of AI applications is critical for battery operated devices. On the other hand, they may require periodic updates based on fresh data, settling specific requirements on the rate at which data are fetched and processed. The standard metric to this respect is the Age of Information (AoI), denoting the freshness of information when received at the destination [4]. In the literature, AoI performance has been studied for several queuing disciplines determining the average system time of data reads. The standard objective is to maximize some long term reward for the AoI of retrieved information [5, 6]. In the context of this paper, it is the long term reward for processing data, also called age of processing [7].

Energy harvesting from renewable sources such as solar cells or piezoelectric generators has becoming available on devices used for Internet of things (IoT) applications. In the literature, AoI optimal policies under energy harvesting have been studied using Markov decision theory both in continuous and in discrete time [8-12]. Most such studies focused on optimal policies to optimize AoI subject to energy causality constraints. The problem is to reserve the battery charge for moments when it is most needed, e.g., for alarm generation events [12].

The scenario studied in this paper considers the uncertainty of the battery consumption and the uncertainty of the amount of energy harvested over time. In fact, the energy cost of AI tasks depend inherently on the statistical distribution of input data and on the preference for lightweight or heavyweight models. Thus, when such a computing task is launched on a batch of data, the energy required to finish processing may exceed the available battery charge. In this event, a further delay for data processing is needed in order to harvest enough energy to complete the ongoing task.

Finally, in edge computing, task offloading to edge-servers mitigates the problem of energy consumption to run AI tasks on mobile devices [7, 13-19]. Thus, a device can delegate computing tasks to edge servers. However, in a realistic scenario, offloading is constrained by the availability of the edge-server, which may be serving multiple devices. Not always task offloading proves convenient, depending also on the delays it introduces.

Main contribution. This work develops a markovian modeling framework where multiple battery-operated devices process data in parallel and perform energy harvesting. Furthermore, an intermittent edge-server supports task offloading. Events of battery depletion may occur due to the unknown complexity of data processing and to the randomness of the harvesting process. Hence, an optimal stationary policy prescribes whether or not a device should process a new data batch based on AoI and battery status. By proving that such policy is monotone it is possible to design a lightweight reinforcement learning (RL) algorithm, namely Ordered Q-learning. It applies to the wide class of problems with monotone structure of the optimal

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Table 1: List of symbols used in the paper.

Symbol	Meaning
t	time step $t = 1, 2,$
x <sub>t</sub>	age of information (AoI) $x_t \in \{1, \ldots, M\}$
et	buffer energy level $e_t \in \{0, 1, \ldots, B\}$
$z_t$	server availability $z_t \in \{0, 1\}$
$s_t = (x_t, e_t, z_t)$	device state
γ	discount factor
S	state space
$\mathcal{A} = \{0, 1\}$	action set, action set of state $s$ , $\mathcal{A}(s) \subset \mathcal{A}$
$a_t = \{0, 1\}$	action taken at time t
В	battery capacity
М	saturation bound on AoI
$C_t$	batch processing cost at time $t$
$H_t$	harvested energy at time t
δ	offloading roundtrip time



Figure 1: Device performing local processing of data batches with energy harvesting and intermittent edge-server offloading.

value function with respect to the state-action partial order. In this setting, it outperforms standard baselines including policy gradient methods. Finally, the polling rate vector of the server is optimized by stochastic learning. To the best of the authors' knowledge, this is the first work to study the problem of AoI minimization in a system where multiple battery operated devices perform energy harvesting and rely on an edge-server for task offloading, incurring possibly in battery depletion depending on data and battery status.

The paper is organized as follows. Sec. 2 describes the state of the art. In Sec. 3 the basic server-device Markov decision model is introduced. The monotone structure of the optimal policy is analyzed in Sec. 4. Sec. 5 describes RL algorithms converging the optimal solution. Sec. 6 extends analysis and algorithmic solutions to the case of multiple devices. Numerical results are provided in Sec. 7 and a concluding section ends the paper.

# 2 RELATED WORKS

The term AoI was first used in [4] to denote freshness of data received at the destination. Most works in the literature minimize the average or the peak AoI [5]. AoI minimization has been studied in combination with energy harvesting to charge IoT devices. Energyaware scheduling policies of the type studied in this paper have been introduced to vary the sensing rate based on the instantaneous battery charge [10]. Conversely, the effect of error prone channels of the type studied in [8] and [9] is left for future works. In the proposed model, as in [10, 11], measurements are possible only when the battery level is sufficiently high. As in [12], the model accounts for data-dependent energy consumption. However, in all those works, events of battery depletion are not considered.

Several works combine AoI minimization and task offloading to external devices. In [7], [17] and [18] is studied the performance of a device-server system where a local processor device is supported by a remote server. As the present work, [13] and [14] study the tradeoff between energy consumption and execution delays to offload applications to an edge server. In [15] and [16] deep reinforcement learning (DRL) is used to compute optimal offloading policies. As showed later, knowing the properties of the optimal value function permits much less expensive solutions. With very few exceptions, e.g., [19] and [6], AoI in multi-device systems of the type studied in this paper are not addressed in the literature. In [19] the average AoI is minimized by choosing whether to use the local processor or an edge server according to the network status. [6] provides lower bounds on the average AoI performance achievable for several queuing disciplines on a single-hop network as the one studied in this work. Note that the metric addressed in this work is not the average AoI [20].

The first Markovian model for the control of AoI appeared in [21]. MDP models were later used in [12, 16, 19, 22]. Works [7, 8] use constrained MDPs, while [11, 23] use partially observable MDPs. Including constraints and partial observations are interesting extensions but out of the scope of the present work. Threshold policies for the control of the AoI have been obtained before [7, 9, 11]. In [9] the threshold is one-dimensional while the policy structure described in [7] is only characterized numerically. In [11] the threshold regards two incorrelated control variables: here, transitions for AoI and energy state variables both depend on the chosen action.

#### **3** SYSTEM MODEL

The device-server scheme is represented in Fig. 1: the device can read data batches and process them, while the edge server can poll the device to offload the computation. Thus, data batches are either read and processed locally on the device or read and offloaded. Time steps are discrete with index  $t = 1, 2, \dots$  Processing data at time t on the device requires  $C_t$  energy units, which form a sequence of i.i.d. random variables  $\{C_t\}$  with probability distribution  $p_C(c) = \mathbb{P}\{C = c\}$ . When the device is polled by the edge server, the data processing task is offloaded and it has zero energy cost for the device. However, sending and retrieving processed data from the server incurs in a constant delay of  $\delta \ge 0$  time units. The device has a battery of capacity B > 0 energy units and it can harvest a number of energy units  $H_t$  per timestep t. The energy fetched per timestep, namely the harvesting rate, forms a sequence of i.i.d. random variables  $\{H_t\}$  whose corresponding probability distribution is  $p_H(h) = \mathbb{P}\{H = h\}.$ 

Let introduce the Semi Markov Decision Process (SMDP) which models the system. The state of the device at time t is denoted as  $s_t = (x_t, e_t, v_t)$ , where  $x_t$  represents the age of information



Figure 2: A sample path of the process  $s_t = (x_t, e_t, 0)$  for costant H = 1: in red the vacation periods where the battery is empty. The duration of timeslot starting at time t = 6 is 4 timesteps.

