

# PACKED-BED THERMAL-STORAGE SYSTEM CONFIGURATIONS FOR SHUTTLE KILN WASTE-HEAT RECOVERY

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# ABSTRACT

The ceramic industry, which heavily relies on non-renewable energy sources, significantly contributes to carbon dioxide emissions. In particular, periodic or shuttle kilns, used in sanitaryware production, are batch-type furnaces and are predominantly fueled by natural gas. The produced waste heat is often rejected into the environment. This study compares three distinct system configurations for waste heat recovery, which are namely: a standalone heat exchanger (HEX), crossflow or counterflow, a packed-bed thermal storage (PBTS) system for sensible heat storage, and a combination of both. A validated 1-dimensional numerical model, which solves the energy conservation equation for both solid and fluid phases is used for analyzing the PBTS system. The HEX alone recovers up to 16% of total heat input, achieving a 0.93-year payback period and a levelized cost of heat at  $0.017\epsilon/kWh_{th}$ . Conversely, an optimally dimensioned PBTS system surpasses HEX efficiency, reclaiming 28.5% of the heat input. It boasts a shorter payback period (0.17-0.32 years) and lower levelized cost (0.003-0.006 $\epsilon/kWh_{th}$ ), demonstrating its superiority in enhancing waste heat recovery in industrial settings.

# **1** INTRODUCTION

To meet the goals set by the "Paris Agreement", technologies for the decarbonization of the industrial sector through the recovery of produced waste heat (WH) have received significant scientific attention. Specifically, in the ceramic industry, sanitaryware production contributes to approximately 3% of the ceramic sector's total energy consumption in the EU and this energy is mainly covered by natural gas (NG) consumption (EUROPEAN COMISSION, 2007). Periodic kilns or shuttle kilns (SKs) are batchtype furnaces that operate with distinct heating and cooling trajectories and are commonly used for sanitaryware production. It has been reported that one-third of the waste heat generated in such manufacturing sites can be recovered, as currently the exhaust gas energy (around 90% of the total inlet heat) is rejected into the environment (Cuviella-Suárez et al., 2021). Regarding industrial waste heat recovery (WHR), different technologies can be used, like active systems (heat pumps, power cycles) or passive systems (heat exchangers, thermal energy storage) (Miró et al., 2016). WHR in the context of the ceramic industry has been considered by Kaushik and Shah (2016) where the option of recuperating waste heat of a roller kiln for tiles production for usage at the facility's dryer was examined. Mezquita et al. (2014) furthermore investigated the energy-saving potential of a tunnel kiln through direct gas recirculation. However, studies on WHR through thermal energy storage and specifically for the ceramic industrial sector are generally lacking in the literature.

The novelty of this research is that it focuses on industrial WHR within the context of a SK by integrating a packed-bed thermal storage system (PBTS), designed for sensible heat storage (SHS), in the production process. The PBTS system will be charged with the exhaust gas heat from the SK's cooling phase and discharged for the preheating of the combustion air during the SK's heating phase.

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This study will compare 3 different system configurations for WHR and in all cases the WH will be reused for preheating the combustion air and thus lowering the natural gas (NG) consumption. In Section 2 the SK operation and data processing, as well as the methodologies used to analyze the counterflow and crossflow heat exchangers (HEXs) and the PBTS system, are presented. In Section 3 the performance of each WHR scenario in terms of their technical performance (WHR and thermal operation) and from a techno-economic perspective is examined. Conclusions highlight the effectiveness of the PBTS system in enhancing WHR, suggesting future research directions for optimization and broader application.

## 2 MATERIALS AND METHOD

In this section the SK's basic operation is discussed, emphasizing the WHR potential. Next, the processed data for a SK, based on Cuviella-Suárez *et al.* (2021), are presented. Then, the HEX and PBTS analysis methodology is explained together with underlying assumptions. For the PBTS system, a numerical model is presented, which is validated based on experimental measurements derived from a literature study.

### 2.1 SK Basic Operation

To analyze either the application of a HEX or a PBTS for WHR in a SK, it is necessary to study the mass and energy balance during the SK operation. Based on the study of Cuviella-Suárez et al. (2021) the energy balance of a SK, the NG calorific value ( $H_B$ =53.5GJ), the combustion under 10% of excess air and the stoichiometric air-to-NG mass ratio  $\mu_{NG}$ =17.16 for combusting 1kg of NG, were used. The standard SK configuration includes burners combusting NG ( $\dot{m}_{NG}$ [kg/s]) with combustion air  $(m_{comb}[kg/s])$ , along with diffusion  $(m_{dif}[kg/s])$  and injection  $(m_{inj}[kg/s])$  air to control the temperature inside the SK ( $T_{inside}$  [°C]). In total 52.798GJ of heat ( $Q_{react,tol}$ ) are provided by NG combustion for a single operation cycle. The basic SK operation lacks WHR technology, releasing all waste heat, from both the heating (HP) and cooling phases (CP) into the environment, which are 21.350GJ and 26.855GJ respectively. Three WHR scenarios are investigated. Scenario 1 (Sc.1) involves a recuperative HEX, operating during the HP, for preheating ambient temperature  $(T_{amb})$  combustion air ( $\dot{m}_{comb}$ [kg/s]) up to temperature  $T_{preheated}$  by using heat from the exhaust gases ( $\dot{m}_{exh}$ [kg/s]) of temperature  $T_{exh}$ . Scenario 2 (Sc.2) features only a PBTS storage tank charged with heat from SK's CP, of temperature  $T_{inside} = T_{exh}$ , which is utilized for preheating combustion air during the next HP. Scenario 3 (Sc.3) combines both a PBTS and a recuperative HEX for maximum WHR. A schematic of the SK basic operation and the 3 WHR scenarios is provided in Figure 1.



