

# Simulation & Validation of Performance & Emission Parameters of Jatropha Blends in Diesel with Exhaust Gas Recirculation

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**Abstract:** In the present study, a simulation code is written for diesel engine with various combinations of jatropha blends and with exhaust gas recirculation (EGR), based on steady-state phenomenological model. The emission model was developed by using chemical equilibrium reactions and calculation of thermodynamics properties based on their mixture composition. The model has been developed to predict the inlet and exhaust manifold temperature, pressure, engine torque and exhaust species variation at different percentage of jatropha blends and EGR. The system has been validated with experimental data.

**Keywords:** Exhaust gas recirculation, Jatropha blends, Simulation, Equilibrium combustion

## 1. Notation

$p_{im}$  = Inlet Manifold Pressure

$p_{em}$  = Exhaust Manifold Pressure

$T_{im}$  = Inlet Manifold Temperature

$T_{em}$  = Exhaust Manifold Temperature

$V_{im}$  = Inlet manifold volume

$W_c$  = Mass flow rate at compressor

$W_{egr}$  = Mass flow rate of re-circulated exhaust gas

$W_{ei}$  = Mass flow rate of at engine inlet

$W_{eo}$  = Mass flow rate of at engine outlet

$W_t$  = Mass flow rate at turbine

$R_a$  = ideal gas constant for air

$W_f$  = Mass flow rate of fuel

$\eta_s$  = compensation factor for non ideal cycles

$q_{in}$  = Specific energy content of the charge per unit mass

$x_r$  = Residual gas fraction

$x_{cv}$  = Ratio of fuel consumed during constant volume combustion

$\eta_{vol}$  = Volumetric efficiency

$N$  = Number of cylinders

$V_D$  = Displacement volume

## 2. Introduction

The mathematical modelling is one of the important tools to simulate a diesel Engine using Jatropha bio-fuel to predict: (i) engine performance; (ii) engine emissions (oxides of nitrogen and carbon) at various percentage of EGR. The combustion process is very complex in a diesel engine [1]. Combustion models can be classified as thermodynamic model [2] and detailed model [3]. Thermodynamics and phenomenological models give reasonable prediction of intake and exhaust pressure and temperature, so a phenomenological steady state model has been used and numerical model is validated against experimental data.

## 3. Mathematical Modelling

To develop a simple model, that captures the dominating effects in the mass flows, the following assumptions were made:

- The manifolds are modelled as standard isothermal models.
- All gases are considered to be ideal and there are two sets of thermodynamic properties:
  - a) Air has the gas constant  $R_a$  and the specific heat capacity ratio  $C_a$ .

- b) Exhaust gas has the gas constant  $R_e$  and the specific heat capacity ratio  $C_e$ .
- The EGR gas in the intake manifold affects neither the gas constant nor the specific heat capacity in the intake manifold.
  - No heat transfer takes place in the intake manifold.
  - No backflow can occur in the EGR valve, compressor, turbine or the cylinder.
  - The oxygen fuel ratio is always larger than 1.
  - The mixture of air and fuel is treated as perfect gas.
  - The mixture inside the cylinder is homogeneous.
  - Equivalence ratio is assumed to be 0.5 for analysis with EGR.
  - Effect of radiation is neglected.
  - No chemical change takes place in the mixture of air and fuel before combustion.
  - All properties of gases inside the cylinder is only time dependent.
  - The percentage of re-circulated exhaust gases is ranging from 0 to 35.

The standard isothermal model that is based upon mass conservation and the ideal gas law, gives the differential equations for the manifold pressures as [4]:

$$\frac{dp_{im}}{dt} = \frac{R_a T_{im}}{V_{im}} (W_c + W_{egr} - W_{ei}) \quad (1)$$

$$\frac{dp_{em}}{dt} = \frac{R_e T_{em}}{V_{em}} (W_{eo} - W_{egr} - W_t) \quad (2)$$

The total mass flow from the intake manifold into the cylinders is modelled using the volumetric efficiency

$$W_{ei} = \frac{\eta_{vol} p_{im} \eta_e V_d}{120 R_a T_{im}} \quad (3)$$

The volumetric efficiency is in its turn modelled as:

$$\eta_{vol} = C_{vol1} \sqrt{p_{im}} + C_{vol2} \sqrt{N} + C_{vol3} \quad (4)$$

Where  $C_{vol1}$ ,  $C_{vol2}$  and  $C_{vol3}$  are constants for volumetric efficiency determination.