(AoI),  $e_t$  is the device battery level and  $z_t$  is the server state. Binary variable  $z_t$  indicates whether the server has polled the device for offloading,  $z_t = 1$ , or not,  $z_t = 0$ . The corresponding per slot transition probability is  $p_Z(z'|z)$ , e.g.,  $p_Z(1|0)$  is the probability for the server to poll the device at t + 1 given that it did not at t. Note that  $x_t$  is the age of information of the last data batch processed by a tagged device and the age of information is measured at the end of each time slot. It holds  $x_t = 1$  when the device has just processed fresh data. Afterwards, the AoI increases of one timestep at every time slot t until either the device fetches a new data batch and performs the computation or the server polls the device and offloading occurs. In both cases, at the end of the data processing the AoI is reset to 1.

Freshness function u(x) is the utility for processing of a data batch x time units after the last one was processed;  $u(\cdot)$  is bounded and non-increasing. It is assumed that there exists M > 0 such that u(M + k) = u(M) for all k > 0 so that the age of information takes values in  $\{1, \ldots, M\}$ . The state space is denoted  $S = \{1, \ldots, M\} \times \{0, \ldots, B\} \times \{0, 1\}$ .

Finally, it is possible that the battery available at timeslot *t* is not sufficient to terminate the computation immediately, namely,  $\phi_t := c_t - (h_t + e_t) > 0$ . In this case, a delay is incurred in order to harvest a sufficient amount of energy and complete the processing of the current batch. The corresponding timeslot has a random duration, due to the stochastic nature of energy harvesting. The action set  $\mathcal{A} = \{0, 1\}$ , where 0 means *wait* and 1 *process*. The action taken at time *t* is denoted  $a_t$ . If  $a_t = 1$ , the device fetches a new data batch which is processed at energy cost  $c_t > 0$ . For every state  $s_t$  where  $z_t = 0$  the action set available at each device is to process or not the information, i.e.,  $\mathcal{A}(s_t) = \{0, 1\}$ . On the other hand, when  $z_t = 1$  the action set is  $\mathcal{A}(s_t) = \{1\}$ , i.e., when the device is polled by the server, by default the data unit is processed. Furthermore, the energy to transmit data to the server is assumed to be negligible compared to local data processing. Finally, the dynamics of the age

of information for data batches is

$$x_{t+1} = \begin{cases} 1 & \text{if } a_t = 1\\ [x_t+1]_M & \text{if } a_t = 0 \end{cases}$$

where  $[y]_A := \max\{0, \min\{y, A\}\}$ . The AoI at the renewal instants is also called *peak age of information* [20] in the literature: it corresponds to the processing rate.

It now possible to characterize the transition probabilities for a given action  $a \in \mathcal{A}(s)$ . Hereafter *H* is considered to be deterministic with  $H \equiv 1$  and the time index is omitted for notation's sake. The general case for stochastic harvesting rate is derived in [24].

Let s = (x, e, z) be the current state of the device and let next state s' = (x', e', z') under action *a*.

The transition probability from s to s' is written, for a = 0, as

$$p(s'|s,0) = \begin{cases} s = (x, e, 1), \\ p_Z(z' \mid 0) & s' = ([x+1]_M, [e+1]_B, z'), \\ 0 & \text{otherwise} \end{cases}$$
(1)

Instead, when a = 1, the transition probability from s to s' is

$$(s' \mid s, 1) = \begin{cases} s = (x, e, 0), \\ p_Z(z' \mid 0)p_C(e + 1 - e') & s' = (1, e', z'), \\ 0 < e' < B \\ p_Z(z \mid 0)\sum_{c=e+1}^{\infty} p_C(c) & s = (x, e, 0), \\ p_Z(z \mid 0)\sum_{c=1}^{e+1-B} p_C(c) & s' = (1, 0, z') \\ p_Z(z' \mid 1) & s' = (1, B, z') \\ s = (x, e, 1), \\ p_Z(z' \mid 1) & s' = (1, [e + 1]_B, z'), \\ e + h < B \\ 0 & \text{otherwise} \end{cases}$$

The sojourn times  $\tau(s, s')$  of the SMDP are associated to the transition from an origin state s = (x, e, z) to a destination s' = (x', e', z'). They are unitary except possibly those corresponding to the transition to a renewal state, i.e., where x' = 1. For that transition, if z = 1 the sojourn time is  $\tau(s, s') = 1 + \delta$ , due to the round-trip delay for the edge-server offloading. If z = 0 it is  $\tau(s, s') = 1$  when  $\phi = C - e - H \le 0$ . Otherwise, when C - e - H > 0 it writes

$$\tau(s, s') = \min\left\{k \mid \sum_{r=1}^{k} H_r > C - e - H\right\}$$
(3)

that is the number of recharges required in order to complete the computation. In the case of deterministic energy harvesting with  $H \equiv 1$ , it holds  $\tau(s, s') = (1 - z) \max\{1, C - e - H\} + z(1 + \delta)$ . Note that, in the AoI literature, what is here defined as sojourn time corresponds to the notion of system time. The reward under the state action pair  $(s_t, a_t)$  at time t is proportional to the freshness of information of the data unit when it ends to be processed, as well as on the initial level of energy. Formally it writes

$$r_{t+1}(s_t, a_t) = \begin{cases} u(x_t) - d(c_t - e_t - h_t) \cdot a_t & \text{if } z_t = 0\\ u([x_t + \delta]_M) & \text{if } z_t = 1 \end{cases}$$
(4)

where function  $d : \mathbb{R} \to \mathbb{R}^+$  is a penalty that depends on the amount of energy units required to complete the local processing after a = 1

р

is selected. It is further assumed that d(x) = 0 when  $x \le 0$  and that d is a bounded non-decreasing function.

### 3.1 Discounted Model

The processing policy  $\pi$  for a tagged device is a probability distribution over the state-action space. The value function for policy  $\pi$  is  $v_{\pi}(s) := E_{\pi}[G_t|s_t = s]$  where  $G_t^{\gamma} := \sum_{t=0}^{\infty} \gamma^k r_{t+1+k}(s_t, a_t)$ ; the *Q*-function  $q_{\pi}(s, a) := E_{\pi}[G_t|s_t = s, a_t = a]$  is used later in Section 5.2. In the discounted model presented next, the objective is to maximize  $v_{\pi}(s)$  in the set of stationary policies, where  $\pi = \pi(s)$  is the probability that the device performs action a = 1 in state *s*. As showed in Sec. 4.2, the corresponding ergodic state occupancy probability permits to calculate the sampling rate based on the average peak AoI. The optimal value function  $v_*(s)$  solves the the Bellman optimality equation associated to the discounted reward problem

$$v_*(x, e, 0) = \max \left\{ u(x) - \sum_{c=1}^{\infty} p_C(c) \sum_{h=1}^{\infty} p_H(h) d(e+h-c) + \sum_{c=1}^{\infty} p_C(c) \sum_{h=1}^{\infty} p_H(h) \gamma \mathbb{E}_Z[v_*(1, [e+h-c]_B, z)], u(x) + \gamma \sum_{h=1}^{\infty} p_H(h) \mathbb{E}_Z[v_*([x+1]_M, [e+h]_B, z)] \right\}$$

in the case z = 0. When z = 1 the action set is a singleton, then

$$\upsilon_*(x,e,1) = u([x+\delta]_M) + \sum_{h=1}^{\infty} p_H(h) \mathbb{E}_Z[\upsilon_*(1,[e+h]_B,z)]$$
 (5)

Structural properties. The optimal policy and the corresponding value functions have several properties that are exploited in the next section. First, from (4) it is immediate that, if the harvesting process  $\{H_t\}$  is stochastically larger than  $\{H'_t\}$ , also  $v_H(s) \ge v_{H'}(s)$  for all  $s \in S$ ; similar observations hold w.r.t. to process  $\{C_t\}$ .

