Enhancement the octane number of sudanese gasoline (Nile Blend)

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ABSTRACT

Gasoline is one of most important sources of energy but when manufactured with low octane number may leave great risk that affect the human health and environment impact. Therefore processes of gasoline production must be developed. Zeolite material dependence on their extremely high frame work extended interactive surface, ability to interact with molecule and cations. They show a wide variety of interesting properties which result in remarkable applications in the fields of catalysts, gas separations, selective adsorption and water softening. Zeolite catalyst loaded with finely (Sr) metal have been used in hydro-cracking, catalytic cracking. The zeolite used is the synthetic zeolite (type Y) has high thermal stability, good surface area and its heterogeneous catalyst which can be separated after treatment and may be reactivated. The study of the octane number was carried out for this zeolite. The gasoline was obtained from market (commercial gasoline) it has octane number 91.7 determined by IROX-2000.The gasoline was treated by zeolite in the temperature about 331 Kelvin. The synthetic zeolite (type-Y) raised the octane number (from 91.7 to 92.4).

Keywords: octane number; sudanese gasoline; Nile Blend

1. INTRODUCTION

Sudan has two sorts of crude. They are different in quality and price. Sudan's Nile Blend crude is sold at much higher prices than Dar Blend crude(1). Nile blend is the one of the best crude qualities, it contains a little amount of impurities and high chemical properties, which require low invested and operational costs to produce best quality products. It is of the medium density, which reach 34 API degree, its pour point range between 33-36 ^{O}C centigrade, and sulphur composition content is low, not exceeding 4% weight. However its complicated aromatic complex content can be ignored from the practical aspect (2). As crude oil comes from its well containing a mixture of hydrocarbon compounds and relatively small quantities of other materials such as oxygen, nitrogen, sulphur salt and water. In the refinery most of these non-hydrocarbon substances are removed and the oil is broken down into its various components, and blended into useful products (3). These component can be separated easily as they have different boiling temperatures, by fractional distillation (4). At about 40-70º C gasoline, which it is largely

a mixture of hydrocarbons, can be separated. Some may also contain significant quantities of ethanol and \ or small quantities of additives such as tert-arybutylmethyl ether as anti-knock agents to increase the octane rating. The hydrocarbons consist of a mixture of n-paraffins, naphthenes, olefins and aromatics. Naphthenes, olefins and aromatics increase the octane rating of the gasoline whereas the n-paraffins have the opposite effect. The aromatics are mostly a mixture of benzene, toluene and the xylenes. The benzene content is kept to a minimum (but is not negligible) due to its perceived toxicity. Originally lead tetraethyl was added as an anti-knock agent but is now rarely used and is prohibited in most countries due to its toxicity(5)

By definition an octane number is that percentage of isooctane in a blend of isooctane and normal heptane that exactly match the knock behavior of the gasoline. Thus of octane gasoline matches the knock characteristic of blend containing 90% isooctane and 10% n-heptane. The octane numbers, as defined by ASTM methods, is an empirical property. It cannot be submitted to any analysis that would allow octane optimizing strategies. There fore it must be related to measurable properties inherent to gasoline, such as its molecular composition. The relation between the structure of hydrocarbons and their octane number was studied using a number of topological indices. A physicochemical discussion about the meaning of octane number on the molecular level, however, did not go beyond some empirical rules:

- ON increases with the number of tertiary and quaternary carbon atoms [2,6] respectively it increases with the number of methyl groups. [13-15]
- It decreases with the total number of carbon atoms [2,6] respectively with the length of the chain [13-15]
- ON gets larger as the branching point is moving toward the center of the longest chain in the alkane molecule [6].

Zeolites are micro-porous crystalline solids with well-defined structures. Generally they contain silicon, aluminium and oxygen in their framework and cations, water and/or other molecules within their pores [7]. The natural zeolite faujasite has the same framework (FAU) and similar framework composition to the Type Y synthetic zeolite but is rare in nature [8]. From the early 1960s, use of synthetic zeolites in catalysis and in related adsorption separation processes has dramatically transformed petroleum refining by vastly increasing the yield of high-quality fuels and reducing capital and operating costs, energy requirements, and adverse environmental impact [9] In modern petroleum refineries in the world, gasoil and other heavier fractions from the crude oil fractionation unit are fed to fluid catalytic cracking units, which use small, fluidizable catalyst particles containing Type Y zeolite or other zeolites, or to hydro cracking units, which use fixed beds of larger catalyst particles also containing zeolites [10 -12].

