Modeling a Non-ferrous Melting Furnace

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Abstract: A one-dimensional model employed for the purpose of simulating the thermal characteristics of a non-ferrous melting furnace is presented. The thermal characteristics affect the heat transfer and hence the processing of materials within the furnace. The model employed consists of a governing equation for the combustion chamber and a transient heat conduction equation for the walls and roof. Each furnace component is treated as a one-dimensional conduction medium and the governing equations are solved numerically. The model was tested for the case of melting 3 kg of Aluminum charge in a Furnace designed at the Ambrose Alli University Epoma-Nigeria for melting non-ferrous metals. From the output of the program the maximum temperatures obtained at the end of melting for walls and roof, in-furnace average gas temperature (avt) and stack or exhaust gas were compiled and compared with experimental values. Higher temperature values were obtained when compared with experimental values. The wall temperature simulations are observed to be closely coupled an indication of a uniform temperature distribution within the furnace which was also reflected in the heat flux for the surfaces. The simulations are adjudged comparable with experimental data and the model is thus capable of predicting the key independent variables.

Key words: Furnace modeling, one-dimensional, simulations, temperature

INTRODUCTION

A mathematical model is a set of equations, algebraic or differential, which may be used to represent and predict certain phenomena (Szekely, 1988). Such models may be built from basic physical (including mechanical) and chemical laws or by drawing on the analogy of previous work through consideration of some basic physical situations.

Mathematical models may also be built from experiments with scale down physical models. The use of mathematical models thus makes possible the simulation and modification of system behaviour. System optimization and control are other important advantages of using mathematical models. The different model types vary both in their degree of complexity and the information obtained upon their application. According to Khalil (1982) and Baurkal et al. (2001), three different types exist. These are zero-dimensional, one-dimensional, and 2- and 3-dimensional models.

In Zero-dimensional modeling an overall heat and material balance of the system is done. This type of model does not give any spatial resolution but still gives a reasonable approximation of the overall performance of the system.

One-dimensional modeling considers only one spatial dimension and this greatly simplifies the number of equations, these models may still be fairly complicated and provide many details into the spatial changes of a given parameter.

2- and 3-Dimensional models models allow the determination of the spatial distribution of fluid and heat flow and other properties. Appropriate approximations of the real situation, leads to the solution of simultaneous partial differential equations, which represent the conservation of mass, momentum, energy and species.

A melting Furnace for non-ferrous metals have been previously developed and tested (Ighodalo and Ajuwa, 2010), the aim of the present work is to present the one-dimensional model employed in simulating the furnace thermal characteristics and the results obtained.

MATERIALS AND METHODS

Description of the furnace: This research work was carried out at the Mechanical Engineering department of Ambrose Alli University Ekpoma-Nigeria. The gas-fired melting furnace which was designed and constructed in the same department has been described by Ighodalo and Ajuwa (2010). The furnace walls were made from refractory castable material and are 110mm thick. A 10 mm thick Fibre-glass insulation was sandwiched between the outer face of the refractory walls and the encasing steel which is 1.5 mm thick. The whole frame work is a cubicule box, 700 x 600 x 600 mm with provisions for a burner hole, a chimney hole, a spout opening and charging door.

The burner is fired with butane gas and is a high velocity burner with a heat release rate of 160.3 kW. It is 480 mm in length and is connected with an air/fuel regulator (Ighodalo and Ajuwa, 2006).

The furnace chamber is mounted on a tilting mechanism. The isometric view for the complete assembly of furnace components is as shown in Fig. 1.
**Furnace modeling:** The model adopted is a one-dimensional model which consists of a governing equation for the combustion chamber and a transient heat conduction equation for the walls and roof. Each furnace component is treated as a one-dimensional conduction medium and the governing equations are solved numerically to yield the gas, walls and roof temperature profiles which can be compared with experimental data. A similar one-dimensional model has been used by Bui and Perron (1988) and Davies et al. (2000). The components of the furnace for the purpose of modeling are as represented in Fig. 2.

**Equations governing the combustion chamber:** The furnace chamber is governed by an ordinary differential equation describing the conservation of energy of the gas body (Bui and Perron, 1988). It states that the rate of energy accumulation in the gas equals the heat brought into the gas by the combustion of the fuel, minus the sum of the heat transferred to the furnace structure (walls and roof), through chimney and into the melt:

\[
V_c \cdot C_v \frac{dT}{dt} = Q_1 - \{Q_{11} + Q_{10} + Q_{12}\} \quad (1)
\]

This equation is simply evaluated over time using a backward difference approximation. \(Q_1\) is given by the combustion of natural gas:

\[
Q_1 = \sum_{\rho} n_1 \int_{T_298}^{T_\infty} C_p(T) dT + \Delta H_{298}
\]
where P and R are the products and reactants. The combustion is assumed to be complete and stoichiometric. The constant pressure specific heat terms \( c_p(T) \) are approximated using polynomials of temperature (T) of the form:

\[
C_p(T) = a + b \cdot 10^{-2} T + c \cdot 10^{-4} T^2 + d \cdot 10^{-6} T^3
\]

where a, b, c and d are constants for each gas in the combustion product.