The fuel mass flow  $W_f$  into the cylinders is controlled by  $U_\delta$  which gives the injected mass of fuel in mg per cycle and cylinder:

$$W_f = \frac{10^{-6} U_\delta N \eta_{cyl}}{120} \quad (5)$$

The mass flow  $W_{eo}$  out from the cylinder is given by the mass balance as

$$W_{eo} = W_f + W_{ei} \quad (6)$$

The cylinder out temperature  $T_{em}$  modelled based upon ideal gas Seliger cycle

$$T_{em} = \eta_s \left( \frac{p_{em}}{p_{im}} \right)^{1-(1/\gamma)} \left( \frac{V_2}{V_1} \right)^{\gamma-1} \left( 1 + \frac{q_{in}}{C_v T_1} \left( \frac{V_2}{V_1} \right)^{\gamma-1} x_{cv} \right)^{1/\gamma-1} * \left( q_{in} \left( \frac{1-x_{cv}}{C_p} + \frac{x_{cv}}{C_v} \right) + T_1 \left( \frac{V_1}{V_2} \right)^{\gamma-1} \right) \quad (7)$$

The temperature at inlet valve closing after intake stroke and mixing

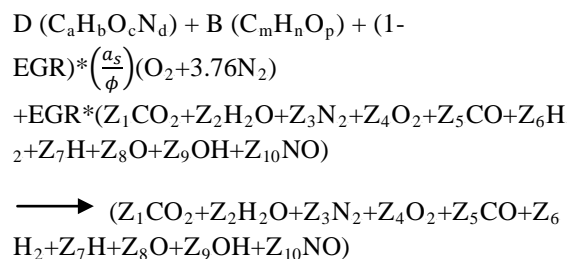
$$T_1 = x_r T_{em} + (1 - x_r) T_{im} \quad (8)$$

Where the residual gas fraction is modelled as

$$x_r = \left( \frac{p_{em}}{p_{im}} \right)^{(1/\gamma)} \left( 1 + \frac{q_{in}}{C_v T_1} \left( \frac{V_2}{V_1} \right)^{\gamma-1} x_{cv} \right)^{-1/\gamma} \quad (9)$$

#### 4. Chemical Equilibrium Combustion and Mole Fraction of Species

When the blend of diesel fuel and Jatropha bio fuel is combusted in presence of air with the percentage of re-circulated exhaust gases, it will replace the amount of oxygen and nitrogen from the fresh air entering the cylinder with the carbon dioxide and water vapour. The considerable 10 species, which are considered, are H, O, N, H<sub>2</sub>, OH, CO, NO, CO<sub>2</sub>, H<sub>2</sub>O and N<sub>2</sub> [5]. The chemical reaction of combustion of diesel and bio-fuel blend, with EGR, is as follows [6]:



#### Carbon balancing:

$$Bm + Da = (1 - \text{EGR}) * (x_1 + x_5) * N_t$$

**Hydrogen balancing**

$$Bn + Db = (1-EGR) * (2x_2 + 2x_6 + 2x_7 + x_9) * N_t$$

**Oxygen balancing**

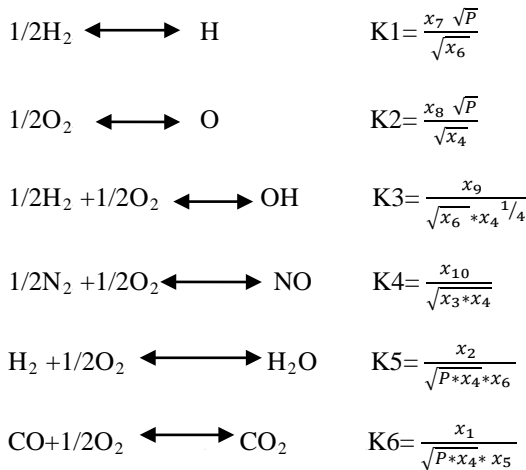
$$Bp + Dc + (1-EGR) * 2 \left( \frac{a_s}{\phi} \right) = (1-EGR) * (2x_1 + x_2 + 2x_4 + x_5 + x_8 + x_9 + x_{10}) * N_t$$

**Nitrogen balancing**

$$Dd + (1-EGR) * 2 \left( \frac{a_s}{\phi} \right) * 3.76 = (1-EGR) * (2x_3 + x_{10}) * N_t$$

Where  $x_i = \frac{Z_i}{Z_t}$  ;  $Z_t = \sum Z_i$  ;  $\sum x_i = 1$

**Dissociation reaction**



Where,  $K_i = \exp(-\frac{\Delta G}{RT})$

$K_i$ = equilibrium constant for species  $i$  [14].

