

Theoretical and experimental ground of the fuel energy efficiency improvement by an activation of the burning reaction molecules-reagents

B. KOVALYSHYN

Department of Electric and Electro Technology, Educational and Research Institute of Energy and Automation, National University of Life and Environmental Sciences of Ukraine, 03041, Kyiv, Geroyiv Oborony 12, Ukraine, tel. 044-527-85-22, bikoval08@mail.ru

Received February 13.2012; accepted March 15.2012

Abstract. In theory and experimentally a method of fuel options energy efficiency rise due to an activation of burning reaction molecules-reagents under high voltage electric field action is grounded. A variant of the technical realization of the offered method is described.

Key words: high voltage, electric field, activation, energy efficiency, RedOx.

INTRODUCTION

The economy of fuel and energy resources due to the rise of efficiency of the use of power equipment is one of the basic tasks of modern science and production. Accordingly, a greater expenditure of power equipment per unit of products has a greater negative specific influence on environment. Therefore, a rise of the power equipment use efficiency is valid both for economy in general and, by necessity, for every industry from the viewpoint of economic and ecological efficiency of their functioning.

At present, traditional fossil power equipment is the basic energy source of practically all the economies of the world. Thermal energy sources are traditional and they supply to the most traditional power equipment. Thermal energy is created in the process of fossil fuels incineration in the oxidizing gas environment, that is at the motion of oxidizing-restoration chemical reactions. Therefore, at the use of power equipments optimization is an important task of the motion of chemical reactions.

The process of molecules-reagents activation at burning exothermic reactions is an object of our theoretical and experimental researches.

The technical and technological researches aim at the optimization of burning reaction of hydrocarbon fuels in the air under the high tension electric field.

The article presents the results of theoretical and experimental studies the target of which is the efficiency of motion of exothermic reactions of burning on the

example of incineration of hydrocarbon gaseous fuel in mid-air under high voltage electric field.

MATERIALS AND METHODS

The fuel heating value and temperature of flame are basic fuel properties. The fuel heating value (specific burning warmth) is the quantity of warmth (MJ/kg), which is selected at the complete burning of 1 kg of hard or liquid fuel, or 1 m³ of gaseous fuel. In practice, it is possible to distinguish higher Q_h^f and lower Q_l^f heating value of fuel [1]. A higher heating value is the description of fuel at its complete burning and condensations of the well-educated aquatic pair during burning. So, as in the real terms an aquatic pair is thrown out in the atmosphere, therefore in the heating engineering computations and reference literature under the term heating value of fuel we understand a lower heating value of fuel, in which the warmth of aquatic pair condensation is taken into account. The important question is whether it is possible to change the heating value of every traditional power equipment, or non-traditional one, of a strictly definite and constant size, by means of the proper technologies in the direction of increase.

Processes of generation of warmth are linked, mainly, with the reducing-oxidizing (RedOx) exothermic burning reactions. For the optimization of these processes it is expedient to consider them from the viewpoint of the chemical kinetics theory. The dependence of reactionary properties of the chemical system on the inlying energy, structure and atomic and molecular reagents composition is the basic question of modern theory of chemical kinetics.

Still in 1889 Arrhenius created the law of chemical kinetics [2], known as the law of his name. This law expresses the possibility of motion of chemical reactions between molecules-reagents. He links the constant of speed

of reaction with energy of activating (E_A), which characterizes the power state of molecule and is written down as:

$$k = k_0 \cdot e^{-\frac{E_A}{RT}}, \quad (1)$$

where: k_0 - preexponential constant; R - gas constant, even 1,987 cal/grad·mol; T - temperature in the degrees of Kelvin scale, °K; e - basis of natural logarithms.

To find an activating energy, reaction speed is caat a different temperature and for every value of speed constant. After taking the logarithm of both parts of formula (1) get a formula:

$$\ln k = \ln k_0 - \frac{E_A}{RT}. \quad (2)$$

For a greater comfort of computations decimal logarithms are used:

$$\lg k = \lg k_0 - \frac{E_A}{4,575T}, \quad (3)$$

where: the number 1/4,575 - module of translation of natural decimal logarithms, increased by $R = 1,987$.

Energy of activation can be defined with the formula

$$E_A = (\lg k_0 - \lg k) 4,575T. \quad (4)$$

From formula (4) it is evident, that the energy of activation of molecules relies straight proportionally on temperature.

A process of activation of the molecular system consists in the transformation of power state of electrons on the higher power level.

The explanation of the given process can be illustrated by means of power diagram of atom (molecules) in Fig.1

A process of excitation of atom (molecules) consists in transition under action of power factor of one or a few valency electrons from the stationary shell in atom (molecule) on higher excitation levels.