Figure 1. Schematic diagrams for the SK basic operation and the investigated 3 WHR scenarios components (indicated by a different color), during the heating (left) and cooling phase (right).

### 2.2 SK Data Processing

Operational data from the SK were analyzed to approximate the  $\dot{m}_{comb}$ ,  $\dot{m}_{exh}$  and exhaust gas temperature  $(T_{exh}[^{\circ}C])$  for each moment during its operation. Based on Cuviella-Suárez *et al.* (2021),

specific heat amounts over distinct time intervals ( $\Delta t_i$ [min]) for both the HP and CP, with 14 and 10 intervals respectively, were utilized. The  $\Delta t_i$  intervals are non-equal time steps. During the HP, given heat amounts are the combustion heat ( $Q_{comb}$ [kJ]), which is a sum of various terms like heat losses to the environment, heat absorbed by the ceramic products, diffusion, and injection heat), and exhaust gas heat ( $Q_{exh}$ [kJ]). The authors assert that  $Q_{comb}$  and  $Q_{exh}$  collectively represent the total heat input from NG consumption ( $Q_{react}$ [kJ]).

NG consumption  $(m_{NG}[kg])$  at each stage is determined based on  $Q_{react}$  and the  $H_B$ . Based on the excess air combustion and  $\mu_{NG}$  the combustion air masses  $(m_{comb}[kg])$  are computed for every  $\Delta t_i$ . Injection  $(m_{inj}[kg])$  and diffusion air masses  $(m_{dif}[kg])$  are calculated assuming that the given heat values for  $Q_{dif}[kJ]$  and  $Q_{inj}[kJ]$  represent the heat needed for inlet masses  $(m_{inj} \text{ and } m_{dif})$  to rise from ambient temperature to the SK inside temperature  $(T_{inside})$ .  $T_{exh}$  is calculated based on the energy conservation equation at the exhaust of the SK, using the  $Q_{exh}$  and  $m_{exh}$ , where  $m_{exh}$  is the sum of  $m_{NG}$ ,  $m_{comb}$ ,  $m_{dif}$  and  $m_{inj}$ . For the CP, masses of cooling, injection, and diffusion air are computed similarly to the HP for  $m_{inj}$  and  $m_{dif}$ . These masses constitute CP  $m_{exh}$ , with  $T_{exh}$  assumed to be equal to  $T_{inside}$  during the CP.

The process concludes with the calculation of the required instant mass fluxes ( $\dot{m}_{comb}$  and  $\dot{m}_{exh}$ ) for 24 time moments ( $i_i$ ) using backward time difference, for the non-equal time step time intervals. Also, for every  $\Delta t_i$ , the calculated  $T_{exh}$  values are taken for the corresponding  $i_t$  time moment. The ambient temperature is assumed  $T_{amb}=20^{\circ}$ C. The calculated air and NG total masses showed an acceptable relative error compared to the values of Cuviella-Suárez *et al.* (2021), i.e. below 2%. Regarding the total combustion air the relative error was around 7%. Finally, the calculated points for the 24  $i_t$  time moments of the SK operation, for the air and NG mass fluxes are shown in Figure 2 a. and in Figure 2 b for the temperatures inside the SK ( $T_{inside}$ ) and for the exhaust gas ( $T_{exh}$ ). The calculated points were interpolated using cubic-b splines to obtain all the time plots of Figure 2.



**Figure 2.** Calculated diagrams of **a**. air and NG mass fluxes and **b**. the temperature inside the SK  $(T_{inside})$  and of the exhaust gas  $(T_{exh})$ , during the operation of the SK (for the cooling phase  $T_{exh}=T_{inside}$ ).

### 2.3 HEXs Designs and Analysis

The HEX operates during the SK HP and preheats the  $\dot{m}_{comb}$  (air) using the heat of  $\dot{m}_{exh}$  (flue gas). The temperatures at the inlet of the HEX for the  $\dot{m}_{comb}$  and  $\dot{m}_{exh}$  are  $T_{c1} = T_{amb}$  and  $T_{h1} = T_{exh}$ , while at the outlet are  $T_{c2}$  (= $T_{preheated}$ ) and  $T_{h2}$ , respectively (see Figure 1). Two HEX designs are considered: a crossflow, as they are used for gas-to-gas applications, and a counterflow to provide a theoretical threshold for the operation of a HEX as it can achieve maximum effectiveness. The product UA of HEX area (A) and heat-transfer coefficient (U) is used as the design variable and considered in the range of 1-20kW/K. Heat exchange between  $\dot{m}_{exh}$  and  $\dot{m}_{comb}$  is determined via the effectiveness-NTU method for both HEX designs (Komninos and Rogdakis, 2018). An operational constraint is set for the HEX, allowing their operation only as long as  $T_{exh}$  is below 1000°C (Kautz and Hansen, 2007). This means that the HEX will operate up to the time moment of 414min ( $T_{exh}$ =1056°C). For the techno-economic analysis, it is essential to calculate both the capital expenditure (*OPEX*) and operational expenditure (*OPEX*) of the investigated components. The *CAPEX* of the HEX is also computed as a

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function of the UA, with OPEX being estimated at 20% of CAPEX (Cocco and Serra, 2015; Georgousis et al., 2022).

