Prediction and Optimization of Fish Biodiesel Characteristics Using Permittivity Properties

M. Zarein¹, M. H. Khoshtaghaza¹*, B. Ghobadian¹, and H. Ameri Mahabadi²

ABSTRACT

The purpose of this research was to predict and optimize the fish biodiesel characteristics using its permittivity properties. The parameters of biodiesel permittivity properties such as ε, dielectric constant, and ε'', loss factor at microwave frequencies of 434, 915, and 2,450 MHz, were used as input variables. The fish biodiesel characteristics, as Fatty Acid Methyl Ester (FAME) content and flash point at three different levels of reaction time 3, 9, and 27 min and catalyst concentrations 1, 1.5, and 2% w/w oil, were selected as output parameters for the models. Linear Regression (LR), the Multi-Layer Perceptron (MLP), and the Radial Basis Function (RBF) as the methods of Artificial Neural Networks (ANN), and the response surface methodology were compared for prediction and optimization of FAME content and flash point. A comparison of the results showed that the RBF recorded higher coefficient of determination at frequency of 2,450 MHz as 0.999 and 0.988 and lower root mean square error as 0.009 and 0.023 for FAME content and flash point, respectively. The optimum condition was obtained using RSM by FAME content of 89.88% and flash point of 152.7°C with desirability of 0.998.

Keywords: Artificial neural network, Fatty acid methyl ester content, Flash point, Optimum condition, Response surface methodology.

INTRODUCTION

Waste Fish Oil (WFO) is produced in large quantity by fish-processing industry as a by-product in Iran. This by-product has similar calorific value to petroleum distillates and is a renewable energy source (Yahyae et al., 2013). Biodiesel is defined as a fuel that contains mono-alkyl esters of long chain fatty acids (Ghobadian et al., 2008). One of the advantages of using biodiesel fuel is the significant reduction in emissions such as CO and HC. Therefore, biodiesel produced from oilseed crops can be used as fuel that is required by diesel engines (Ghazali et al., 2015; Abedin et al., 2014).

Biodiesel has been produced globally from several products such as waste fish oil (García-Moreno et al., 2014), palm oil (Tan et al., 2011), sunflower oil (Yin et al., 2012), vegetable oil (Kouzu and Hidaka, 2012), waste cooking oil (Talebian-Kiaalaieh et al., 2013) and rapeseed oil (Duren et al., 2015). Since Free Fatty Acid (FFA) content is a critical parameter in the conversion of fish oils to methyl esters, the performance of a Fourier Transform InfraRed (FTIR) spectroscopic method was assessed as an alternative to the conventional AOCS titrimetric method (Alberta et al., 2009). The properties of the triglyceride and the biodiesel fuel are determined by the amounts of each fatty acid present in their molecules. Transesterification does not alter the fatty acid composition of the feedstock and this composition plays an important role in some critical parameters of the biodiesel, as viscosity and flash point properties (Mejia et al., 2013; Ramos et al., 2009). Incorporating ultrasonic energy into traditional transesterification reactions can emulsify the reactants to reduce the requirement of catalyst amount, methanol-to-
oil ratio, reaction time and reaction temperature (Parkar et al., 2012; Santos et al., 2009). In biodiesel production using ultrasonic, the maximum reaction time to reach the maximum conversion rate for castor, palm, and fish oil were obtained as 9, 63, and 27 minutes, respectively (Maghami et al., 2015).

Dielectric spectroscopy has been successfully used for production and characterization of biodiesel (Sorichetti and Romano, 2005). It is used to characterize feedstocks from different origins (Corach et al., 2014), to detect alcohol in the light phase after the completion of the transesterification process (Ramos et al., 2009), during the purification process and in the final product (Gonzalez et al., 2008), and also for the characterization of FAME (Corach et al., 2012). Based on geographic limitations and oil prices, biodiesel can be produced from a variety of feedstocks and different technologies can be applied for biodiesel production (Oliveira et al., 2008; Balabin et al., 2010). Consequently, the final product can have different properties, so, the quality control of biodiesel is very important. The EN 14214 mandates 25 parameters that have to be analyzed to certify biodiesel quality and these analyses are expensive and time consuming (Flores et al., 2012; Silva et al., 2017).