Many zeolites can be synthesized with SiO_2 higher or lower than in nature for the same framework type. Higher SiO_2 generally gives greater hydrothermal stability, stronger-acid catalytic activity, and greater hydrophobicity as adsorbents. Conversely, lower SiO_2 gives greater cation exchange capacity and higher adsorbance for polar molecules. Controlling the synthesis process optimizes a zeolite for different applications [13].

The fluid catalytic cracking and hydrocracking units convert higher-molecular-weight hydrocarbons to lighter ones suitable for gasoline, light fuel oils, olefins, and other uses [14]. Zeolite Y with Si/Al = 2.5 - 3 stability is increased by ion exchange with rare earth cations [15] ZSM - 5 with pore size $\approx$ 1.2 nm was reported to have a high reaction selectivity for n-alkanes cracking over acidic centers by forming a secondary or tertiary carbonium ions from a beta C-C bond. This process leads to production of propenes or

butenes from long chain n-alkanes [15].

Unfortunately, tetraethyl lead and other metal and non-metal additives for octane boosting or octane improvement of gasoline such as other lead compounds, manganese compounds, boron compounds, phosphorous compounds, etc., deactivate or poison the catalyst in the catalytic converters so that the converter cannot reduce the noxious emissions to acceptable levels [16]. Therefore the use of zeolites as safe natural material was done in this study. Zeolites were prepared from clays and materials available in Sudan. The synthesized zeolite was employed in enhancement of octane number of fuel.

2. MATERIALS AND METHODS

2.1. Materials

The chemical below were obtained from natural sources as in table 1. They were used as received without further purification:

Table 1. Raw material for the synthesis of zeolite.

Name	Source	Purity
Clay	River Nile	Natural
Sodium carbonate	Alfasher	Natural
Strontium nitrate	Loba India	Analytical Reagent
K-Feldspar $\underline{KAlSi_3O_8}$	Bioda desert	Natural
Sudanese gasoline	Refinaery	

2.2. Instruments

Thermal Gravimetric Analysis (TGA), FT-IR (SHIMADZU 84005), IROX-2000, Density meter (Anton par), Sensitive balance, water bath, Electric Furnace (1273 K – 1437 K Carbolite)

2.3. Preparation of synthetic zeolite

The procedure of preparation of synthetic zeolite consists of the following steps:

a-The K feldspar was extracted from rocks with jaw and crusher, then crushed till powder,

b- The clay was crushed with mortar and pestle.

c- The trona (sodium carbonate) was crushed till became powder.

d- The mixture of clay, K- feldspar, sodium carbonate and strontium nitrate were placed in the clean crucible of porcelain as in the following ratios clay: K-feldspar: sodium carbonate: strontium nitrate (39.2 : 39.2 :19.6 : 1.96).

e- The mixture was heated in the electric furnace at a temperature of 1373 K for six hours.

f- After that the crucible was taken away from the furnace and cooled. Finally the synthetic zeolite was crushed and was ready to use. Synthesized zeolite was characterized by XRD, TGA (thermal gravimetric analysis) and FT-IR.

XRD, FT-IR and TGA analysis of zeolite

X-ray diffraction (XRD) for zeolites A was analyzed following the standard method [17]. Experiment was performed using a Philips PW1732/10 diffractometer and CuKα radiation and Ni filter. The working conditions of this experiment were 40 kV and 20 mA. Also zeolites A was analyzed by Fourier Transform Infra Red(FT-IR), in which a mixture of 0.005g of zeolite A and 0.2g of KBr were pressed into disk for IR measurement [18]. Thermo gravimetric analyzer instrument manual, the STA PT1600 is a highly versatile research to controlled temperature environment (TG). Temperature range 323 K to 723 K and heating rate (273.1 to 373 K/min), max sample weight 25g , resolution 0.5 μg. in this research we can analyzed at 723 K and rate 10° C min.