### 2.4 PBTS Storage Tank Geometry, Materials and Assumptions

The PBTS system under investigation consists of a cylindrical storage tank and uses air as the HTF. Spherical Electric Arc Furnace (EAF) steel slag packing elements (PEs) are used to store heat with a density of  $\rho_s$ =3430kg/m<sup>3</sup> and heat capacity of  $c_{p,s}$ =933J/kgK=const (Kocak *et al.*, 2021). The following PBTS design variables are considered constant: the aspect ratio (a=H/D=1), the PEs sphere diameter ( $d_p$ =1cm) and the PBTS porosity ( $\varepsilon$ =0.39). Diameter D(=H) is taken in the range of 2.25m-3.5m and the charging time ( $t_{ch}$ ) varies between 150 and 375 minutes. During the PBTS charging phase (SK CP), hot exhaust air at  $T_{exh}$  of the SK charges the PBTS. The heat remains stored in the PBTS in the subsequent stand-by phase, which lasts  $t_{stb}$ =530min. The final discharging phase (SK HP) has a time duration of  $t_{dis}$ =535min. The aforementioned 3 consecutive phases correspond to 1 PBTS operational cycle, which lasts  $t_{ch}+t_{stb}+t_{dis}$ =24hr. Temperature  $T_{exh}$  keeps decreasing during the charging phase (Figure 2 b). For that reason, the HTF inlet in the PBTS for both the charging and discharging is from the same point (Z=0m). This way the stored heat of higher temperatures will be discharged at the beginning of the discharge phase.

The PBTS walls are considered adiabatic. Also, buoyancy-induced flow effects due to differences in air density are not considered (Schwarzmayr *et al.*, 2023). Therefore, the temperature distribution inside the PBTS can be considered 1-dimensional (1D). The calculation of the *CAPEX* includes the storage tank cost ( $C_{ST}$  [€]) and PEs specific cost ( $C_{PE}$ =80 €/tn) (Kocak *et al.*, 2021).  $C_{ST}$  is calculated using a constant cost per m<sup>3</sup> of the storage tank ( $C_V$  [€/m<sup>3</sup>]). Literature provides varying  $C_V$  values and the two, lower and upper limits that were chosen are  $C_{V,I}$ =745€/m<sup>3</sup> (Cocco and Serra, 2015) and  $C_{V,2}$ =1580€/m<sup>3</sup> (Strasser and Selvam, 2014). *OPEX* was calculated as in Section 2.3.

### 2.5 Numerical Model for PBTS System and Validation

For analyzing the performance of the PBTS, the temperature profiles during the charging, stand-by and discharging phases for both the HTF and solid phases need to be calculated. For the computation of the temperature profiles a 2Phase–1D (2P-1D) numerical model is used, which is based on the energy conservation equation for the solid and HTF phases using a Darcy-type approach. For air-solid PBTS systems the density of the HTF ( $\rho_f$ ) and the heat capacity of the solid phase ( $c_{ps}$ ) are taken temperature dependent, whilst the other properties are assumed constant and computed at a reference temperature. The pressure drop is considered to have a negligible effect on the HTF density. The set of equations describing the problem consists of the mass and 2P-1D energy conservation equations (Ortega-Fernández and Rodríguez-Aseguinolaza, 2019):

HTF mass conservation: 
$$\varepsilon \frac{\partial \rho_f}{\partial t} + \frac{\partial (\rho_f u_D)}{\partial x} = 0$$
 (1)

HTF energy conservation:  $\varepsilon (\rho c_p)_f \frac{\partial T_f}{\partial t} + \rho_f c_{p,f} u_D \frac{\partial T_f}{\partial x} = k_{f,eff} \frac{\partial^2 T_f}{\partial x^2} + a_p h_p (T_s - T_f)$  (2)

PEs energy conservation: 
$$(1 - \varepsilon)\rho_s \frac{\partial (c_p T)_s}{\partial t} = k_{s,eff} \frac{\partial^2 T_s}{\partial x^2} - a_p h_p (T_s - T_f)$$
 (3)