Measurements of dielectric properties at microwave frequencies offer several advantages for characterization (Romano and Sorichetti, 2011). Particularly in the microwave range, the technique is fast, accurate, simple, cheap, and non-destructive (Liptak, 2003).

Artificial Neural Network (ANN) and Response Surface Methodology (RSM) have been widely applied for modeling and optimization of biodiesel synthesis from different feedstocks using different methods such as classical base catalyzed transesterification (Stamenkovic et al., 2015; Rajkovic et al., 2013), heterogeneous base catalyzed transesterification (Betiku and Ajala, 2014), ultrasound assisted base catalyzed transesterification (Moorthi et al., 2015), infrared irradiation assisted esterification (Chakraborty and Sahu, 2014). Different ANN models based on multi-layer feed forward, radial base, generalized regression, and recurrent network were developed for predicting the Cetane Number (CN) of biodiesel fuel (Ramadhas et al., 2006). The biodiesel production from rapeseed soap stock and methanol in the presence of the candida rugosa lipase immobilized on chitosan was analyzed. Using ANN showed desirable correspondence between predicted and experimental values of the FAME yield (Ying et al., 2009).

The aim of this study was, therefore, to develop ANN models to predict the Fatty Acid Methyl Ester (FAME) content and flash point during biodiesel production using permittivity properties of biodiesel as the ANN inputs and optimize the conditions of these characteristics by RSM.

**MATERIALS AND METHODS**

Waste Fish Oil (WFO) was purchased from fishmeal plant in Semnan, Iran. WFO properties like acid value, density, kinematic viscosity, flash point, Free Fatty Acid (FFA) content, color and mean molecular weight were measured (Table 1). All chemicals, such as methanol, KOH, H₂SO₄ with purity of 99.99%, were analytical grade and purchased from Merck (Germany). The ultrasonic processor UP400S (400W, 24 kHz) with probe diameter of 40 mm and probe length of 100 mm was made by Hielscher (Germany).

**Experimental Procedure**

Esterification was applied to reduce FFA content. Methanol and waste fish oil were transmitted to two-layer 250 mL glass reactor which was equipped with temperature control (±1°C) and condenser. For the temperature control, external heating was applied with an accuracy of ±1°C. Experiments were carried out in the temperature range of 40-60°C. Sulfuric acid
(1% wt wt⁻¹) was used as catalyst and reaction continued for 1 h to reduce FFA content. Transesterification is a reversible reaction that requires more Methanol Ratio (MR) than the stoichiometric values to reach the higher product yields (Vicente et al., 2007; Leevijit et al., 2008). The optimum condition of biodiesel production is related to the type and amount of catalysts and oils (Zheng et al., 2006). For acid catalysis reactions, always higher ratios were suggested with longer reaction times. After FFA content reduction, the reactor content was removed for separating funnel and allowed to separate layers methanol-reach phase and esterified oil. Homogenous base catalysts are used widely for the biodiesel production. These catalysts have more conversion when compared to the acid, enzyme, or base heterogeneous catalysts. For this reason, potassium hydroxide as a homogeneous catalyst was used in this work for the biodiesel production from waste fish oil. Catalyst type and loading amount are the main parameters that affect the biodiesel production. Lower catalyst loading caused minor conversion and higher loading caused side reactions like saponification in transesterification process. The properties of the produced biodiesel in comparison with the ASTM D6751 standard are described in Table 2. FAME content of each sample was calculated according to equation (1):

\[
\text{FAME content (\%)} = \frac{\sum A_i - A_i}{A_i} \times \frac{m_i}{M} \times 100
\]

(1)

Where, \(i\) is standard sample and \(A_i\) and \(A\) are the peak Areas of standard and total sample in GC reports, respectively. \(M\) is the weight of sample and \(m\) is the weight of standard in the sample. Weight of product was estimated after separation of glycerin phase and removal of impurities by washing and drying of biodiesel phase in each experiment. The flash point test was carried out by FLPH (CCCFP) Continuously Close Cup (Grabner, Austria) according to ASTM D93. For measuring acid number and Free Fatty Acid (FFA) content of WFO, 1 g of waste oil was dissolved in 10 mL 2-propanol and then titrated by KOH 0.1M (5.61 mg KOH in 1 mL ethanol) in the presence of phenolphthalein as indicator (Maghami et al., 2015). Acid number is mg KOH which is used for titration of 1 g oil and FFA content was calculated by Equation (2).

\[
\text{FFA\%} = \frac{V.C.M}{10.m} \times \frac{m_i}{M}
\]

(2)

Where, \(V\) is KOH solution Volume (mL), \(C\) is solution molarity (0.1 mol L⁻¹), \(M\) is WFO Molecular weight (g mol⁻¹) and \(m\) is WFO weight (g).