\( Q_{71}, Q_{70} \) and \( Q_{72} \) are the heat flow through the walls, metal and extraction system respectively as shown in Fig. 2. These are calculated using radiation and convection models.

**The radiation and convection models:** Radiosity approach is employed in modeling thermal radiation in the furnace chamber. The radiosity expression for each surface in an enclosure with combustion gas is as given by Davies *et al.* (2000):

\[
J_i = \varepsilon_i E_{bi} + (\sum \Sigma F_{ij} \tau_j J_j) + \varepsilon_i E_{bs}
\]

where \( \varepsilon = 1 - \varepsilon_i \), \( E_{bi} \) is emissivity of gas medium, \( \tau_j \) is transmissivity of gas evaluated at the temperature of the \( j^{th} \)wall:

\[
Q_i = \varepsilon A_i / (1 - \varepsilon) (E_{bs} - J) + h A_i (T_g - T_s)
\]

The emissivity and absorptivity of the combustion gases used in Eq. (21) are obtained from the mixed grey gas model (Tucker, 2003),

\[
e_g(T_g) = \sum_{n} a_n(T_g) [1 - \exp(-K_{g,n}P L_m)]
\]

For \( n = 1, N \) (6)

\[
a_g(T_g) = \sum_{n} a_n(T_s) [1 - \exp(-K_{g,n}P L_m)]
\]

For \( n = 1, N \) (7)

\[
\sum_{n} a_n(T_g) = 1 \quad \text{and} \quad \sum_{n} a_n(T_s) = 1
\]

where \( a_n \) are the weighting coefficients, \( T_g \) and \( T_s \) are gas and wall surface temperatures respectively, \( K_{g,n} \) is absorption coefficient. \( L_m \) is the mean beam length, which for a particular geometry can be approximated from (Holman, 1992):

\[
L_m = 3.6 * V / A
\]

where V is the total volume of the gas and A is the total surface area.

The weighting coefficients are simply represented by (Tucker, 2003):

\[
a_g(T_g) = b_{1,n} + b_{2,n} T_g
\]

\[
a_g(T_s) = b_{1,n} + b_{2,n} T_s
\]

where, \( b_{1,n} \) and \( b_{2,n} \) are constants.

The radiation shape factors \( F_{ij} \) that are also needed in Eq. (4), were obtained from the expressions for two rectangles which are either parallel or at right angles were used. For parallel rectangles with equal sides of lengths a and b spaced a distance c apart, the edge distance ratios are \( X = a/c \) and \( Y = b/c \). The view factor is given by (Siegel and Howell, 1972):

\[
F_{ij} = \frac{2}{\pi X Y} \left[ \ln \left( \frac{1 + X^2}{1 + Y^2} \right) \right]^{\frac{1}{2}}
\]

\[
+ X \sqrt{1 + Y^2} \tan^{-1} \frac{X}{\sqrt{1 + Y^2}}
\]

\[
+ Y \sqrt{1 + X^2} \tan^{-1} \frac{Y}{\sqrt{1 + X^2}}
\]

\[
- X \tan^{-1} X - Y \tan^{-1} Y
\]

For two rectangles \( hl \) and \( lw \) at right angles with a common edge l, the length ratios are \( H = h/l \) and \( W = w/l \). The view factor as given by Siegel and Howell (1972) is:

\[
F_{ij} = \frac{1}{\pi W} \left[ W \tan^{-1} \frac{1}{W} + H \tan^{-1} \frac{1}{H} \right]
\]

\[
- \sqrt{H^2 + W^2} \tan^{-1} \frac{1}{\sqrt{H^2 + W^2}}
\]

\[
+ \frac{1}{4} \ln \left( \frac{1 + W^2}{1 + H^2} \right)
\]

\[
\]
The body, i.e., the wall is divided into a number of elements, the finite difference approximation is obtained for the governing equation and the appropriate form for internal and boundary nodes are established so that the temperature-time history can be obtained.