Before proving the next result, a few definitions are introduced. First, a policy is monotone if the action is either increasing or decreasing in the state, given an assigned partial order on the state space S and on the action space A. In particular, the following partial order is defined on the state space

$$(x, e + 1, z) \ge (x, e, z')$$
  
 $(x - 1, e, z) \ge (x, e, z')$  (6)

whereas the obvious order is imposed on the action space  $A = \{0, 1\}$ . Let  $s^+$  be *larger* than  $s^-$  if  $s^+ \ge s^-$ . The real function  $\psi : S \times \mathcal{A} \to \mathbb{R}$  is *supermodular* if  $\psi(x^+, y^+) + \psi(x^-, y^-) \ge \psi(x^+, y^-) + \psi(x^-, y^+)$ , also equivalent to say that  $\psi(x^+, y) - \psi(x^-, y)$  is decreasing in y. The *tail transition probability*  $w(s'|s, a) := \sum_{\zeta \ge s'} p(\zeta|s, a)$  defines the probability of making a transition to states larger than s taking action a.

LEMMA 1. i.  $v_*(x, e, z)$  is non increasing in x ii.  $v_*(x, e, z)$  is non decreasing in e.

PROOF. The proof is made by verifying that the system model adheres to the following three assumptions [25].

*i.* Non decreasing rewards. Let consider  $s^+ = (x, e, z)$ ,  $s^- = (x + 1, e, z)$  for a fixed value of *c*. For a = 1,  $r(s^+, 1) = u(x) - d(c - e - h) \ge u(x + 1) - d(c - e - h) = r(s^-, 1)$ , as *u* is nonincreasing. For a = 0 the

same property holds since  $r(s^+, 0) = u(x) \ge u(x+1) = r(s^-, 0)$ . Now assume  $s^+ = (x, e+1, z)$  and  $s^- = (x, e, z)$ . For a = 0 the reward is equal for both states. For a = 1,  $r(s^+, 1) = u(x) - d(c - e - 1 - h) \ge u(x) - d(c - e - h) = r(s^-, 1)$  since *d* is by assumption nondecreasing. *ii. Supermodular rewards.* This property can be easily verified for both cases considered before;

*iii. Increasing tail transition probability.* First, let consider the case  $s^+ = (x, e, z)$  and  $s^- = (x + 1, e, z)$ . When a = 1,

$$w(s'|s^+, 1) = \sum_{\zeta \ge s'} p(\zeta|s^+, 1) = \sum_{\zeta \ge s'} p(\zeta|s^-, 1) = w(s'|s^-, 1)$$

as the final level of age of information is always 1 after action a = 1, while the final energy level only depends on the initial energy *e*. If a = 0, then  $w(s'|s^+, 0) = \mathbb{1}\{(x + 1, e + h, z) \ge s'\} \ge \mathbb{1}\{(x + 2, e + h, z) \ge s'\} = w(s'|s^-, 0)$  where the inequality is induced by the partial order of states.

Now, consider  $s^+ = (x, e + 1, z)$  and  $s^- = (x, e, z)$ . When a = 1, the inequality is verified because the final energy obtained with a transition that starts in  $s^+$  is larger or equal than the one obtained when starting in  $s^-$  w.p.1. Conversely, when a = 0,  $w(s' | s^+, 0) = \mathbb{1}\{(x + 1, e + h + 1, z) \ge s'\} \ge \mathbb{1}\{(x + 1, e + h, z) \ge s'\} = w(s' | s^-, 0)$  and the inequality follows from (6).

From Lemma 1 it follows a property of optimal *Q*-function  $q_*(s, a)$ , that will later be used to develop suitable learning algorithms. From Lemma 1 it follows easily

COROLLARY 1. Fixed  $a \in \{0, 1\}$ , then for  $s = (x, e, z) \in S$ i.  $q_*((x, e, z), a) \le q_*((x, [e + 1]_B, z), a)$ ii.  $q_*(([x + 1]_M, e, z), a) \le q_*((x, e, z), a)$ 

In all numerical tests, function Q appears indeed superadditive in  $S \times \mathcal{A}$ , a property that would be sufficient to prove the monotonicity of the optimal policy. The classic assumption for the supermodularity of the tail probabilities (see Prop. 4.7.3 in [25]) which grants the existence of a monotone policy can be easily disproved.

THEOREM 1 (OPTIMAL POLICY STRUCTURE). A stationary deterministic policy  $\pi$  is optimal if and only if it is monotone. In particular, there exist thresholds T(e),  $0 \le T(e) \le M$  such that : i.  $\pi(e, x, 0) = 1$  if and only if  $x \ge T(e)$ ; ii.  $T(e) \ge T(e + 1) + 1$ ; iii. if  $\gamma = 0$  then  $T(e) \ge T(e + 1)$ 

PROOF. i. At first it is showed the existence of a threshold T(e) for every level of energy e. Consider the function  $\Delta q_*(x, e, z) = q_*((x, e, z), 1) - q_*((x, e, z), 0)$ : it holds

$$\Delta q_*(x, e, 0) = -\mathbb{E}_C[d(e+h-c)] + \gamma \Big\{ \sum_{c=1}^{\infty} p_C(c) \mathbb{E}_Z[v_*(1, [e+h-c]_B, z)] \\ -\mathbb{E}_Z[v_*([x+1]_M, [e+h]_B, z)] \Big\}$$
(7)

The increment with regard to x is nonnegative, due to Lemma 1:

$$\Delta q_*(x+1,e,0) - \Delta q_*(x,e,0) = -\gamma \left( \mathbb{E}_Z [\upsilon_*([x+2]_M, [e+h]_B, z] - \mathbb{E}_Z [\upsilon_*([x+1]_M, [e+h]_B, z)] \right)$$

ii. Observe that all states  $([T(e) + k]_M, [e + 1]_B, z)$  are transient for any integer k > 1. Hence, one could replace the policy for those

states with no change in the state occupancy probability. iii. Follows from the structure of the instantaneous reward.

From Thm. 1 it follows that an optimal policy is deterministic and monotone with a threshold structure.

### **4 OPTIMAL POLICY PERFORMANCE**

The optimal policy  $\pi^*$  can be obtained using standard dynamic programming methods, under the assumption of having full information on the transition probabilities; in the following  $p_{\pi^*}$  is the stationary distribution given by  $\pi^*$ . Efficient variants of value iteration or policy iteration algorithms for monotone policies do exist; alternatively, the optimal solution can be found solving a suitable linear program [25]. This section introduces an evaluation method for any policy  $\pi$ , using a formulation that will be used throughout the work.