### Table 1. Chemical and physical properties of the used WFO.

<table>
<thead>
<tr>
<th>Property</th>
<th>Units</th>
<th>Measured value</th>
</tr>
</thead>
<tbody>
<tr>
<td>Acid value</td>
<td>mg KOH g⁻¹ oil</td>
<td>9.98±0.02</td>
</tr>
<tr>
<td>Density (@15°C)</td>
<td>g cm⁻³</td>
<td>0.912±0.007</td>
</tr>
<tr>
<td>Kinematic viscosity (@25°C)</td>
<td>mm² s⁻¹</td>
<td>44.551±2.685</td>
</tr>
<tr>
<td>Flash point, Closed cup</td>
<td>°C</td>
<td>196±1</td>
</tr>
<tr>
<td>FFA content</td>
<td>%</td>
<td>5.47±0.01</td>
</tr>
<tr>
<td>Color</td>
<td></td>
<td>Dark brown</td>
</tr>
<tr>
<td>Mean molecular weight of WFO</td>
<td>g mol⁻¹</td>
<td>903</td>
</tr>
</tbody>
</table>
Figure 1. (a) Gas Chromatography (GC) analyzer and (b) The yields of WFO compounds according to the carbon chain by GC.

Permittivity Measurement

The permittivity properties of a substance may be described in the frequency domain by the complex permittivity, \( \varepsilon \). The real part (dielectric constant), \( \varepsilon' \), represents electrical polarization and the imaginary part (loss factor), \( \varepsilon'' \), is related to energy dissipation as the following equation (Corach et al., 2015):

\[
\varepsilon = \varepsilon' - j\varepsilon''
\]  

(3)

Permittivity properties of the mixtures were determined using an Agilent ENA series E5071C Network Analyzer with Agilent 85070E dielectric probe kit (Agilent Technologies, Inc. Santa Carla, CA) using coaxial line method at frequency of 434, 915 and 2,450 MHz according to ISM standard. Frequency values were selected based on the limits of measurement system, because of Agilent 85070E dielectric probe kit is suitable for frequencies above 200 MHz (Muley and Boldor, 2013). The network analyzer was controlled by Agilent 85070E dielectric kit software and calibrated using the 3-point method (short-circuit, air and water at 25°C).

Linear Regression Model

Regression analysis is a statistical technique for examining and modeling the association between variables. It is a widely used statistical technique that operates by building mathematical equations relating the response (variable) to set the predictors or independent variables. Linear Regression (LR) was used to model the values of a dependent variable based on a linear relationship with one or more predictors. A LR model assumes that there is a linear relationship (denoted by a straight line) between the dependent variables and the independent variables (Rodriguez et al.,...
The LR model is often formulated as Equation (4):
\[ y = b_0 + b_1 x_1 + \cdots + b_n x_n + e \]  
(4)
Where, \( y \) is the value of the dependent variable, \( x_n \) is the predictor variable, \( b_n \) is the coefficient value, and \( e \) is the observed error (uncontrolled factors and experimental error). The model parameters \( (b_i) \) were estimated using a regression model.

Artificial Neural Networks (ANN) Models

Neural networks may be used as a direct substitute for auto correlation, multivariable regression, linear regression, trigonometric and other statistical analysis and techniques (Singh et al., 2003). Multi Layer Perceptron (MLP) and Radial Basis Function (RBF) are two of the most widely used neural network architecture in literature for classification or regression problems (Cohen and Intrator, 2003). Both types of neural network structures are good in pattern classification problems. The output of a MLP is produced by linear combinations of the outputs of hidden layer nodes in which every neuron maps a weighted average of the inputs through a sigmoid function. In one hidden layer RBF network hidden nodes map distances between input vectors and center vectors to outputs through a nonlinear kernel or radial function (Balabin et al., 2011). In this study, the two different architectures of ANN (MLP and RBF) were also used to estimate the FAME content and flash point. All data were first normalized and divided into three data set such as training (70% of all data), test (15% of all data) and verification (15% of all data). The STATISTICA12 software was used in neural network analysis having a three-layer feed-forward network that consists of an input layer (2 neurons), one hidden layers (2 neurons for MLP, 15 neurons for RBF) and two output layer (Figure 2). Neuron numbers in hidden layers were selected from a series of trial runs of the networks having 1 neuron to 20 neurons in order to obtain the neuron number in the network having minimum error. Variable learning rate with momentum (trainLm) as networks training function and tangent sigmoid (tansig) was used as an activation (transfer) function for all layers (Yilmaz and Kaynar, 2011).