Air state equation:

$$\rho_f = \frac{P_{ref}}{\bar{R}\,T_f} \tag{4}$$

Here  $u_D$  [m/s] is the HTF Darcy velocity,  $\varepsilon$  [-] is the porosity of the PBTS, subscripts "s" and "f" correspond to the solid PEs and HTF, respectively, T [K] is the temperature of either the PEs or HTF,  $k_{f,eff}$  and  $k_{s,eff}$  [W/mK] are the effective axial thermal conductivities of the 2 phases,  $\rho$  [kg/m<sup>3</sup>] and  $c_p$  [J/kgK] are the density and specific heat,  $h_p$  [W/mK] is the convective heat transfer coefficient between the 2 phases and  $a_p=6(1-\varepsilon)/d_p$  [1/m] is the "shape factor" of the PEs. The  $k_{f,eff}$  and  $k_{s,eff}$  are calculated based on the stagnant HTF effective conductivity ( $k_{stag}$ ), while the calculation of  $k_{f,eff}$  also includes the HTF thermal dispersion conductivity ( $k_{f,dis}$ ) due to the flow through the porous structure (Ismail and Stuginsky, 1999, Yang and Nakayama, 2010). Depending on the Reynolds number of the HTF flow,

defined as  $Re = u_D \rho_f d_p / \mu_f$  where  $\mu_f$  [Pa/s] is the HTF's dynamic viscosity, different Nusselt correlations were used for the calculation of  $h_p$  (Ortega-Fernández and Rodríguez-Aseguinolaza, 2019). Relevant air properties were calculated using the Coolprop package available in Python 3.11.

In PBTS systems, during their operation, a temperature gradient zone, called "thermocline" (TC), is formed for the HTF and the PEs respectively. Inside the TC, of each phase, the temperature varies from the "high" temperature ( $T_H$ ) to the "cold" temperature ( $T_C$ ) level. During the (dis)charge of the PBTS the TC zones dynamically propagate inside the storage tank. The progression of the TC zones for the charging phase is illustrated in Figure 3 over a time duration  $\Delta t$  (the solid blue lines correspond to the HTF temperatures and the dashed yellow to the PEs). The "thermocline velocity" ( $u_{TC}$ ) can be defined as the characteristic velocity at which the center of the region, enclosing both the HTF and PEs TCs, moves. The  $u_{TC}$  definition equation for incompressible HTF flow and constant properties can be found in literature (Yang and Garimella 2010). For air-solid PBTS systems where the  $\rho_f$  and  $c_{p,s}$  are temperature-dependent, a first-order approximation of  $u_{TC}$  can be derived by assuming constant mass flow rate ( $\dot{m}_f = \rho_f u_D A = const$ ), as done by Yang and Garimella (2010):



**Figure 3.** Schematic illustration of charging a PBTS and the thermocline progression, over a time scale  $\Delta t = t^{m+1} - t^m$ . HTF temperature: solid blue lines, PEs temperature: dashed yellow lines.

Here  $\dot{m}_f$  [kg/s] is the HTF mass flux and A [m<sup>2</sup>] is the PBTS cross sectional area. Based on this relationship and considering that  $\rho_s \approx 10^3 \rho_f$ , it can be easily deduced that  $u_{TC} \approx 10^{-3} u_D$ . The distance that the TC crosses based on  $u_{TC}$  over a time duration of  $\Delta t$  is equal to  $\Delta x_{TC} = u_{TC} \times \Delta t$ . Using as a reference the temperature measurements from Esence *et al.* (2019) for charging, it can be computed that  $u_{TC} = 1.29 \ 10^{-4} \text{m/s}$ . This means that the TC propagates by  $\Delta x_{TC} = 0.31 \text{m}$  in a 40min duration. This aligns with the experimental temperature distribution measurements (Figure 4 a.), validating Equation (5) as an initial  $u_{TC}$  estimate for PBTS systems using air/flue gas as HTF and variable  $c_{p.s}$ .



Figure 4. a. Experimental measurements of PEs temperatures for "CLAIRE" (Esence *et al.* (2019)) (only the points) and b. 2P-1D numerical model PEs temperature prediction for "CLAIRE" for 6 different time moments (triangle points: experimental measured solid temperatures of corresponding time moment, 1<sup>st</sup> approach: solid lines, 2<sup>nd</sup> approach: dashed lines).

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Considering the HTF mass conservation (Equation (1)) and a point in the TC region (see Figure 3) the first term T1 is of the order:

$$T1: \varepsilon \frac{\partial \rho_f}{\partial t} \approx \varepsilon \frac{\Delta \rho_f}{\Delta t} \approx \varepsilon \, u_{TC} \frac{\Delta \rho_f}{\Delta x_{TC}} \approx \ 10^{-4} u_D \frac{\Delta \rho_f}{\Delta x_{TC}} \tag{6}$$

The second term T2 in the HTF mass conservation (Equation (1)) is of the order:

T2: 
$$\frac{\partial(\rho_f u_D)}{\partial x} = u_D \frac{\partial(\rho_f)}{\partial x} + \rho_f \frac{\partial(u_D)}{\partial x} \approx u_D \frac{\Delta(\rho_f)}{\Delta x_{TC}} + \rho_f \frac{\Delta(u_D)}{\Delta x_{TC}}$$
(7)

It can now easily be seen from Equations (6) and (7) that T1 is negligible compared to T2, so Equation (1) becomes:

$$\frac{\partial(\rho_f u_D)}{\partial x} = 0 \Leftrightarrow \rho_f u_D = \frac{\dot{m}_f}{A} = const$$
(8)