#### 4.1 Optimal Peak AoI

Once determined the optimal policy, the average peak AoI can be computed by considering the renewal process  $\{N(t), t \ge 0\}$  defined by the instants of visit to a state with age of information x = 1 for the corresponding Markov reward process. Let  $Y_r$ , r = 1, 2, ... be the random variable describing the length of the *r*th renewal and the corresponding renewal process [26]. Denote as  $\Delta N = \mathbb{E}[N(t+1)] - \mathbb{E}[N(t)]$  the average number of renewal events, i.e., of processing events, per time unit. From Blackwell theorem [27],  $\frac{\Delta N}{t} \rightarrow \frac{1}{\mathbb{E}[Y]}$ . Let denote *L* the number of transitions during renewal period *Y*: it is the mean return time to a state with x = 1, i.e., the inverse of he stationary probability of being in a state *s* where  $\pi^*(s) = 1$ :

$$L = \frac{1}{\sum_{s:\pi^*(s)=1} p_{\pi^*}(s)}$$
(8)

By recalling the form of the sojourn times  $\tau(s, s')$  of the SMDP, the expected value of the renewal cycle *Y* is computed as

$$\mathbb{E}[Y] = \sum_{\{s':x'=1\}} \mathbb{E}_{\substack{s \sim p_{\pi^*} \\ \pi^*(s)=1}} \left[\tau(s,s')\right] p_{\pi^*}(s') + \frac{1}{L} - 1$$

where the last equality applies a partition of the possible renewal states. For notation's sake, let

$$\chi(e') := \mathbb{E}_{\substack{s \sim p_{\pi^*} \\ \pi^*(s)=1, z'}} \left[ \tau(s, (1, e', z')) \right]$$

the expected sojourn time when the transition ends in a state s' = (1, e, ', z'). Computing  $\chi(e')$  accounts for the possible sample paths to renew to an energy state e'. Two possible cases are possible. If z = 1, then e + h = e' with  $p_H(\Delta e)$  where  $\Delta e = e' - e$ . Else, if z = 0, the possibility that  $\phi > 0$  requires to account for the recharges needed to get to the final level of energy e', which depends on the distribution of both H and C.

$$\chi(e') = p_Z(1)(1+\delta) \sum_{e=0}^{B} p_E(e) p_H(\Delta e) + p_Z(0) \left( \sum_{e=0}^{B} p_E(e) \sum_{r=1}^{\infty} \Gamma_{1,r} p_C(h_1 - \Delta e) + \sum_{k=2}^{\infty} k \sum_{e=0}^{B} p_E(e) \sum_{r=1}^{\infty} \Gamma_{k-1,r} \sum_{c=e+\sum_{j=1}^{k-1} h_j + 1}^{\infty} p_C(c) p_H\left(\Delta e + c - \sum_{j=1}^{k-1} h_j\right) \right)$$

where  $p_E(e)$  indicates the probability that action a = 1 is done when the energy level is equal to e (can be easily obtained from the stationary distribution under a given policy) and  $\Gamma_{k,r}$  is the probability to harvest *r* energy units in *k* steps, namely  $\Gamma_{k,r} = \mathbb{P}\left(\sum_{j=1}^{k} h_j = r\right)$ . Note how it has been assumed that  $\forall z_1, z_2 \mathbb{P}(z_1 | z_2) = p_Z(z_1 | z_2) = p_Z(z_1)$ . Without loss of generality, let  $k \leq B$ :  $\Gamma_{k,r}$  is defined by the following iteration

$$\Gamma_{k,r} = \begin{cases} \sum_{j=1}^{B} \Gamma_{k-1,r-j} p_H(j) & k \le r \le B\\ 0 & \text{otherwise} \end{cases}$$

where  $\Gamma_{1,r} = p_H(r) \mathbb{1}\{1 \le r \le B\}$ . Finally, the expected value of *X* is

$$\mathbb{E}[X] = \sum_{z=0}^{1} \sum_{e'=0}^{B} \frac{p_{\pi^*}(1, e', z)}{\sum_{j=0}^{B} p_{\pi^*}(1, j, z)} \chi(e') + \frac{1}{L} - 1$$

In Section 7 numerical results will explore how the model parameters influence the Peak Age of Information.

# 4.2 Reward of a policy

The learning procedure for multi-device model presented in Sec.6 pivots on the dependence of the discounted reward  $R_{\gamma,\pi}$  of a tagged policy  $\pi$  on the server polling rate $p_Z$ . The performance measure for a policy used by a device is the discounted reward given a certain initial state distribution. In particular, sample paths initiate on renewal states with age of information x = 1, energy *e* chosen uniformly at random and expected server availability  $p_Z$ .

Fixed  $\pi$ , let  $P_{\pi}$  be the transition matrix of the corresponding Markov Chain and let vector  $r_{\pi}$  be the instantaneous reward attained at each state by following the action indicated by the policy. Let further define vector  $d_0$  describing the initial distribution of states. The discounted reward of a device can be computed as

$$R_{\gamma,\pi} = d_0 \left( I + \gamma P_{\pi} + \gamma^2 P_{\pi}^2 + \dots \right) r_{\pi}$$
(9)

Basic facts of MDP theory imply that the geometric series generated by  $\gamma P_{\pi}$  is always converging so that

$$R_{\gamma,\pi} = d_0 \left( I - \gamma P_{\pi} \right)^{-1} r_{\pi} \tag{10}$$

To analyze this quantity it is useful to explicitly express the parametric dependence  $d_0 = d_0(p_Z)$  and  $P_{\pi} = P_{\pi}(p_Z)$ . In particular,

$$d_0(p_Z) = \begin{cases} \frac{1}{B+1}p_Z & x = 1, z = 0\\ \frac{1}{B+1}(1-p_Z) & x = 1, z = 1\\ 0 & \text{otherwise} \end{cases}$$

The bijective function  $f : \{1, ..., M\} \times \{0, ..., B\} \rightarrow \{1, ..., M(B+1)\}$  maps each couple (x, e) to an integer value and two matrixes  $P^0, P^1 \in \mathcal{M}^{(B+1)M \times (B+1)M}$  such that

$$(P^0)_{ij} = \mathbb{P}\left\{f(x', e') = j \mid f(x, e) = i, z = 0, a = 0\right\}$$

and

$$(P^1)_{ij} = \mathbb{P}\left\{f(x', e') = j \mid f(x, e) = i, z = 1, a = 1\right\}$$

Both matrix will depend on the distribution of the random variables C and H, as well as on the fixed policy  $\pi$ . The explicit dependence of the transition matrix  $P_{\pi}$  on the server's polling control writes

$$P_{\pi}(p_Z) = \begin{pmatrix} (1 - p_Z)P^0 & p_Z P^0 \\ (1 - p_Z)P^1 & p_Z P^1 \end{pmatrix}$$

which provides an explicit form to

$$R_{\gamma,\pi}(p_Z) = d_0(p_Z) \Big( I - \gamma P_\pi(p_Z) \Big)^{-1} r_\pi$$

*Remark.* For  $\delta > 0$  the function  $R_{\gamma,\pi}(p_Z)$  appears concave on the interval [0, 1] for all tested stationary policies and parameters of the Markov decision process. The proof of the convergence of the algorithm introduced in Section 6 assumes this fact to be true. A formal proof can be provided for a specific choice of parameters (M = B = 2) and a particular policy [24]. Additionally,  $R_{\gamma,\pi}(p_Z)$ is indeed continuous and differentiable since each block matrix appearing in its explicit form has such property, as showed in [24].