Response Surface Methodology (RSM)

Response Surface Methodology (RSM) has an important application in the design, development, and formulation of new products, as well as in the improvement of existing product design. It defines the effect of the independent variables, alone or in combination, on processes. In addition, to analyze the effects of the independent variables, this experimental methodology generates a mathematical model which describes the chemical or biochemical processes (Halim et al., 2009). In order to obtain the optimum value, Equation (5) is used:

![Figure 2. MLP and RBF neural network structure used in the study.](image)
\[ Y_i = \beta_0 + \sum \beta_i X_i + \sum \beta_{ij} X_i X_j + \sum \beta_{ij} X_i^2 + \varepsilon \tag{5} \]

Where, \( \beta_0, \beta_i, \beta_{ij}, \beta_{ij} \) are regression coefficients for intercept, linear, interaction, and quadratic coefficients, respectively, while \( X_i \) and \( X_j \) are coded independent variables and \( \varepsilon \) is the error. In the present study, Box-Behnken design with 3 central points was used.

The validation and performance of the LR, MLP, RBF, and RSM models were compared using \( R^2 \), MAPE, and RMSE as follows, equations (6-8) (Amid and Mesri Gundoshmian, 2016):

\[ R^2 = 1 - \left( \frac{\sum_{i=1}^{N}(z_i - \bar{z}_i)^2}{\sum_{i=1}^{N}(z_i - \bar{z})^2} \right) \tag{6} \]

\[ MAPE = \frac{1}{N} \sum_{i=1}^{N} \left| \frac{z_i - \bar{z}_i}{z_i} \right| \times 100 \tag{7} \]

\[ RMSE = \sqrt{\frac{\sum_{i=1}^{N}(z_i - \bar{z}_i)^2}{N}} \tag{8} \]

Where, \( z_i \) is the measured value, \( \bar{z}_i \) is the predicted value, and \( N \) is the total Number of observations (Yilmaz and Yuksek, 2009; Garg et al., 2015). \( R^2 \) is a descriptive measure between zero and one and indicates the ability of a parameter to predict another parameter. The highest value for MAPE (100) and the lowest value for RMSE (0) denote the highest values for model performance. MAPE usually gives accuracy as a percentage and a low percentage for this parameter denotes good performance of the model (Khoshnevisan et al., 2014). This research was carried out using SPSS 24.0, STATISTICA12 and Design Expert 7 software.

**RESULTS AND DISCUSSION**

**Evaluation of Linear Regression Model**

In this paper, FAME content and flash point are assumed to be a function of dielectric constant (\( \epsilon \)) and loss factor (\( \epsilon'' \)) at three different frequencies. The obtained LR models for prediction of FAME content and flash point are presented in Table 3.

\[ R^2 \] was used to ascertain the predictive performance of the model for the measured and predicted values. For example, it shows that \( R^2 \) for the relationships between the measured and predicted values obtained at frequency of 2,450 MHz from the LR models for FAME content and flash point demonstrate suitable correlation (Figure 3).

**Evaluation of MLP Model**

Input parameters of dielectric constant and loss factor were selected as input variables and FAME content and flash point were chosen as model outputs. The results showed that the best model was that with one input layer and two input variables, one hidden layer with 18 neurons, and one output layer and one output variable (2–18–1 structure). Cross-correlation of the predicted and target values at frequency of 2,450 MHz (Figure 4) indicated that the ANN MLP model was most acceptable. The \( R^2 \), RMSE and MAPE for the output variables are shown in Table 4.