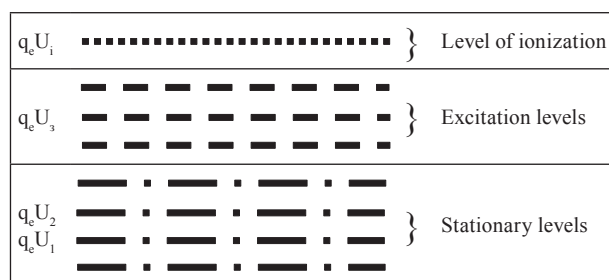


Fig. 1. Power diagram of atom

To release an electron from a molecule of methane, one should get the quantum energy lower than 12.6 eV [3], for ethane and propane – 11.5 eV, for oxygen molecules - 15.5 eV [4] - the quantum of energy lower than the energy of ionization, but sufficient to transfer an electron in an excited state. For computations we take the value of quantum energy of activation molecules about 5% lower than the energy of ionization.

Molecules activation by light is the result of inelastic collision of molecules and quantum of light [5]. With sufficient quantum energy of electromagnetic radiation, electrons move from molecule stationary energy levels to excitation levels. This quantum energy of light (E_l) is calculated by the formula:

$$E_l = h_0 \nu, \quad (5)$$

where: h_0 - Planck's constant ($6,626 \cdot 10^{-34} \cdot \text{J} \cdot \text{s}$); ν - electromagnetic wave frequency (Hz).

The frequency was calculated of electromagnetic radiation needed to activate the aforementioned molecules and the results were put into Table 1.

As seen from the obtained results, for the activation of molecules the so called vacuum ultraviolet rays can be used with a wavelength within 100 nm or less. Getting radiation with such parameters in normal states is problematic. Therefore, photoactivation of molecules without special training parameters of the environment is not effective.

After our working hypothesis, for activating of molecules-reagents, except for thermal, it is possible to use other outsourcings with energy. If external energy sources are constant and secure the source of pulsating energy, frequency of which answers resonance frequency of molecules of reagents, it is possible to attain an effect of their translation in the active voice with the substantially lower power expenditure. Complementing the formula (4) by the effect of action on molecules-reagents we will get other factors:

$$E_A = E_{AT} - E_{A3} = [(\lg k_0 - \lg k) 4,575T] - Wb, \quad (6)$$

where: W - energy obtained from outsourcings; b - coefficient of the use of external energy by molecules-reagents.

From formula (6) it is evident that thermal energy of activation can be lowered due to the use of other sources of molecules activation.

Table 1. Options of electromagnetic radiation activation

Molecule	Ionization energy, eV	Activation energy			Light frequency, Hz	Electromagnetic wave length, nm
		eV	J / molecule	J / mol		
Methane	12,6	11,91	19,08·10 ⁻¹⁹	11,48·10 ⁵	2,87·10 ¹⁵	104
Ethane	11,5	10,87	17,41·10 ⁻¹⁹	10,48·10 ⁵	2,63·10 ¹⁵	114
Propane	11,5	10,87	17,41·10 ⁻¹⁹	10,48·10 ⁵	2,63·10 ¹⁵	114
Oxygen	15,5	14,65	23,47·10 ⁻¹⁹	14,12·10 ⁵	3,54·10 ¹⁵	85

There were three repetitions of the experiment. The research results are given in Table 2.

The obtained experimental results show that electro-activation of burning exothermic reaction components of propane in mid-air reduces the heating time of 1 litre of water in an old tub practically in all variants of research. The most positive effect is observed at the action on both of the components of reaction of burning of high-voltage pulsating electric field with frequency 100-120 Hertz.

CONCLUSIONS

1. Activation of molecules-components of burning reaction by the high-voltage pulsating electric field results in the reduction of expenditure of fuel.

2. The highest efficiency of electro-activation (about 22 %) is obtained at the frequency of high-voltage pulsating electric field 120 Hertz and the activation of both the components of burning reaction.

REFERENCES

1. **Korchemnyj M.** Energozberezhennia w agropromyslowomu kompleksi. W. – Ternopil: Pidruchnyky i posibnyky. 2001 – pp. 976.

2. Fizicheskaja Izhimia. Pod red. K. S. Krasnowa. M.: Wyszchaja shkola. 2001. t. 1. pp. 512; t. 2. pp. 319.

3. The Open University/ <http://openlearn.open.ac.uk/mod/ou-content>

4. Science Encyclopedia / <http://science.jrank.org/pages/3677/Ion-Ionization-energy.html>

5. Patent Nr 17392 Ukraine. Sposibotrymannia molekul z zadanyimi vlastywestiami khimichnykh zwiazkiw Melnykow W.M., Usachow A.W.

6. Patent Nr 52845. Pristrij pidhotowki okisluwacha do spalennia paliwa. Maltsew W.O., Kushniv W.M., Pedos W.A., Nikołajew M.M.

7. **Bagriatyshwili W.N.** Unogofotonnyje protsessy w molekulakh z infrakrosnom lazernom pole. M.: izd-wo "Nauka" 1988. pp. 245.

8. Patent Nr 24193 Ukraina MKP sposib pidhotuwannia okisluwacha do spaluwannia to prustrij dla joho zdijsnennia, Sklab W.S., Owsij O.W.

9. Patent Nr 37572 Ukraina MKP Sposib pidwushchennia efektywnosti i podiwnykh ustanowok na wuhlewodnewomu paliwi ta prustrij dla joho realizatsii, Kowalysyn B.M.