By using this simplified mass conservation equation for the HTF (Equation (8)), Equation (2) and (3) can be solved independently, they are de-coupled from Equation (8). The axial dimension – height of the PBTS (*H*) is discretized in an equidistant grid from node i=1 to i=N, where the total number of nodes N was taken in the range of 650-1000. To discretize the 2P-1D energy conservation equations, an implicit Euler scheme was used for the time derivatives and central differences for the spatial derivatives. The time step  $\Delta t$  is considered between the present time moment and the previous time moment, which are indicated by the superscripts "m+1" and "m" respectively.  $\Delta t$  was taken in the range of 5-10s. For many solid phase materials,  $c_{p,s}$  highly depends on the temperature. The same holds for the  $\rho_f$  of the HTF. To model the temperature dependency 2 different approaches were compared for the temperature dependent  $c_{p,s}$  term of Equation (3), which are:

$$\frac{\partial(c_p T)_s}{\partial t} = \frac{\int_{T_{s,i}}^{T_{s,i}^{m+1}} c_{p,s} dT}{\Delta t} \approx \begin{cases} c_{p,s}^{m+1} \frac{T_{s,i}^{m+1} - T_{s,i}^m}{\Delta t} \left[ 1^{st} approach \right] \\ c_{p,s}^m \frac{T_{s,i}^{m+1} - T_{s,i}^m}{\Delta t} \left[ 2^{nd} approach \right] \end{cases}$$
(9)

Where *i* is a random node in the axial direction of the PBTS. In the 1<sup>st</sup> approach the heat capacity  $c_{p,s}$  is evaluated at time level m+1 using an iterative process until convergence, in the 2<sup>nd</sup> approach at time level m. Regarding temperature dependency of the HTF density ( $\rho_f$ ), which influences the volumetric term of Equation (2), when using the 1<sup>st</sup> approach (Equation (9)),  $\rho_f$  is evaluated at time level m+1 through the conducted iterative process and with the 2<sup>nd</sup> approach at time level m. With respect to the boundary conditions (BC), the inlet HTF node i=1 uses a known/given temperature ( $T_f = T_{given}$ ) (Dirichlet type BC), while at the outlet node i=N, a 0-value Neumann type BC ( $\partial T_f/\partial z = 0$ ) is implied. The PEs have a 0-value Neumann BC ( $\partial T_s/\partial z = 0$ ) at the PBTS top and bottom nodes. The stand-by phase features 0-value Neumann BC for both HTF and PEs (so no heat losses to the environment).

The charging phase temperature measurements of the PBTS experimental set up called "CLAIRE" of Esence *et al.* (2019) are used for the validation of the 2P-1D numerical model. Initial conditions temperature measurements (0min-black points in Figure 4 b.) were interpolated using cubic-b splines over the PBTS height (*H*) assuming also that  $T_{f,i}=T_{s,l}$  for i=1,...,N for the initial condition. The values produced from the interpolation were then used as initial condition for the 2P-1D numerical model. Figure 4 b. depicts the comparison of the PEs temperature distribution obtained from the 2P-1D numerical model with the two different approaches for the heat capacity of the PEs (Equation (9)), as well as the  $\rho_f$ , with the experimental measurement results of Esence *et al.* (2019). Six different experimental time moments are presented, ranging from 40min up to 235min. Both approaches of Equation (9) result in almost equal PEs temperature profiles between them and with good comparison to the experimental measurements can be seen only at the later stages (200min and 235min). Conclusively, the 2P-1D numerical model gives reliable results and the  $\rho_f$ .

#### 2.6 Parameters and KPIs Used

The following three key performance indicators (KPIs) were used to evaluate the thermal performance of the WHR equipments (HEX and/or PBTS) and their techno-economic performance. The first KPI is the recovered waste heat for  $\dot{m}_{comb}$  preheating, per SK operation cycle (*WHR*<sub>cycle</sub>[GJ]). The other 2 KPIs are the equipment's payback period (*PBP*[years]) and levelized cost of recovered heat (*LCOH* [€/kWh<sub>th</sub>]) (Cocco and Serra, 2015). The *PBP* was calculated based on:

$$PBP = \frac{ln(\frac{CF}{CF - r CAPEX})}{ln(1+r)}$$
(10)

Where r is the discount rate (7%), and CF is cash flow due to the reduction of the NG consumption. The calculation of the CF is done with the equation:

$$CF = WHR_{cycle}N_{annual}C_{NG} - OPEX = WHR_{annual}C_{NG} - OPEX$$
(11)

Where  $N_{annual}$ =232 are the operational annual SK cycles,  $C_{NG}$ =0.0659€/kWh<sub>th</sub> the NG heat cost for nonhousehold consumers in Netherlands and  $WHR_{annual}$ [GJ](= $WHR_{cycle}N_{annual}$ ) is waste heat that is being recovered annually. For the *LCOH* an estimation regarding the lifetime of the equipment is needed (N). A value of N=20 years was assumed. The *LCOH* was defined as:

$$LCOH = \frac{CAPEX + \sum_{n=1}^{N} \frac{OPEX}{(1+r)^n}}{\sum_{n=1}^{N} \frac{WHR_{annual}}{(1+r)^n}}$$
(12)