**Evaluation of RBF Model**

The performance of RBF networks for estimated performance characteristics modeling used a single-layer neural network. Input directly entered the hidden layer cells and the output of the hidden layer cells multiplied by the weights entered a collector as the output of the neural network. Input-output data for network training were the same as that used to generalize the MLP network. The error value was zero and the network attained this error with 40 cells in the hidden layer. Figure 5 shows scatter plots of the

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**Table 3.** LR models for prediction of FAME content and flash point.

<table>
<thead>
<tr>
<th>Frequency (MHz)</th>
<th>FAME Content (%)</th>
<th>( R^2 )</th>
<th>Flash Point (°C)</th>
<th>( R^2 )</th>
</tr>
</thead>
<tbody>
<tr>
<td>434</td>
<td>( FC= 41.7\epsilon+50.6\epsilon''-110.6 )</td>
<td>0.995</td>
<td>( FP= -19.8\epsilon-34.5\epsilon''+257.8 )</td>
<td>0.933</td>
</tr>
<tr>
<td>915</td>
<td>( FC= 34.9\epsilon+52.6\epsilon''-86.5 )</td>
<td>0.968</td>
<td>( FP= -16.9\epsilon-35.3\epsilon''+246.7 )</td>
<td>0.920</td>
</tr>
<tr>
<td>2450</td>
<td>( FC= 43.6\epsilon+51.8\epsilon''-111.8 )</td>
<td>0.994</td>
<td>( FP= -21.1\epsilon-36.2\epsilon''+264.5 )</td>
<td>0.941</td>
</tr>
</tbody>
</table>
predicted FAME content and flash point based on the actual values.

$R^2$, RMSE and MAPE are shown in Table 4. The results show that the RBF model had high acceptability for prediction of output parameters at frequency of 2,450 MHz. The MLP, RBF, and RSM were applied for prediction of biodiesel production from crude mahua (*Madhuca indica*) oil and compared the results with those from a Multiple Regression (MR) model (Sarve et al., 2015). Comparison of factors indicated that RBF had the lowest MAPE and RMSE and the highest $R^2$. It can be concluded that the RBF performed much better than MLP and MR for predicting the FAME content and flash point.

**Comparison between the Models**

This study was developed and compared LR and ANN (MLP-RBF) methods for determining FAME content and flash point. The results of simple regression analysis showed a statistically suitable relationship between the output and input characteristics. The prediction models were developed with two inputs and outputs and the networks were evaluated for each output separately. The results of prediction of FAME content and flash point indicated that the equations obtained from the ANN-MLP model...
Table 4. Performance indices ($R^2$, RMSE and MAPE) for models at various frequencies.

<table>
<thead>
<tr>
<th>Frequency (MHz)</th>
<th>Model</th>
<th>FAME content</th>
<th>Flash point</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>$R^2$</td>
<td>RMSE</td>
<td>MAPE (%)</td>
</tr>
<tr>
<td>434</td>
<td>LR</td>
<td>0.995</td>
<td>0.011</td>
</tr>
<tr>
<td></td>
<td>ANN-MLP</td>
<td>0.977</td>
<td>0.008</td>
</tr>
<tr>
<td></td>
<td>ANN-RBF</td>
<td>0.998</td>
<td>0.005</td>
</tr>
<tr>
<td></td>
<td>LR</td>
<td>0.968</td>
<td>0.027</td>
</tr>
<tr>
<td>915</td>
<td>ANN-MLP</td>
<td>0.996</td>
<td>0.015</td>
</tr>
<tr>
<td></td>
<td>ANN-RBF</td>
<td>0.998</td>
<td>0.008</td>
</tr>
<tr>
<td></td>
<td>LR</td>
<td>0.994</td>
<td>0.012</td>
</tr>
<tr>
<td>2450</td>
<td>ANN-MLP</td>
<td>0.997</td>
<td>0.009</td>
</tr>
<tr>
<td></td>
<td>ANN-RBF</td>
<td>0.999</td>
<td>0.003</td>
</tr>
</tbody>
</table>

Figure 5. Cross-correlation of predicted and actual values of (a) FAME content and (b) Flash point for ANN (RBF) model at frequency of 2,450 MHz.

provided more suitable and reliable prediction than LR model. Besides, the RBF with two inputs and one output exhibited greater reliability of prediction than the LR and MLP models. Also, the values for $R^2$, RMSE and MAPE indicated that the prediction performance of the ANN-RBF model was better than LR and ANN-MLP (Table 4).