Finally, evaluating the optimal reward as a function of  $p_Z$  requires to consider different optimal policies and to determine

$$R_{\gamma}(p_Z) = \max R_{\gamma,\pi}(p_Z) \tag{11}$$

As it is a point-wise maximum, the continuity of  $\gamma(\cdot)$  is true. However, even if the conjecture about the concavity of  $R_{\gamma,\pi}$  is verified,  $R_{\gamma}(p_Z)$  is not necessarily concave as it is the maximum of a set of concave functions.

# **5 LEARNING THE OPTIMAL POLICY**

In a realistic setting, transition probabilities (1) and (2) may not be available. Furthermore, rewards (4) depend on the model used for the data-batch computation and also on data properties which may be only learned at runtime. A new model-free algorithm, namely *Ordered Q-learning* (OQL) is introduced next. It is a version of Qlearning (QL) [28] adapted for MDPs with monotone value functions under a given partial order imposed on state space. As showed in Section 7, OQL outperforms other reinforcement learning algorithms such as policy gradient methods and yet retains the convergence guarantees of QL. Two variants of OQL based on the structure of the optimal policy are introduced next; they leverage two different partial orders on the states to find the optimal policy.

# 5.1 Ordered Q-learning

In the Ordered Q-learning the vector of the estimates of the *Q*-function are forced to comply to the monotonicity properties of the optimal *Q*-function. The update rule (12) for each state-action pair  $(s, a) \in S \times A$  hence writes

$$Q_{t+1}(s,a) = (1 - \alpha_t)Q_t(s,a) + \alpha_t \left( r_{t+1}(s,a) + \gamma \max_b Q_t(s',b) \right)$$
$$Q_{t+1}(s',a) = \Pi_s \left( \overline{Q}_{t+1}(s',a) \right) \quad \forall s' \in \mathcal{S}$$
(12)

where projection  $\Pi_s(\cdot)$  at state *s* is a projection adapted to specific partial order imposed on the state space. Let  $f : S \to \mathbb{R}$ : if s' > s, then  $\Pi_s(f(s')) = \max\{f(s'), f(s)\}$  and if  $s' \leq s$ , then  $\Pi_s(f(s')) = \min\{f(s'), f(s)\}$ .

Thus, the basic Q-learning iteration is followed by a state-dependent projection on the set of allowed estimates. Alg. 1 reports on the complete algorithm for the version of the algorithm with a constant stepsize  $\alpha_t = \alpha$ . The convergence of the Ordered Q-Learning (OQL) can be ensured under the assumption that  $\alpha_t = \alpha_t(s, a)$  appearing in (12) is a standard stepsize i.e.,  $\sum_{t=1}^{+\infty} \alpha_t(s, a) = +\infty$  w.p.1. and  $\sum_{t=1}^{+\infty} \alpha_t(s, a)^2 < +\infty$  w.p.1.; it holds the following result

THEOREM 2. Consider the Ordered Q-learning algorithm described by the update rule in (12). Let  $\gamma < 1$ . Let  $q_*$  be monotone, i.e., if  $s_1 \leq s_2$  according to some order on the states, then  $q_*(s_1, a) \leq q_*(s_2, a)$ . Then  $q_{*,t}(s, a)$  converges to  $q_*(s, a)$  with probability 1 for every state  $s \in S$  and for every action  $a \in \mathcal{A}(s)$ .

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Rec	<b>quire:</b> step size $\alpha \in (0, 1]$ , discount factor $\gamma \in [0, 1)$		
1:	$Q(s, a) \in \mathbb{R}, \forall s \in \mathcal{S}$		
2:	$Q(s', a) \leftarrow 0$ for s' terminal state		
3:	3: <b>for</b> each episode		
4:	initialize $s \in S$		
5:	for each step of the episode		
6:	choose $a \in \mathcal{A}(s)$ given by $Q$ (e.g., $\epsilon$ -greedy)		
7:	take action $a$ and observe reward $R$ and $s'$		
8:	$Q(s, a) \leftarrow (1 - \alpha) Q(s, a) + \alpha [R + \gamma \max_a Q(s', a)]$		
9:	$Q(\tilde{s}, a) \leftarrow \Pi_s(Q(\tilde{s}, a)),  \forall \tilde{s} \in \mathcal{S}$		
10:	$s \leftarrow s'$		
11:	end for		
12:	end for		

PROOF. A sketch of the proof of Theorem 2 is based on the Policy Improvement theorem's [29]. Let consider a generic step of the Ordered Q-Learning algorithm. Let  $\pi$  be the current policy and  $\pi'$  be the policy obtained at the iteration of Ordered Q-learning. Let assume that  $\pi'$  improves the current policy, i.e.,  $q_{\pi}(s, \pi'(s)) \ge v_{\pi}(s)$  for all  $s \in S$ . As showed in [29], this implies that  $v'_{\pi}(s) \ge v_{\pi}(s)$  for all  $s \in S$ .

Now let consider any state  $s^+ \geq s$ . After applying the projection, it holds  $q_{\pi}(s^+, \pi'(s)) \geq q_{\pi}(s, \pi'(s))$ . Finally, since both  $\pi'(s)$  and  $\pi'(s^+) \in \mathcal{A}(s^+)$ , it is possible to write

$$v_{\pi'}(s^+) = q_{\pi'}(s^+, \pi'(s^+)) = \max\{q_{\pi'}(s^+, \pi'(s)), q_{\pi'}(s^+, \pi(s^+))\}$$
  
$$\geq q_{\pi'}(s^+, \pi'(s)) \geq q_{\pi'}(s, \pi'(s)) = v_{\pi'}(s)$$

which concludes the proof.

The above proof sketch is based on the simplifying assumption that the true values of the *Q*-function are available at each iteration; this assumption permits to use dynamic programming arguments. The complete proof of Theorem 2 is rooted in the original argument of convergence for Q-learning, developed by Tsitsiklis in [30], which is based on stochastic approximations. It is presented in the extended version of this paper [24]. For Ordered Q-learning, the technical difficulty is to account for the use of state-dependent projection  $\Pi_s$ .

It is further possible to define the *n*-step version of OQL, where the first equation in (12) is replaced by the one of *n*-step QL [29].

### 5.2 Stairway Q-Learning

Stairway Q-learning (SQL) is defined imposing on the state space the natural partial order considered in Cor. 1 so that the assumptions of Thm. 2 automatically hold for the system at hand. For every pair of states  $s_1 = (x_1, e_1, z)$  and  $s_2 = (x_2, e_2, z)$  and for every vector  $v \in \mathbb{R}^{|S|}$ , the explicit form of the projection operator on the set of vectors that suit the partial order considered is

$$\Pi_{s_1}(v(s_2)) = \begin{cases} \max(v(s_1), v(s_2)) & \text{if } x_1 < x_2 \text{ and } e_1 \ge e_2 \text{ or} \\ x_1 \le x_2 \text{ and } e_1 > e_2 \\ v(s_1) & \text{if } x_1 = x_2 \text{ and } e_1 > e_2 \\ v(s_1) & \text{if } x_1 = x_2 \text{ and } e_1 = e_2 \\ \min(v(s_1), v(s_2)) & \text{if } x_1 > x_2 \text{ and } e_1 \le e_2 \text{ or} \\ x_1 \ge x_2 \text{ and } e_1 < e_2 \\ v(s_2) & \text{otherwise} \end{cases}$$

In Section 7 it will be showed that the performance of SQL improves significantly compared to the standard QL, both in the 1-step version here introduced and in the *n*-step version.