To compare the performance of the investigated HEXs with a PBTS, an optimal UA-value needs to be determined for either the crossflow or counterflow HEX. To achieve that, a multi-objective optimization criterion is used to optimize the KPIs defined above (i.e. maximize  $WHR_{cycle}$  and minimize PBP and LCOH minimize). First, the theoretical optimum design point or Ideal Point (IP) of the HEX is determined, characterized by max: $WHR_{cycle}$ , min:PBP and min:LCOH. The optimal solution is obtained based on the minimization of a multi-objective function "F", defined as the Euclidian distance between a HEX Design Point (DP) and the IP, based on the following equation (Georgousis *et al.*, 2022):

$$F(UA) = \sqrt{\left[\frac{WHR_{cycle}(UA) - WHR_{cycle,max}}{WHR_{cycle,max} - WHR_{cycle,min}}\right]^2 + \left[\frac{PBP(UA) - PBP_{min}}{PBP_{max} - PBP_{min}}\right]^2 + \left[\frac{LCOH(UA) - LCOH_{min}}{LCOH_{max} - LCOH_{min}}\right]^2$$
(13)

Where the subscripts min/max indicate the respective minimum/maximum values. An additional useful parameter is the average preheated temperature ( $\bar{T}_{preh}$ ) of the total combustion air ( $m_{comb,tot}$ =19.7tn), either when using a HEX or a PBTS or both, during the SK HP.  $\bar{T}_{preh}$  is defined as:

$$WHR_{cycle} = m_{comb,tot}\bar{c}_{p,air}(\bar{T}_{preh} - T_{amb}) \tag{14}$$

## **3 RESULTS**

### 3.1 Scenario 1: Implementation of Crossflow and Counterflow HEX

Starting with a parametric analysis of varying the UA value in the range of 1-20kW/K for a crossflow and counterflow HEX, the resulting values of the WHR<sub>cycle</sub>, CAPEX, PBP and LCOH are calculated (Figure 5). To select an optimum UA value for the crossflow and counterflow HEX, the multi-objective optimization criterion of Equation (13) is used. The optimal solution for crossflow is F=0.528, where UA=3.49kW/K and {WHR<sub>cycle</sub>, PBP, LCOH}={8.17GJ, 1.00years, 0.0182€/kWh<sub>th</sub>}. For the counterflow, the optimal solution is F=0.018, where UA=3.72kW/K and {WHR<sub>cycle</sub>, PBP, LCOH}={8.54GJ, 0.93years, 0.0170€/kWh<sub>th</sub>}.

For the optimum solutions of UA=3.49kW/K (crossflow) and UA=3.72kW/K (counterflow), the temperature diagrams of the  $\dot{m}_{comb}$  and  $\dot{m}_{exh}$  during the SK HP are presented in Figure 6. The diagrams show the preheated  $\dot{m}_{comb}$  temperature ( $T_{preheated}=T_{c2}$ ) and the  $\dot{m}_{exh}$  linet ( $T_{exh}=T_{h1}$ ) and outlet ( $T_{h2}$ )

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temperatures during the SK HP. The  $T_{c2}$  and  $T_{h2}$  lines that correspond to the optimum UA counterflow HEX (darker colors) are over and below the respective  $T_{c2}$  and  $T_{h2}$  lines of the optimum UA crossflow HEX (lighter colors). This result is due to the better heat exchange achieved with the counterflow HEX than crossflow HEX, which also leads to higher total WHR<sub>cycle</sub>.



**Figure 5. a.** *PBP* (blue lines) and *WHR*<sub>cycle</sub> (red lines) as function of *UA* and **b.** *LCOH* (blue lines) and *CAPEX* (red line-shared cost for both HEX types) as function of *UA* for crossflow HEX: solid lines, counterflow HEX: dashed lines.



**Figure 6.** Temperature diagrams for optimum crossflow (UA=3.49kW/K) and counterflow (UA=3.72kW/K) HEXs. Dashed orange line: inlet temperature ( $T_{h1}$ ) of  $\dot{m}_{exh}$ , solid orange lines: outlet temperature ( $T_{h2}$ ), solid blue lines: preheated  $\dot{m}_{comb}$  temperature ( $T_{c2}$ ).

# 3.2 Scenario 2: Implementation of a PBTS System

In Sc.2 the PBTS system is charged during the SK CP by  $\dot{m}_{exh}$  and discharged during the HP by  $\dot{m}_{comb}$ . A parametric investigation is conducted for variable storage tank diameters D(=H) ranging from 2.25-3.5m ( $t_{ch}$ =375min). Figure 7 shows the *PBP*, *WH*<sub>stored</sub>, *WHR*<sub>cycle</sub>, *LCOH* and *CAPEX* as a function of diameter *D*. Optimal performance is achieved at *D*=3.3m, maximizing *WHR*<sub>cycle</sub> to 13.637GJ and minimizing *PBP* to 0.47 years (for  $C_{V,1}$ ) and to 0.91 years (for  $C_{V,2}$ ), alongside achieving the minimum *LCOH* of 0.009 /kWh<sub>th</sub> (for  $C_{V,1}$ ) and 0.017€/kWh<sub>th</sub> (for  $C_{V,2}$ ). Figure 8 presents the temperature distribution of the HTF and PEs (solid and dashed lines respectively) at the end of the charging, standby and discharging phases inside the PBTS for 4 different diameters. The HTF and PEs temperatures are quite close at the end of the mentioned 3 operation phases.