Treatment process of gasoline by zeolite

In the clean bottle the 2.5 liters of gasoline was taken, and 119 g of zeolite were added to this gasoline the mixture was shaken gently and it was put in a water bath at ≈ 331 K till boiling of gasoline, and time was recorded. After that the mixture was cooled and the octane number, density, copper strip corrosion of gasoline was measured.

Determination of the octane number

For determination the research octane number (RON) the American Society for Testing and Materials (ASTM) method (ASTM D2699) was used. About 150 of gasoline was taken to measure the octane number of gasoline after and before treatment, by using IROX-2000 (is an extremely compact, robust and user friendly Mid-FTIR spectrometer for the automatic measurement of the concentration of the most important components of gasoline). This instrument was operated in the following conditions; display in the large graphics display, back lit. warm up time 10 min response time 3 minutes and the retention time of the treated sample is 7 days.

Determination of density

For determination of density the ASTM method (D-1298) was used. A small volume (Approximately 0.7 ml) of gasoline was introduced into an oscillating sample tube and the change in oscillating frequency caused by the change of the mass of tube which was used in conjunction with calibration data to determine the density of the sample.

Copper strip corrosion

A polished copper strip is immersed in a specific volume of the sample being tested and heated under conditions of temperature and time (3h/50°c) that are specific to the class of material being tested. At the end of the heating period, the copper strip is removed, washed and the color and tarnish level assessed against the ASTM (D130) Copper Strip Corrosion standard.

Fig. 1. Copper strip corrosion instrument.

3. RESULTS AND DISCUSSION

The XRD pattern of the synthesized is shown in figure (2). The characteristic peaks of Y zeolite at $2\Theta = 6.5°$ (2), 10.1° (3), 15.6° (6) and 23.7°(7) [19] indicate the its formation. The pattern is matching with that of commercial zeolite Y (Fig. 2b).
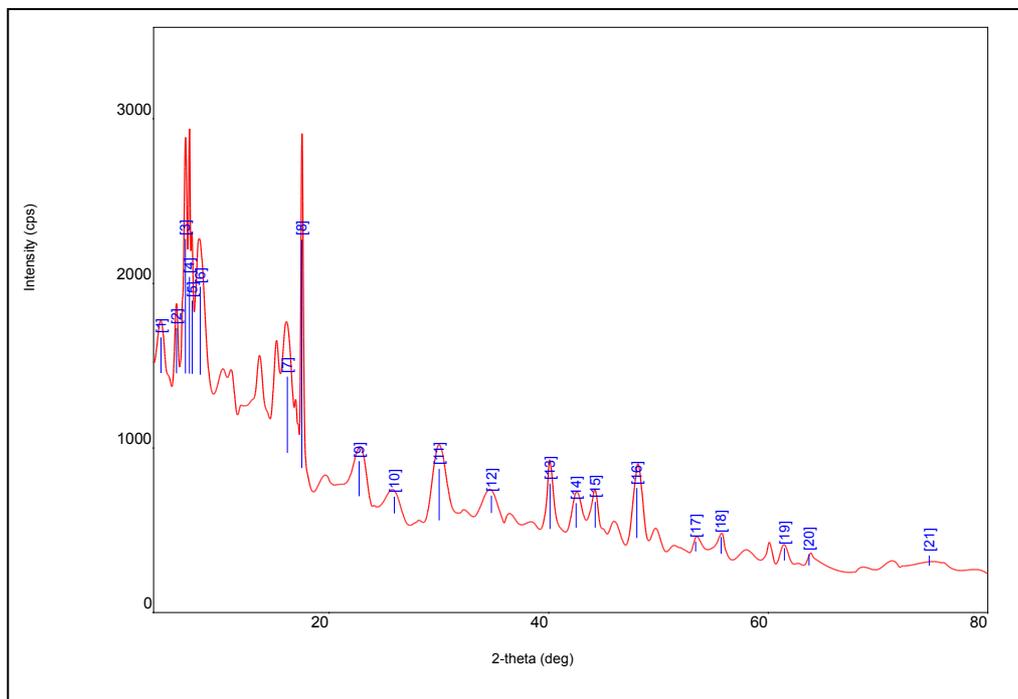
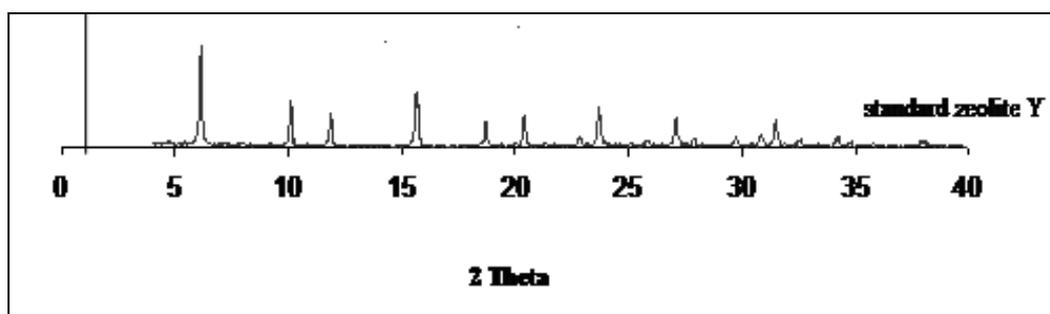**Fig. 2a.** XRD pattern of synthesized Y zeolite.