# 5.3 Threshold Q-learning

*Threshold Q-learning* (TQL) considers a simplified partial order of the type  $(x, e, z) \ge (x + 1, e, z')$  for x = 1, ..., M - 1: it imposes the partial order only w.r.t. energy. The monotonicity condition writes

$$q_*((x, e, z), a) \le q_*(([x+1]_M, e, z), a) \quad \forall x, e, z, a$$
(13)

The convergence follows from Theorem 2. The projection function with regard to a certain state s,  $\Pi_s : \mathbb{R} \to \mathbb{R}$ , can be written as

$$\Pi_{s_1}(v(s_2)) = \begin{cases} \max(v(s_1), v(s_2)) & \text{if } x_1 < x_2 \\ v(s_1) & \text{if } x_1 = x_2 \\ \min(v(s_1), v(s_2)) & \text{if } x_1 > x_2 \\ v(s_2) & \text{otherwise} \end{cases}$$
(14)

for every pair of states  $s_1 = (x_1, e, z)$  and  $s_2 = (x_2, e, z)$ . Note that in this case only states with same energy and server availability are compared: this significantly reduces the number of projection operations compared to SQL, especially when *B* is large. Despite its reduced complexity, in many cases this algorithm has showed similar convergence speed as SQL. Note how for both SQL and TQL it is not possible to guarantee the corresponding policy, e.g., an  $\epsilon$ -greedy policy, to have the threshold structure at each timesteps. However, the numerical results show that policy displays a threshold structure after a rather small number of iterations.

#### 5.4 Reinforce

One of the baselines algorithms used for comparison is Reinforce (RF) [31]. RF performs a basic policy-gradient iteration in the form

$$\theta_{t+1} = \theta_t + \alpha_t G_t \frac{\nabla_\theta \pi(a_t | s_t, \theta_t)}{\pi(a_t | s_t, \theta_t)}$$
(15)

where  $0 < \alpha_t < 1$  is a stepsize and  $G_t$  is the policy reward estimation. The gradient is operated on the following policy parametrization based on the threshold structure of the optimal policy

$$\pi (x, e, z) = \begin{cases} 0 & x < \lfloor \theta_e \rfloor \\ \frac{1}{1 + \exp\{k(\theta_e - x - 0.5)\}} & x = \lfloor \theta_e \rfloor \\ 1 & x > \lceil \theta_e \rceil \end{cases}$$
(16)

where k is a suitable hyperparameter.

# 6 MULTI-DEVICE MODEL

The model is now extended to the case of *N* devices connected to the edge server. Each device  $k \in \{1, ..., N\}$  follows the device-server model described in Section 3. At each timeslot the server polls device *k* with probability  $p_{Z,k}$ , where  $p_Z \in P_Z = \{p_Z \in [0,1]^N : \sum_{k=0}^N p_{Z,k} = 1\}$  and k = 0 denotes the null device, i.e., no polling. All the devices are assumed to be independent of each other: by doing so, the optimal policy of each device is not influenced by the state of the other devices at a given timestep; the interesting case of correlated harvesting or data samples is left as part of future works. Moreover, the polling decision of the server feedforward, i.e., the state of a device is only known once the device is polled, so that the polling action is independent of the state of devices.

Algorithm 2 Pseudocode of APPI

**Require:**  $\epsilon > 0$ 

1: initial polling distribution  $p_{Z,\text{new}}$ 

2: while  $|p_{Z,\text{old}} - p_{Z,\text{new}}| > \epsilon$ 

3:  $p_{Z,\text{old}} \leftarrow p_{Z,\text{new}}$ 

- 4: **Optimal policy learning:** find optimal policy  $\pi_k^*(p_{Z,\text{old}})$  for  $\forall k$
- 5: **Polling optimization:** find  $p_{Z,\text{new}} \ge 0$  such that

$$\begin{cases} p_{Z,\text{new}} = \arg \max_{p_Z} \sum_k R_{\gamma_k, \pi_k^*(p_{Z,k,\text{old}})}(p_Z) \\ \sum_k p_{Z,\text{new}} = 1 \end{cases}$$

6: end while

7: return  $p_{Z,\text{new}}$ 

The server seeks the optimal random polling distribution  $p_{Z,k}$  that maximizes the sum of the rewards of the devices, i.e.,  $R(p_Z) := \sum_{k=1}^{N} R_{\gamma_k}(p_{Z,k})$ , where  $R_{\gamma_k}(p_{Z,k})$  obeys to (11), subject to the polling capacity constraint, i.e.,

maximize: 
$$R(p_Z) = \sum_{k=1}^{N} R_{\gamma_k}(p_{Z,k})$$
 (MD)

subj. to: 
$$\sum_{k=0}^{N} p_{Z,k} = 1$$
 (17)

$$p_{Z,k} \ge 0 \quad k = 0, \dots, N$$

In order to solve problem (MD), since the discounted reward (11) cannot be computed (unless all the transition probabilities are known), stochastic approximation might be used to find a solution. Furthermore, as already observed in Section 4.2,  $R_Y(\cdot)$  cannot be assumed differentiable or concave, ruling out stochastic approximation methods requiring differentiable objective functions, such as SPSA [32]. Other direct methods with less strict requirements, such as Enhanced Localized Random Search do exist [33].

The proposed algorithm, namely the Alternating Polling and Policy Improvement (APPI) algorithm, is summarized in Alg. 2. It alternates two steps: 1) a *policy learning* step (line 4) for a given polling vector  $p_Z$ , and 2) a *polling optimization* step (line 5) optimizing  $p_Z$  for a given policy using stochastic approximation methods. For the policy optimization step it is sufficient to use one of the learning methods proposed in Sec. 5, e.g., Stairway Q-learning for the efficiency's sake. Conversely, as observed before, if the policy for each device is fixed, the objective function is  $\sum_k R_{Yk}, \pi_k (p_{Z,k})$  with  $\pi_k = \pi_k^*(p_{Z,k,old})$ , i.e., the optimal policy according to the previous value of the polling vector. As showed in Section 7.3, this algorithm is both faster and more accurate than general methods for the problem studied.

In order to discuss the polling optimization step, recall that the objective function is concave in  $p_Z$  (and so a.e. continuously differentiable) suggesting the use of stochastic gradient ascent methods of the Kiefer-Wolfowitz family [34]. In particular, the iterates of the algorithm write

$$p_Z^{n+1} = \Pi_\Theta \left( p_Z^n - \alpha_n \hat{g}_n \right) \tag{18}$$

where  $p_Z^n$  is the *n*th iterate of the parameter,  $\hat{g}_n$  represents an estimate of the gradient of the objective function,  $\{\alpha_n\}_n$  is a sequence converging to 0 and  $\Pi_{\Theta}$  is a projection on the N + 1-dimensional

simplex  $\Theta = \{p_Z \ge 0 | \sum p_{Z,k} = 1\}$ , i.e., the space of possible polling vectors, including the null action when polling is not active. The projection activates when  $p_Z \pm c_n \Delta_n$  lies outside the constraint set and reverts to the nearest point in  $\Theta$ .