Analysis of Figure 7 a. and Figure 8 b.-d., reveals that for PBTS systems with D(=H)>2.40m, full recovery of the stored heat ( $WH_{stored}$ ) is not feasible due to the insufficient air mass available for discharge ( $m_{dis,tol}=m_{comb,tol}=19.7$ tn). This limitation is visualized through the TC region not crossing the entire PBTS height (H) and remaining inside the PBTS at the end of the discharge phase (Figure 8 b.-d.). This causes some stored heat to remain unused by the end of the discharge phase. Moreover, increasing D(=H) increases  $WH_{stored}$ , as more heat can be stored, yet not  $WHR_{cycle}$ , as shown in Figure 7a. Initially, as the storage tank size increases from D=2.25m,  $WHR_{cycle}$  rises because higher-temperature heat is stored and recovered by  $m_{comb}$  (Figure 8 a. and b.). At the optimal design point

(D=3.3m), most of the TC is flushed out by the end of the charging phase (Figure 8b.), and also  $WHR_{cycle}$  peaks. For D>3.3m, despite increasing  $WH_{stored}$ ,  $WHR_{cycle}$  decreases (Figure 7a). Two reasons account for this. First,  $m_{dis,tot}$  cannot fully discharge such PBTS storage tank sizes, as was mentioned before. Second, when D=H increases, the part of the TC stored inside the PBTS (Figure 8 c, d) and eventually discharged also increases. This leads to preheating  $m_{comb}$  at lower temperature. Especially, for D=H=3.75m, although all WH produced during the CP could be stored ( $WH_{stored}$ =26.855GJ), only 2.12GJ could be recovered, given that only part of the TC was eventually discharged.



**Figure 7. a.** *PBP* (blue lines) for  $C_{V,2}=1580$  (m<sup>3</sup> (solid line) and  $C_{V,1}=745$  (m<sup>3</sup> (dashed line) and heat amounts (red lines) as a function of *D* and **b.** *LCOH* (blue lines) and *CAPEX* (red lines) for  $C_{V,2}=1580$  (m<sup>3</sup> (solid lines) and  $C_{V,1}=745$  (m<sup>3</sup> (dashed lines) as a function of *D*.



Figure 8. PBTS temperature profiles at the end of each phase for diameters (D=H) a. 2.25m, b. 3.30m, c. 3.50m, and d. 3.75m. Solid lines for HTF, dashed for PEs.

For optimal operation of the PBTS system the max: $WHR_{cycle}$ , min:*PBP* and min:*LCOH* must be achieved and the PBTS storage tank is designed/sized so that it can be fully discharged by  $\dot{m}_{comb}$ . The design criterion to achieve a complete discharge of the PBTS is that the TC center fully crosses the PBTS height (*H*) during discharge. The average discharge thermocline velocity ( $\bar{u}_{TC,dis}$ ) can be calculated based on Equation (5) using the average  $\bar{m}_{dis} = \bar{m}_{comb} = 0.58$ kg/s based on Figure 1 a. The condition for ensuring a fully discharged PBTS is that  $D(=H)=\bar{u}_{TC,dis} t_{dis}$  and combining this condition

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with Equation (5) it is calculated that: D=H=2.43m. For this storage tank, the TC during the discharging can cross the PBTS height (H). The charging duration ( $t_{ch}$ ) is kept as the free variable.

In Figure 9 the *PBP*, *WH*<sub>stored</sub>, *WHR*<sub>cycle</sub>, *LCOH* and *CAPEX* are presented as a function of  $t_{ch}$  [150-250min]. The optimum solution is for  $t_{ch}$ =212.5min, resulting in max:*WHR*<sub>cycle</sub>=15.046GJ, min:*PBP*=0.17years (for  $C_{V,1}$ ), 0.32years (for  $C_{V,2}$ ) and min:*LCOH*=0.003€/kWh<sub>th</sub> (for  $C_{V,1}$ ), 0.006€/kWh<sub>th</sub> (for  $C_{V,2}$ ). In Figure 10 a. the temperature profiles at the end of the charging, stand-by and discharging operations for the PBTS with D=*H*=2.43m and  $t_{ch}$ =212.5min are presented. At the end of the discharge phase (Figure 9 a.-orange lines) the TC center has almost crossed the entire PBTS height H. In Figure 10 b. the preheated temperature during the SK HP is presented, where  $\overline{T}_{preh}$ =723°C.



**Figure 9. a.** For D=H=2.43m: PBP (blue lines) ( $C_{V,2}=1580$ €/m<sup>3</sup>: solid line and  $C_{V,I}=745$ €/m<sup>3</sup>: dashed line) and heat amounts (red lines) as function of  $t_{ch}$  and **b.** *LCOH* (blue lines) and *CAPEX* (red lines) ( $C_{V,2}=1580$ €/m<sup>3</sup>: solid lines and  $C_{V,I}=745$ €/m<sup>3</sup>: dashed lines) as a function of  $t_{ch}$ .



Figure 10. For the PBTS of D=H=2.43m and  $t_{ch}=212.5min$ : a. Temperature profiles inside the PBTS at the end of each operation phase. b. Preheated combustion air temperature ( $T_{preheated}$ ) during PBTS discharge phase (SK HP).