Fig. 2b. XRD pattern of commercial zeolite Y.

The product was characterized by Fourier transform infra-red (FT-IR) spectroscopy. The IR of zeolite after treatment (Fig. 3) shows strong IR absorption in the spectra region bellow 1200 cm^{-1} vibrations frequencies of the zeolite lattice which results from stretching and bending modes of the T-O units which is observed in the range 400 – 1100 cm^{-1} this indicates that SiO_2 or Al_2O_3 are linked. The band at the 1018 and 773 referred to antisymmetrical and symmetrical stretching of Si-O-Si respectively, and the band centered at 3411 cm^{-1} due to the water molecules in the zeolite [20]. The band at 2356 cm^{-1} attributed to the organic compounds (mercabtanes) which adsorbed by zeolite. The FTIR spectra of synthesized zeolite Y is agrees with the typical FTIR absorption peaks of zeolite Y reported earlier [21].
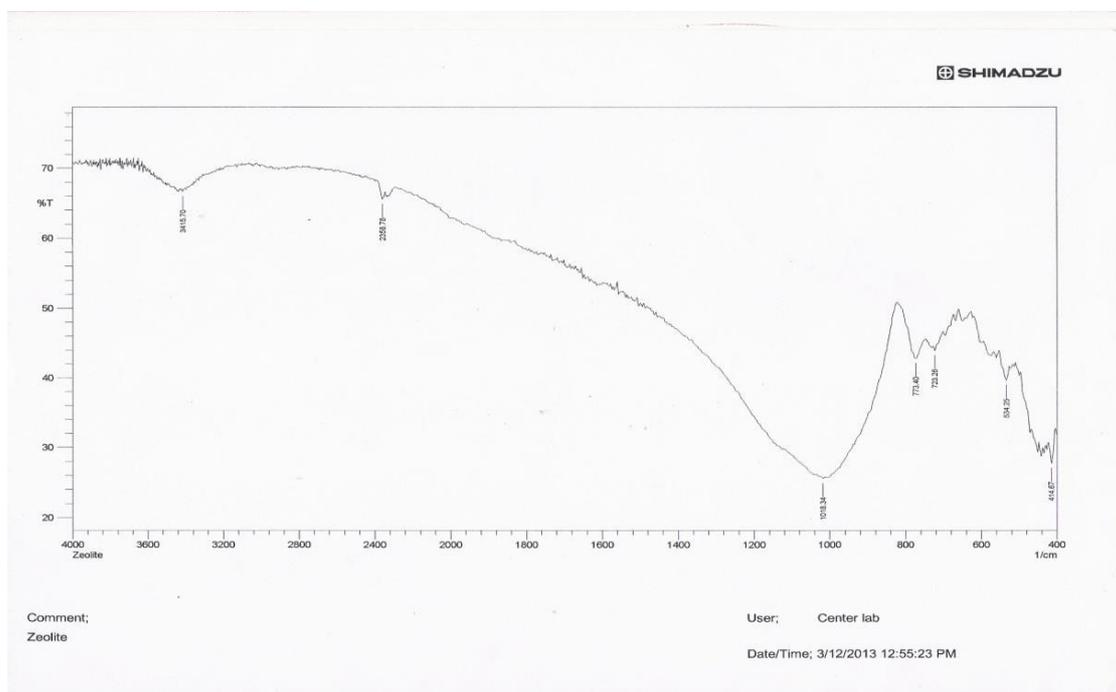

Fig. 3. FT-IR of zeolite after treatment.