The specific technique to determine  $\hat{g}_n$  in (18) is based on simultaneous random increments (SPSA) [32]. This technique estimates two increments, one in the positive and one in the negative direction. Due to the independence between different devices, each component of the gradient depends only on the parameter  $p_{Z,k}$  of the device considered. Let  $\hat{R}_{\gamma_k,\pi_k}$  denote the approximated reward for device k. The corresponding component of the gradient estimate writes

$$(\hat{g}_n)_k = \frac{\hat{R}_{Y_k,\pi_k} \left( p_{Z,k} + c_n \left( \Delta_n \right)_k \right) - \hat{R}_{Y_k,\pi_k} \left( p_{Z,k} - c_n \left( \Delta_n \right)_k \right)}{2c_n (\Delta_n)_k}$$
(19)

where  $\{c_n\}$  is a sequence converging to 0 and  $\{\Delta_n\}$  is an i.i.d. vector sequence of perturbations of i.i.d. components  $\{(\Delta_n)_i, i = 1, ..., N\}$  with zero mean and where  $\mathbb{E}\left[|(\Delta_n)_i^{-2}|\right]$  is uniformly bounded.

Before proving the convergence of the APPI algorithm, let show that fixed the policies of each device, the SPSA iteration converges to a global maximum of the corresponding total discounted reward. Actually, the convergence of depends on some condition on the objective function, on the step-size sequence  $\{\alpha_n\}$  and on the gradient estimates  $\hat{g}_n$ . The following holds:

PROPOSITION 1. Let  $\{\alpha_n\}$  and  $\{c_n\}$  be such that  $\sum_n \alpha_n = \infty$ ,  $\sum_n \left(\frac{\alpha_n}{c_n}\right)^2 < \infty$ . Moreover, assume that  $\mathbb{E}\left[|(\Delta_n)_i|^{-2}\right]$  is uniformly bounded on  $P_Z$  and that  $R_{Y_k,\pi_k}(p_{Z,k})$  is concave. Then the iteration (18) converges to the optimal parameter  $p_{Z,new}$  w.p.1.

PROOF. The desired result is obtained by verifying the assumptions of Proposition 1 in [35], whose proof is a consequence of Theorem 5.3.1 in [34].

i. the objective function is differentiable and either concave or unimodal: differentiability has been described in Remark **??**.

ii.  $b_n = \mathbb{E}[\hat{g}_n | p_Z, s_n] - \nabla J(p_Z) \rightarrow 0$  w.p.1: holds from Lemma 2 in [32] and condition i.;

iii.  $\sum_{n=1}^{\infty} \alpha_n^2 \cdot \mathbb{E}\left[e_n^2\right] < \infty$  w.p.1, where  $e_n := \hat{g}_n - \mathbb{E}\left[\hat{g}_n \mid \theta_n, s_n\right]$ ; this condition holds by Lemma 2 in [32] and by the fact that  $R((p_Z)_n, \omega)$ , representing the reward of sample path  $\omega$  has second moment uniformly bounded in  $p_Z$ : we can notice that it is independent on the parameter. Indeed, given a path, since  $x \leq M$  and  $p_Z \in [0, 1]$ , the reward is upper bounded by a constant, therefore its second moment is uniformly bounded, which concludes the proof.

Hence, fixed the policy for each device, the optimal polling vector is attained. Thanks to the convergence properties of the learning methods (see Section 5) for a fixed vector  $p_Z$ , and since the number of policies is finite, this proves that APPI converges w.p.1.

Unfortunately, this does not imply the convergence to the globally optimal polling vector, as the function to optimize is not concave, but rather just the pointwise maximum of a set of concave functions. Due to its particular structure, only convergence to the local optimum of the objective function can be ensured.

#### 7 NUMERICAL RESULTS

The numerical experiments are divided into 3 parts. The first one describes the performance of the device-server system tested on a

Parameter	Subsec. 7.1	Subsec. 7.2 and 7.3
М	15	$\{10, \ldots, 25\}$
В	15	$\{10, \ldots, 25\}$
γ	0.95	$\{0.9, \ldots, 0.99\}$
pz	0.05	$\{0.01, \ldots, 0.1\}$
δ	1	1
Reward	r(x) = M - x	r(x) = M - x
Discharge Penalty	$d(e) = e^4 \cdot \mathbb{1} \{ e \ge 0 \}$	$d(e) = \mathbb{1}_{e < 0} (-e)^k,$
		with $k \in \{1, 2, 3, 4\}$
Harvesting	Poisson ( $\lambda = 1$ )	Poisson ( $\lambda = 1$ )
Processing Cost	uniform, binary and	symmetric with
	symmetric for $\mu \in$	$\mu \in \{1, \ldots, B/2\}$
	$\{1, 4, 8.5\}$	$\sigma \in \{1, 2, 3, 4\}$

Table 2: The system parameters used in the numerical experiments.

range of system parameters. The second one compares the Ordered Q-learning methods introduced in Sec. 5 with baseline RL algorithms. Finally, the performance of the APPI algorithm introduced in Sec. 6 is assessed against alternative general methods. The devices and server parameters, namely  $M, B, \gamma, p_Z, r(\cdot), d(\cdot), p_H$  and  $p_C$ , constitute the *environment*. For the data processing cost *C* three possible probability distributions will be considered: the *uniform* one in  $\{1, \ldots, B+1\}$ , the *binary* one where  $p_C(1) = p_C(B+1) = 1/2$  and finally *symmetric* one obtained as  $p_C(c) = A \cdot \exp\{-\frac{1}{2}((c-\mu)/\sigma)^2\}$  for  $c \in \{1, \ldots, B+1\}$ :  $\mu$  and  $\sigma$  shape the distribution, whereas *A* is an appropriate normalization constant. In Table 2 the first column reports the sets from which the system parameters are drawn uniformly at random in the experiments of Subsec. 7.2 and Subsec. 7.3.

#### 7.1 Device-server performance evaluation

Fig. 3a describes the impact of the server availability  $p_Z$  onto the average (peak) AoI. Excluding the case  $\mu = 1$ , i.e., for lower values of the average batch processing cost, the average AoI is maximal for  $p_Z = 0$  and attains its minimum close to  $p_Z = 1$  (the optimal  $p_Z$  is highlighted with dots in Fig. 3a). In fact, for higher processing cost, a larger server polling rate permits to select action a = 1 more often, as the possible battery discharge due to data processing is compensated by frequent task offloading events. For lower cost figures, instead, it is the roundtrip delay  $\delta$  that renders the offloading less convenient: when  $\delta = 0$ , actually, the corresponding curve (not reported for the space's sake) becomes strictly decreasing.