# 3.3 Scenario 3: Implementation of both a PBTS System and a HEX

In Sc.3 both a cylindrical PBTS and a HEX are implemented. Regarding the PBTS system, the optimum storage tank of Sc. 2 is considered, meaning the storage tank with D=H=2.43m for  $t_{ch}=212.5$ min. Again, both a crossflow and counterflow HEX are investigated using UA as their design variable [1-20kW/K]. For the PBTS only the cost of  $C_{V,2}=1580$ €/m<sup>3</sup> is considered in Scenario 3. The total values of  $WHR_{cycle}$ , *PBP*, *LCOH* and *CAPEX* are calculated including both the PBTS and the HEX values. The contribution of the PBTS to the total  $WHR_{cycle}$  and CAPEX is 15.046GJ and 19.692 k€ respectively. An additional operational constraint for Sc.3 is that the HEX operates when the preheated  $\dot{m}_{comb}$  temperature  $T_{preh}$  is lower than  $T_{exh}$ , so that  $\dot{m}_{comb}$  can be preheated. KPIs  $WHR_{cycle}$ , *PBP*, *LCOH* and *CAPEX* as a function of UA for both HEX designs are presented in Figure 11 a. and b. To select an optimum *UA* value for the crossflow and counterflow HEX, the multi-objective optimization criterion of Equation (13) is used.

The resulting optimum for the crossflow HEX occurs at minimum F=0.70 for UA=2.86kW/K and corresponds with { $WHR_{cycle}$ , PBP, LCOH}={15.749GJ, 0.813years, 0.0148€/kWh<sub>th</sub>}. For the counterflow HEX, minimum F is 0.58, obtained for UA=2.89kW/K, and results in { $WHR_{cycle}$ , PBP, LCOH}={15.817GJ, 0.811years, 0.0148€/kWh<sub>th</sub>}. For the optimum UA values the  $\dot{m}_{comb}$  preheated temperatures ( $T_{preheated}$ ) as a function of SK HP time duration (535min) are presented in Figure 12 a. for the crossflow HEX and Figure 12 b. for the counterflow HEX.



**Figure 11. a.** *PBP* (blue lines) and *WHR*<sub>cycle</sub> (red lines) as a function of *UA* and **b.** *LCOH* (blue lines) and *CAPEX* (red line-common cost for both HEX types) as function of *UA*. (Crossflow HEX: solid lines, counterflow HEX: dashed lines)



**Figure 12.** Preheated  $\dot{m}_{comb}$  temperature diagram during the SK HP. Light blue line  $(T_{cl})$ : outlet from the PBTS, dark blue line  $(T_{c2}=T_{preheated})$ : final outlet from the HEX, dashed orange line: inlet temperature  $(T_{hl})$  of  $\dot{m}_{exh}$  in the HEX and the solid orange line: the outlet  $(T_{h2})$ . Diagram for optimum **a.** crossflow HEX (*UA*=2.86kW/K) and **b.** counterflow (*UA*=2.89kW/K).

### 4 CONCLUSIONS

This study investigates the potential usage of a PBTS system for WHR in a shuttle kiln as widely used in the ceramic industry. Three distinct system configurations for WHR, namely a standalone HEX (Scenario 1), crossflow or counterflow, a single PBTS storage tank (Scenario 2) and their combination (Scenario 3), are compared in terms of  $WHR_{cycle}$ , PBP, and LCOH. To this end a 2P-1D numerical model, properly adjusted and validated based on experimental measurements for air as HTF, is used. The study's key findings are the following. For Scenario 1 the optimum UA for the crossflow (UA=3.49kW/K) and counterflow (UA=3.72kW/K) HEXs leads to  $WHR_{cycle}$  of 8.17GJ and 8.54GJ, PBPof 1.0 and 0.93yrs and LCOH of  $0.0182E/kWh_{th}$  and  $0.0170E/kWh_{th}$  respectively, with the counterflow outperforming the crossflow HEX at all the three KPIs used. Scenario 2 emphasizes on the proper sizing of the PBTS to preheat the total mass of the combustion air ( $m_{comb,tol}=19.7$ tn), through using the average thermocline velocity during the PBTS discharge phase ( $\bar{u}_{TC,dis}$ ). Optimal PBTS performance for a storage tank that can be full discharged by  $\dot{m}_{comb}$  was noted at D=H=2.43m and  $t_{ch}=212.5min$ , maximizing  $WHR_{cycles}=15.046$ GJ, minimizing PBP=0.17-0.32yrs and  $LCOH=0.003-0.006E/kWh_{th}$ 

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the max  $WHR_{cycle}$  from all 3 scenarios, with 5% enhancement, but significant *PBP* and *LCOH* increases (154%) compared to Scenario 2 alone. Overall, of the 3 scenarios, the HEXs (Scenario 1) appear to have the worst performance, leading to the minimum  $WHR_{cycle}$  and maximum *PBP*, *LCOH*. Future research can include exploring variable PBTS aspect ratios (a), PEs diameter (d<sub>p</sub>) and shape, utilizing multi-objective optimization for max  $WHR_{cycle}$  and min *PBP*, *LCOH*, as well as investigating WHR in other industries and either for on-site or off-site heat usage.

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