The thermal gravimetric analysis (TGA) thermo grams of sample of zeolite before and after treatment are shown in figures (4a) and (4b) the zeolite before treatment underwent

two weigh losses, the endothermic loss between 50 ^{o}C and 100°C is assigned to water adsorption, while for zeolite sample the water molecule found in its structure can be lost in the range between 130° C and 250° C. These losses are the same for the two samples of zeolite (before and after treatment) but the difference between them is the loss of the organic compounds which are adsorbed by zeolite and it is shown in figure (3-2) in the range between 90° C and 200° C. Zeolite has high thermal stability and it is stable until 500° C , this is the desired property for any heterogeneous catalyst, and it can be separated and recovered easily after treatment of gasoline. In the case of the type II zeolite, the increment of the octane number seems to be due to the high aromatics content [22].
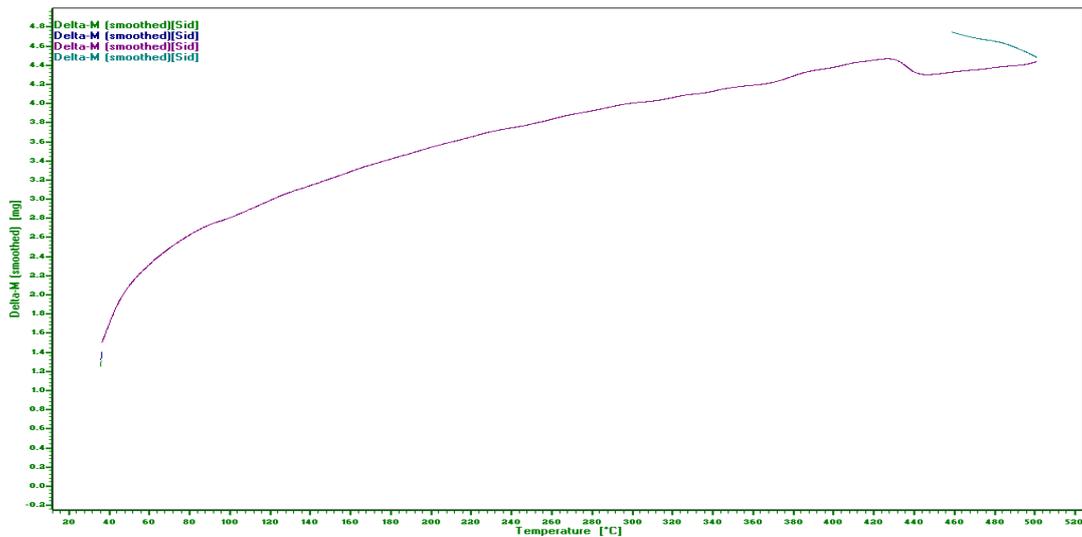