**Optimization Using Response Surface Method (RSM)**

In this paper, Box-Behnken design was employed to develop a relationship between biodiesel characteristics (FAME content and flash point) and independent variables (dielectric constant and loss factor) in order to maximize the FAME content and minimize the flash point. The FAME content and flash point varied from 43.1 to 90.8% and 147 to 183°C, respectively. The application of the advanced multiple regression analysis was employed to obtain the polynomial equation at frequency of 2,450 MHz. The equations of the significant terms obtained from the model in its coded form are as follows (Equations 9-10):

**FAME content**
\[
FAME\ content= -312.83 + 143.78 \varepsilon + 129.19 \varepsilon'' - 19.27 \varepsilon \varepsilon'' - 12.76 \varepsilon^2 - 8.92 \varepsilon'^2
\]
\[R^2= 0.998\]  \hspace{1cm}  \hspace{1cm}  \hspace{1cm}  \hspace{1cm}  (9)

**Flash Point**
\[
Flash\ Point= +245.32 - 27.46 \varepsilon + 48.65 \varepsilon'' - 22.03 \varepsilon \varepsilon'' + 3.25 \varepsilon^2 - 7.03 \varepsilon'^2
\]
\[R^2= 0.948\]  \hspace{1cm}  \hspace{1cm}  \hspace{1cm}  \hspace{1cm}  (10)

From the above equations, the coefficient with one factor signifies the effect in an individual form, while the coefficient which has two factors and second order form signifies the interaction between them and their fourth route effect. The suffix symbols positive or negative (+/-) signify...
the synergy and antagonistic effects, where the positive stands for synergistic effect and the negative stands for antagonistic effect (Joshi et al., 2008). Then, the model was analyzed by Analysis Of Variance (ANOVA) for obtaining the fitness of the model employing least square method. The effective statistical technique which bifurcates into individual roots which allows user to understand the sum of all the variation of the data in the model with specific sources of variation is ANOVA (Yatish et al., 2016). Thus, the obtained models variations are presented in Table 5.

Figure 6 shows an acceptable correlation between the predicted and experimental values of FAME content and flash point at frequency of 2,450 MHz, with a high value of coefficient of determination \( R^2 \) 0.998 and 0.948, respectively. The surface plots in three dimensions represent the graphical interface about the regression equation of reaction variables. Figure 7 represents the surface plots of biodiesel characteristics. Plots “a” and “b” illustrate the interaction between two independent variables on the dependent parameter. The plots are drawn with the aid of the regression equation and represent the interactions of each independent variable on the response variables. Plot a represents the significant interaction between dielectric constant \( \varepsilon \) and loss factor \( \varepsilon'' \), the variation in the FAME content is depicted well in the plot, i.e. the FAME content at first increases significantly by increasing both dielectric constant and loss factor. Plot “b” shows the nature of dielectric constant and loss factor on the flash point. Here, it could find a proportional increase in the flash point as the dielectric constant decreases, whereas a slight increase can be seen when the loss factor is decreased too.

Finally, to reach the optimization of biodiesel characteristics such that to maximize the FAME content and minimize the flash point values at frequency of 2,450 MHz, the conditions were suggested as in Table 6.

The optimum condition with desirability of 0.998 was obtained as dielectric constant of 3.72 and loss factor of 0.89 for FAME content of 89.88% and flash point of 152.7°C.

### CONCLUSIONS

Prediction of biodiesel characteristics based on its permittivity properties can help estimate the quality and simulate the production process. Production systems assess the conditions during processing and is complex because the data is generally inconsistent. This research work was applied to the LR, MLP and RBF models which were developed using these initial conditions to predict FAME content and flash point. The results showed for FAME content and flash point, \( R^2 \) values at frequency of 2450 MHz were calculated as 0.994 and 0.941 for LR model, 0.997 and 0.967 for MLP of ANN model, 0.999 and 0.988 for RBF of ANN model, respectively. The RBF model was exhibited high performance rather than the MLP and LR models for prediction the output data. The use of the RBF model may provide new