$\Delta G$ =Gibbs function at specified temperature

$R$ = universal gas constant

$T$ = combustion temperature

**5. Simulation Code**

A computer code has been written to solve the equations representing the models of gas species, reaction chemistry, and thermodynamic properties of mixture in the cycle. Effective iteration methods have been used for simulation [12]. The time step used in the calculation, especially when combustion takes place, was very small.

**6. Results, Validation and Conclusions**

The results were presented for full cycle for four strokes DI diesel engine. The runs were made using specification of Kirloskar diesel engine ( $B = 110$  mm,  $S = 110$  mm, speed = 1500 rpm) [7]. The simulation results have been validated with experimental data.

For intake manifold pressure, the deviation is about 0.48-3.58 %. The deviation in exhaust manifold pressure and engine torque is about 0.25-4.31 % and 0.95-5.86 % respectively [8].

**6.1 Intake Manifold Pressure**

The fig.1 shows the comparison between the simulated data and the experimental data. This variation of intake manifold pressure is for the engine running at 1500 rpm, the fuel injection rate is 114.4 mg/cycle, EGR rate 50 % [13].

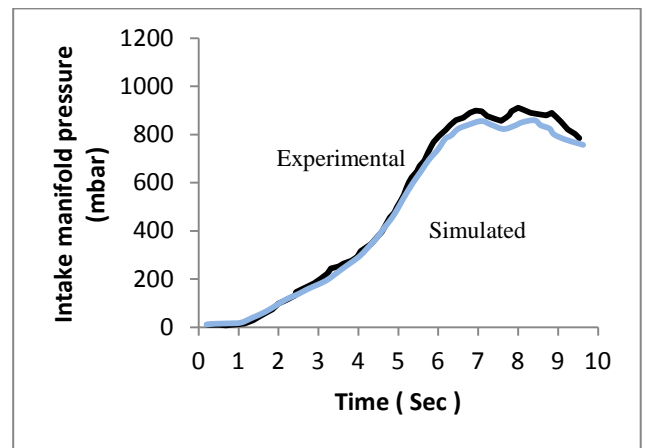


Fig.1 Variation of intake manifold pressure with time

**6.2 Exhaust Manifold Pressure**

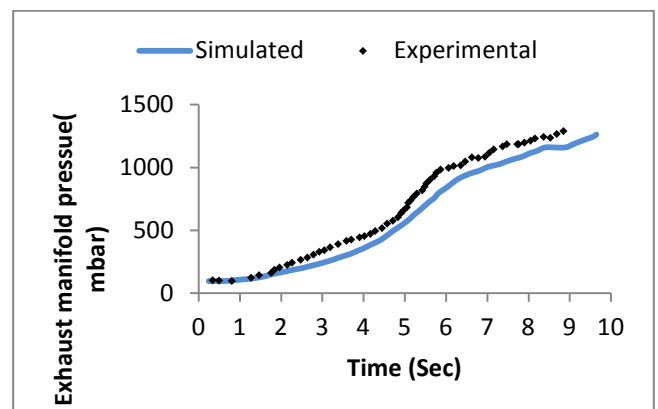


Fig. 2 Variation of exhaust manifold pressure with time

**6.3 Engine Torque**

The variation of torque produced by the engine with time and the comparison of simulated data and experimental data is shown in fig. The maximum engine torque produced by the diesel engine in the present study is about 230 N-m as shown in fig 3.

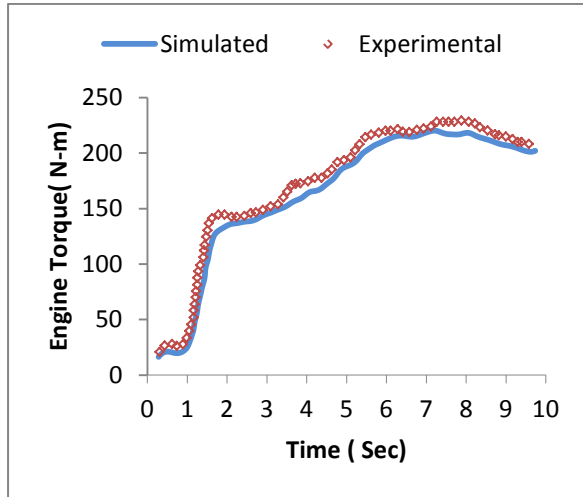


Fig. 3 Variation of engine torque with time

#### 6.4 Nitrogen Oxide Formation

The reactions forming NO are very temperature dependent, as the rate of dissociation of nitrogen is directly proportional to the temperature increase.