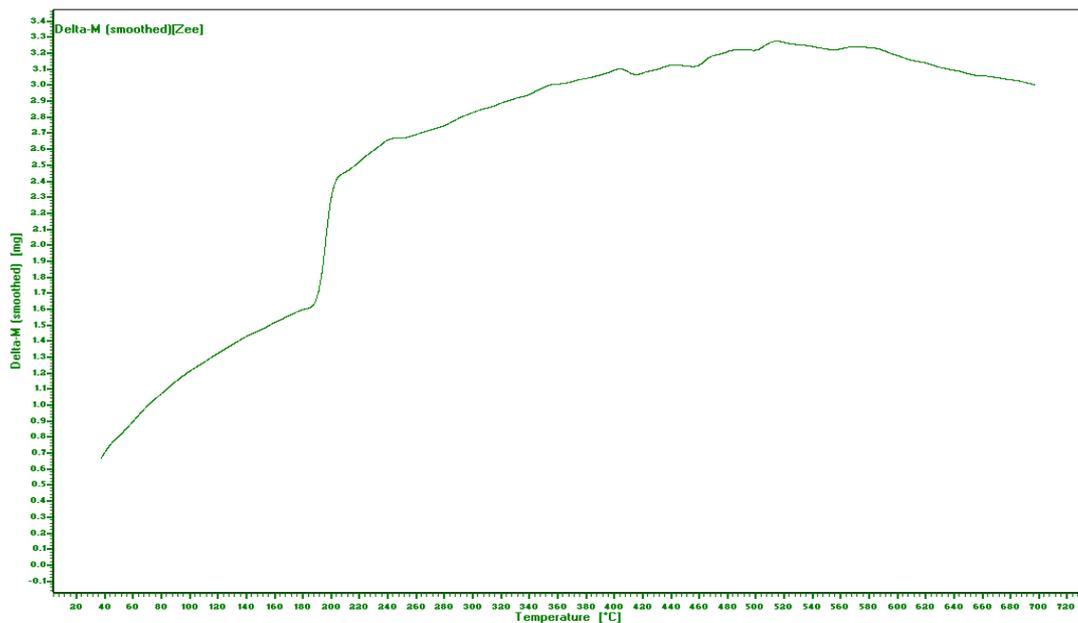**Fig. 4a.** TGA diagram of zeolite before treatment.

Fig. 4b. TGA diagram of zeolite after treatment.

3.1. Octane number, density and corrosion rate

Table 2. Octane number, density and corrosion rate of gasoline before and after treatment with zeolite.

Sample	Research Octane Number	Density $(g.cm^{-3})$	Copper strip corrosion rate
Gasoline without treatment	91.7	0.7320	(3h/50 °C) 1a
Gasoline after treatment	92.4	0.7285	(3h/50 °C) 1a

The use of zeolite in refinery is very important to solve many problems of fuel, and it had many uses in refinery processes to produce friendly products. As the result of treatment the octane number of gasoline increased from 91.7 to 92.4 by zeolite. The empirical rules of octane number dependence on the structure of alkanes can be amended to become as follows: octane number decreases with the increasing of number of CH2 groups and arises with the increasing of the number of CH3 groups; the number of adjacent CH2 groups has the highest but sigmoid influence; octane number decreases with the separation between branches; − ON increases with the more central position of branches; − ON increases with the bulkiness of the branched structure. The zeolite destroys the alkane side chain by absorbing the CH2 which yields short side chain. This reason leads to enhancement of the octane number by zeolite which can be separated after treatment. This method is simple and economic.

The treatment process of gasoline by zeolite involves taking heavy oil such as kerosene or diesel and heating it to a high temperature in the presence of a catalyst. The large molecule breaks down into several smaller ones, some saturated, some unsaturated. The unsaturated products are used as feedstock for the polymer industry while the saturated products are usually high-octane branched chain alkanes suitable for making gasoline. In this process the density of gasoline was decreased after treatment from (0.7320 to 0.7285 g/cm^3) this result lead to say when the octane number increase the density was decreased.

Crude petroleum contains sulfur compounds, most of which are removed during refining. However, of the sulfur compounds remaining in the petroleum product, some can have a corroding action on various metals and this corrosivity is not necessarily related directly to the total sulfur content. The effect can vary according to the chemical types of sulfur compounds present. The copper strip corrosion test is designed to assess the relative degree of corrosivity of a petroleum product. There was no deference between the blank and sample according to the corrosive effect of the copper strip. According to this result the copper strip corrosion of the sample was found in the normal range.

4. CONCLUSION

Zeolite are stable, non-toxic and preventing contamination of valuable feed stock's. zeolite is used as heterogeneous catalyst due to stability to cavity, its thermal at high temperature and secretive even at unfavorable reactant ratio and the reaction is eco-friendly. the synthetic zeolite is high activity more than natural zeolite because it is have high purity. Use of zeolites as a catalyst in the manufacture of some fine chemicals should

expand. New zeolite catalyst should be developed with improve application selectivity and new functionalities, perhaps for strong base and oxidation catalysis. New ion exchange of new microporous oxide also may be expected. Experience has taught that availability of new material normally precedes by many years the discovery of all of their useful properties and the conception and development of new uses.

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