### Table 5. Experimental process obtained for biodiesel characteristics at frequency of 2,450 MHz.

<table>
<thead>
<tr>
<th>Run</th>
<th>Dielectric constant (( \varepsilon ))</th>
<th>Loss factor (( \varepsilon'' ))</th>
<th>FAME content (%)</th>
<th>Flash point (°C)</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>Predicted</td>
<td>Obtained</td>
<td>Predicted</td>
<td>Obtained</td>
</tr>
<tr>
<td>1</td>
<td>3.91</td>
<td>0.68</td>
<td>86.76</td>
<td>87.3</td>
</tr>
<tr>
<td>2</td>
<td>3.84</td>
<td>0.75</td>
<td>87.51</td>
<td>87.7</td>
</tr>
<tr>
<td>3</td>
<td>3.77</td>
<td>0.82</td>
<td>88.23</td>
<td>86.8</td>
</tr>
<tr>
<td>4</td>
<td>3.82</td>
<td>0.8</td>
<td>88.97</td>
<td>89.1</td>
</tr>
<tr>
<td>5</td>
<td>3.74</td>
<td>0.85</td>
<td>88.54</td>
<td>89.6</td>
</tr>
<tr>
<td>6</td>
<td>3.68</td>
<td>0.91</td>
<td>89.13</td>
<td>88.6</td>
</tr>
<tr>
<td>7</td>
<td>3.71</td>
<td>0.92</td>
<td>90.50</td>
<td>90.4</td>
</tr>
<tr>
<td>8</td>
<td>3.65</td>
<td>0.97</td>
<td>90.67</td>
<td>90.8</td>
</tr>
<tr>
<td>9</td>
<td>3.57</td>
<td>1.02</td>
<td>90.17</td>
<td>90.1</td>
</tr>
</tbody>
</table>
Figure 6. Actual values vs. predicted values of (a) FAME content and (b) Flash point at frequency of 2,450 MHz.

Figure 7. Surface plots of (a) FAME content and (b) Flash point at frequency of 2,450 MHz.

Table 6. Suggested conditions to reach optimum biodiesel characteristics.

<table>
<thead>
<tr>
<th>Number</th>
<th>Dielectric constant (ε)</th>
<th>Loss factor (ε&quot;)</th>
<th>FAME content (%)</th>
<th>Flash point (°C)</th>
<th>Desirability</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>3.72</td>
<td>0.89</td>
<td>89.88</td>
<td>152.7</td>
<td>0.998</td>
</tr>
<tr>
<td>2</td>
<td>3.72</td>
<td>0.88</td>
<td>89.52</td>
<td>152.9</td>
<td>0.997</td>
</tr>
<tr>
<td>3</td>
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Optimization of Fish Biodiesel Characteristics

approaches and methodologies, and minimize potential inconsistency of correlation. A comparison of the RBF network and MLP models indicates that RBF can be used efficiently to model and predict output data based on input parameters with similar accuracy. The optimum condition was obtained using response surface methodology as dielectric constant of 3.72, loss factor of 0.89 at frequency of 2450 MHz lead to reach the FAME content of 89.88% and flash point of 152.7°C with desirability of 0.998.

Nomenclature

\( A_p, A \) Peak area of standard and total sample (V s μA)

ANN Artificial Neural Network

\( b_j \) Model parameter

\( b_n \) Coefficient value

\( C \) Solution molarity (0.1 mol L\(^{-1}\))

\( CN \) Cetane Number

\( e \) Observed error

\( f \) Frequency (MHz)

FAME Free Acid Methyl Ester

FFA Free Fatty Acid

GC Gas Chromatography

LR Linear Regression

\( M \) Weight of sample (mg)

\( m \) WFO weight (g)

MAPE Mean Absolute Percentage Error

\( m_s \) Weight of standard in the sample (mg)

MLP Multi Layer Perceptron

MR Methanol Ratio

\( N \) Total number of observations

RBF Radial Basis Function

RMSE Root Mean Square Error

RSM Response Surface Method

\( V \) KOH solution volume (ml)

WFO Waste Fish Oil

\( X_i, X_j \) Coded independent variables

\( x_p \) Predictor variable

\( Z_i, Z' \) Measured and predicted value

\( \beta \) Regression coefficient

\( \varepsilon \) Complex dielectric

\( \varepsilon' \) Dielectric constant

\( \varepsilon'' \) Loss factor

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