The explicit form of the discretised equation for both internal and boundary nodes respectively are:

\[ T_{i,j+1} = 1/R \left( T_{i+1,j} + T_{i-1,j} \right) - T_{i,j} \left( 1/R - 1 \right) \quad (19) \]

where, \( T_{i,j} \) is the temperature of an unknown mesh point (at time period \( j + 1 \)), and it is now expressed in terms of the known temperatures on the previously calculated time period \( j \).

\[ Q_{o.2} \Delta t / (\Delta x)^2 \cdot \alpha \Delta t / (\Delta x)^2 \cdot 2(T_{i}^{n+1} - T_{i}^{n}) = T_{i}^{n+1} - T_{i}^{n} \quad (20) \]

for convergence or stability of the numerical solution \( \Delta t \) and \( \Delta x \) should be chosen in such a manner that:

\[ R = (\Delta x)^2 / \alpha . \Delta t \geq 2 \]

**Boundary and initial conditions:** The initial condition for Eq. (1) is obtained by assuming that at the start of the simulation the gas body has previously reached a steady state and therefore the the accumulation term is nil:

\[ V_{x}C_{v} \frac{dT_{x}}{dt} = 0 \]

The boundary conditions for the walls and charge are of the known boundary heat flux type, as the radiation and convection heat fluxes are calculated separately.

At the start of simulation all surfaces are assumed to be at ambient temperature.

The initial condition is:

\[ T_{o} = 30^\circ C \text{ at } t = 0 \]

The heat received by radiation and convection is conducted into the walls and charge so that the boundary condition is:

\[-k \frac{\partial T}{\partial x} = Q_{rad} + Q_{conv} \quad (21)\]

**Model algorithm:** A computer program written in MATLAB was developed based on the models presented. The flow chart for the simulation is shown in Fig. 3. For the given conditions of furnace geometry, fuel properties, furnace gas composition, charge and wall refractory properties, the adiabatic temperature and heat of gas combustion is estimated. The wall surface and charge surface temperature is initialized and the gas emissivity and absorptivity are calculated, shape factors for the surfaces are determined which leads to the determination.
Table 1: Experimental and Simulated temperatures for walls, roof, in-furnace, exhaust gas, and metal respectively, in °C

<table>
<thead>
<tr>
<th></th>
<th>wall1</th>
<th>wall2</th>
<th>wall3</th>
<th>wall4</th>
<th>Roof</th>
<th>Avtg</th>
<th>Tstac</th>
<th>metal</th>
</tr>
</thead>
<tbody>
<tr>
<td>Experiment</td>
<td>770</td>
<td>745</td>
<td>735</td>
<td>700</td>
<td>780</td>
<td>1025</td>
<td>593.8</td>
<td>660</td>
</tr>
<tr>
<td>Simulated</td>
<td>986.5</td>
<td>987.6</td>
<td>985.4</td>
<td>965.9</td>
<td>986.5</td>
<td>1692.4</td>
<td>833</td>
<td>756.9</td>
</tr>
</tbody>
</table>

RESULTS AND DISCUSSION

The model was tested for the case of furnace operation melting 3 Kg of Aluminium charge. From the output of the program the maximum temperatures obtained at the end of melting for walls and roof, in-furnace average gas temperature (avtg) and stack or exhaust gas were compiled and are compared with experimental values as shown in Table 1. The simulated
From the simulations conducted, higher values of wall temperatures were obtained when compared with experimental values as shown in Table 1. The average difference is 236.4°C for the inner wall surface temperatures.

The wall temperature simulations are observed to be closely coupled as shown in Fig. 4 and 5, this is a confirmation of a uniform temperature distribution within the furnace. The simulated and experimental temperature curves follow a very similar profile in all cases increasing linearly as can be seen from the Figures with the roof having the maximum temperature from Table 1. Nodal temperatures across the walls thicknesses also display an almost linear decreasing distribution as can be seen from Fig. 7. Higher values for the in-furnace gas, exhaust gas and metal surface temperatures were also produced by the simulation. The discrepancies between experimental and simulated values can be attributed to air in-leakage into the furnace, experimental errors and simplifying assumptions employed in the simulations.

The results obtained from the simulations can be said to be in fairly good agreement with the experimental values, thus validating the model employed.

From Table 2, the heat fluxes to the walls are also observed to be uniform, all within 21 kW/m², this is not surprising since a uniform temperature obtains within the furnace. The heat flux on the metal charge is lower than for the walls and is adequate for melting of the charge. The thermal efficiency calculated by the model is 47.52%.

**CONCLUSION**

The one-dimensional model employed in modeling the Furnace, have produced results which are comparable with experimental data.

From the foregoing the model is capable of predicting the key dependent variables, the average inside wall surface temperature, the net heat transfer rate to the exposed wall surfaces, the net heat transfer rates to the metal, the average in-furnace gas temperature, and the metal surface temperature.

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**REFERENCES**