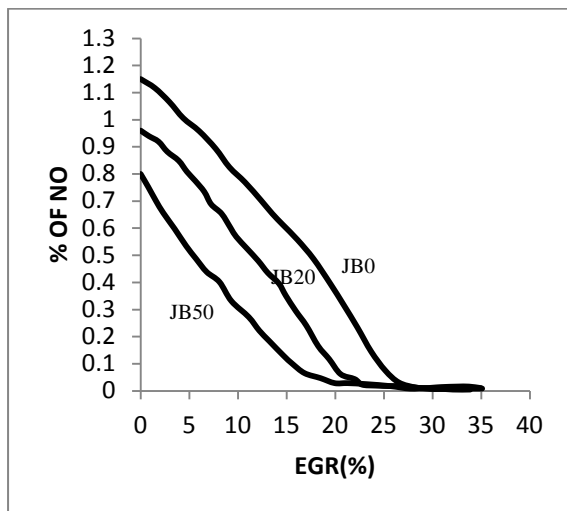


Fig. 4 Variation of Nitrogen oxide with EGR %

JBD (Jatropha biodiesel) blends emitted NOx was slightly lower than that of DF, at 0% EGR. At 0% EGR, the NOx emission of JB20 and JB50 was lower than that of diesel fuel. The NOx emission of all fuels decreased linearly, when EGR was operated [9]. As the percentage of EGR increases,

the amount of NO decreased rapidly for different blends.

#### 6.5 Carbon Monoxide Formation

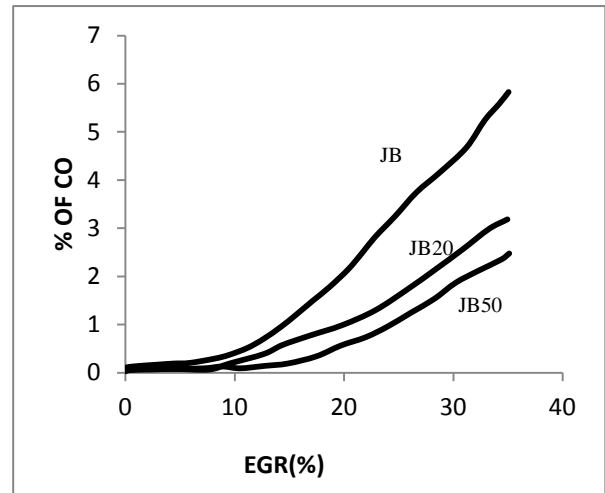


Fig. 5 Variation of Carbon monoxide with EGR %

The CO emission increased with increasing EGR rates. The amount of CO in the engine exhaust increases slowly till 10% EGR and after that the amount of CO is increased very rapidly. The amount of CO is lower for 50% Jatropha blend than that for 0% and 20% Jatropha blends [10].

#### 6.6 Carbon Dioxide Formation

The CO<sub>2</sub> emission increased with increasing biodiesel amount in the blends. The CO<sub>2</sub> emission of DF increased slightly, when EGR was operated. While, the CO<sub>2</sub> emission of JBD blends increased rapidly; especially at over 20% EGR. Beyond 20% EGR, the CO<sub>2</sub> emission from JBD blends was higher than that of DF [11].

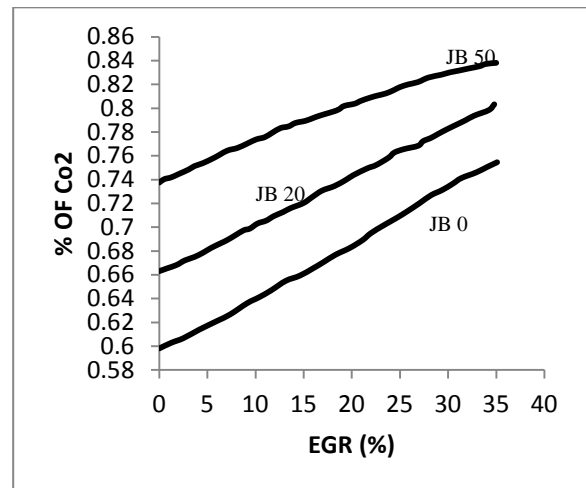


Fig. 6 Variation of Carbon dioxide with EGR %

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## Biographies

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