Fig. 3b plots the value of the optimal polling rate  $p_Z$  for increasing values of  $\delta$ . The minima are obtained by means of an appropriate stochastic optimization. The results show, as expected, that the optimal value of  $p_Z$  decreases as  $\delta$  increases irrespective of the distribution of the processing cost. From Fig. 3b offloading is detrimental for larger values of  $\delta$ . This is more apparent for lower batch processing costs ( $\mu = 1$  and  $\mu = 3$ ). In those cases, there exists a threshold value of  $\delta$  above which offloading is detrimental, i.e.,  $p_Z = 0$ . In fact, for low processing costs, the risk of emptying the battery is sufficiently low to encourage to perform action a = 1 very frequently. In turn, the presence of the server, and the additional delay required by task offloading, becomes a penalty.



Figure 3: a) Average peak AoI for increasing polling probability  $p_Z$ ; b) optimal  $p_Z$  for increasing round-trip delay  $\delta$ ; c) Average peak AoI for increasing average harvesting rate  $\lambda$ .

Finally, Fig. 3c depicts the average AoI for increasing average harvesting rates, i.e., for increasing values of  $\lambda$  (note that here  $\lambda = 0$  indicates the deterministic case H = 1). Irrespective of the processing cost distribution  $p_C$ , the average Peak AoI is monotone decreasing, as expected. Interestingly, the AoI values appear rather insensitive to the cost distribution.

	250 episodes	1000 episodes	3000 episodes
QL	$0.723 \pm 0.254$	$0.796 \pm 0.220$	$0.810\pm0.208$
TQL	$0.752 \pm 0.255$	$0.857 \pm 0.181$	$0.880 \pm 0.117$
SQL	$0.850 \pm 0.209$	0.943 ± 0.051	0.964 ± 0.060
Reinforce	0.893 ± 0.225	$0.893 \pm 0.227$	$0.893 \pm 0.224$

 Table 3: RL tests: discounted reward for increasing number of episodes; best performance results are marked bold.

### 7.2 Learning the optimal policy

The RL algorithms introduced in Section 5 are compared with two baselines, namely QL and RF. Each test is repeated over 50 different environments where the system's parameters are drawn uniformly at random as reported in Table 2. Each experiment consists in a sequence of 1000 episodes. At the end of each episode, the discounted reward is calculated using a policy evaluation step. The evaluation is truncated when the weight discount falls below 0.001. The estimation of the optimal reward (10) is collected starting from a state  $s_0 = (1, e, z)$ , with *e* and *z* drawn uniformly at random. In order to compare results of different runs, the rewards have been normalized agains the baseline stationary policy which chooses action a = 1 if and only the state is such that either z = 1 or s = (M, B, 0).

Finally, for each tested algorithm several sets of hyperparameters are probed to determine the configuration ensuring the highest discounted reward; in fact, such parameters are observed to influence heavily the performance of the different algorithms. For the Q-learning type of algorithms (QL, TQL and SQL), the hyperparameters consist of *n*, determining the adopted *n*-steps variant, and  $\phi$  and  $\beta$ , which determine the learning rate  $\alpha_t(s_t) = \phi$ . (#visits in state  $s_t)^{-\beta}$ . For Reinforce the hyperparameter are *k*, appearing in (16), as well as *n* and *m* defining stepsize  $\alpha_t = n/(t+1)^m$ .

The results of the experiments are summarized in Table 3. In particular, RF appears to converge quickly, but to a suboptimal policy. In order to avoid this, in policy gradient RL it is possible to introduce multiple simultaneous state-action perturbations, e.g., trust-region techniques [36], at the price of increased complexity. On the other hand, SQL outperforms all other methods, attaining a gain of 10% on the average discounted reward.

#### 7.3 Multi-device system

The two last experiments evaluate the performance of the APPI learning procedure introduced in Sec. 6. To this aim, three alternative algorithms to be compared to APPI have been implemented. They replace the polling improvement step with Naive Random Search (NRS), Enhanced Random Search (ENRS) and SPSA, respectively. Details on those procedures are found in [33].

The first test covers a scenario with N = 3 devices. For each run of the experiment, a set of parameters is generated at random for each device (i.e., M, B,  $\gamma$ ,  $p_Z$ ,  $r(\cdot)$ ,  $d(\cdot)$ ,  $p_C$  and  $p_H$ ) and an initial polling distribution is imposed on the edge server. The device parameters are drawn uniformly at random from the sets described in Table 2, second column. For each algorithm it is recorded 1) the total discounted reward and 2) number of policy learning iterations performed. The results are averaged over over 50 runs, by considering each time different initial polling distributions and different environments. Also, all the reward values are normalized to fall into an interval  $[R_{\min}, R_{\max}]$ .  $R_{\max}$  is the reward obtained by an ideal APPI implementation which finds the optimal policy using value iteration, while R<sub>min</sub> is the reward obtained averaging 10 runs under a random polling distribution and the corresponding device's optimal policies. The numerical results are collected in Table 4: APPI consistently outperforms the other algorithms by attaining higher discounted reward and requiring much fewer policy improvement steps as well. The last experiment tests the scalability of APPI by increasing the number of devices N. The outcomes are reported in Table 5. Those are the average of 50 simulations per value of N. As before, in each simulation the environment is drawn at random using the distributions indicated in Table 2. The normalized reward

Algorithm	Discounted reward	Policy learning steps
NRS	$0.423 \pm 0.237$	$4.98 \pm 1.378$
ENRS	$0.678 \pm 0.184$	$10.06 \pm 3.331$
SPSA	$0.449 \pm 0.276$	$5.54 \pm 1.846$
APPI	0.747 ± 0.143	2.98 ± 1.086

**Table 4:** Comparison of different polling improvement algorithms for N = 3.

Ν	Discounted reward	Policy learning steps
1	0.794 ± 0.222	2.18 ± 0.77
3	$0.747 \pm 0.143$	$2.98 \pm 1.086$
5	$0.699 \pm 0.192$	$3.78 \pm 1.221$
7	$0.724 \pm 0.222$	$4.08 \pm 1.074$
10	$0.714 \pm 0.237$	$5.06 \pm 1.302$

 Table 5: Performance of APPI for increasing number of devices connected to the edge server.

attained by APPI appears marginally affected by the number of devices, whereas the number of required policy improvement steps is in the order of a few units and increases linearly.

### 8 CONCLUSIONS

Based on the peak AoI this work has provided a theoretical framework to model edge devices running AI applications whose energy footprint depends on the data being processed. It factors in energy harvesting and edge-server task offloading to determine the optimal processing policy on edge devices. The solution approach builds on the properties of the Markovian model to derive the key structural features of optimal policies. This is the basis for a model-based, specialized reinforcement learning method, namely Ordered Q-learning which is extended to the multidevice setting. There, the server's offloading rate vector is optimized by alternating stochastic gradient ascent and optimal policy learning steps. Numerical tests performed against the ground-truth, i.e., the actual optimal policy, show that existing baselines are systematically outperformed both in accuracy and convergence speed. This confirms that by rooting reinforcement learning techniques into the structural properties of the optimal solution, important performance margins against generic RL solutions are attained. Several extension of the proposed models are indeed possible in the multi-device setting. In particular, it is possible to include capacity constraints in the Markovian model, e.g., to model non-ideal communication channels. This in turn requires to study and learn optimal randomized stationary policies in the context of constrained MDPs. Also, in order to improve the polling learning procedure at the edge-server, one could use the history on the device state which can be retrieved at polling instants.

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