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The phosphate series of Benguerir (Maastrichtian - Ypresian, Morocco): Mineralogy and mine planning

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Abstract

Morocco is the world leader in phosphates. Exploration and reserves evaluation of deposits in terms of tonnage and quality are charged to the geological service of the Office Cherifien of Phosphates (OCP). The Maastrichtian-Ypresian exploitable phosphates from the northen deposit of Benguérir, are transgressive on a Paleozoic basement. The coordinates of the samples and their physicochemical characteristics allowed building a database. The processing of these data by the ArcGIS software allowed us to propose maps of the different properties of the deposit, such as BPL (Bone Phosphate of Lime (BPL = 2, 18 * % Phosphorus pentoxideP2O5) and CO₂ contents, as well as structural characters such as the thicknesses and structures of the phosphate layers. This work offers a quick and efficient method for the evaluation of resources and the definition of deposit structures to assist the operator in the extraction procedures.

1. Introduction

Morocco has the largest deposits of phosphates in the world with estimations of more than three quarters of the world's reserves. Their geographical location, marketability and diversity give Morocco a special place in international market. As the world's largest exporter of phosphates in all its forms and the third largest producer behind the United States and China [1-3], its production reaches approximately 26.6 million tons per year [4,5]. It is the Cherifien Phosphates Office (OCP), which holds the monopoly of the exploitation and extracts the raw phosphates from the Moroccan undergrounds thanks to open-pit sites [6]. Almost half of the ore is exported as raw material, and the other half is delivered to the group's chemical industries for processing into marketable by-products such as basic phosphoric acid, purified phosphoric acid and solid fertilizers [7].

Beyond this very constructive work, we have assigned ourselves the task of proceeding, by our own methods; to the processing of mining data of a well-defined area, using techniques of Geographic Information Systems in order to have tools helping geologists to better visualize the structure of the phosphate deposit and evaluate its resources. The ArcGIS software has been used for this purpose. The problem defined consists of:

- Conducting a geological study of the sector to capture the sedimentary and structural characteristics and link them to the paleogeography of the time of their deposit;
- Studying the petrography of representative phosphate layers and specifying their chemomineralogical properties;

• Collecting and compiling the mining data to form a database that can be processed by the software for geomineral purposes.

2. Material and Methods

2.1. Presentation of the study area

The Ganntour basin corresponds, on the surface, to a 120 Km long and 20 to 30 Km wide east-west oriented rectangle (Figure 1). In the Benguerir deposit, the phosphate series range from the Maastrichtian (last Cretaceous) to the Lutetian (Middle Eocene) with different levels in the form of either phosphate beds (layer, furrows) or sterile to slightly intercalated phosphate levels [8]. The phosphate series of the Benguerir deposit spreads from Maastrichtian to Lutetian with different levels forming phosphate beds (layers and furrows), phosphate-reduced (spacer) or sterile.

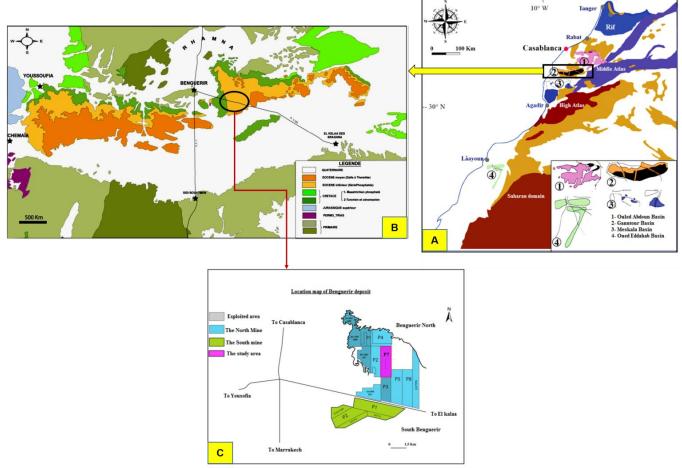


Figure 1: Situation of the study area: A- Structural scheme of Morocco and location of Moroccan phosphate basins [9] modified by EL Haddi [10, 11]. B- Geological map of the Ganntour basin extracted from the geological map of Morocco at 1:500000. [12]. C- Location of panel 7 of the Benguérir deposit [13].

2.2. Contribution of X-ray diffraction

Two samples were subject to an X-ray diffractometry mineralogical analysis. The objectives were, besides confirming and possibly completing the mineralogical study, to define the mineralogy of two reference levels: the yellow level and the C4-C3 intercalary (Spacer), both sterile.X-ray diffraction is a non-destructive analysis technique used for the determination of single-crystal structure as well as the study of qualitative and quantitative analysis of mineral phases.For X-ray diffraction, Panalytical X 'Pert Pro MPD diffractometer has been used for the identification and study of polycrystalline samples (pulverulent or solid). With the following characteristics: an X-ray tube source with copper anode of $\lambda = 1.54 \text{ A}^\circ$; a vertical O-O configuration goniometer; a rotating sample holder called spinner; a fast linear solid detector X 'Celerator; a 45-position autosampler.

2.3. Morphological characterization by scanning electron microscopy

These observations were made using a 20 kV voltage scanning electron microscope (Oxford ISIS 1300 ESM) coupled to a EDXA microanalysis system (Energy Dispersive X-ray Analysis) recording the X-ray fluorescence lines emitted after electronic excitation. The spatial resolution in imaging mode (secondary electrons and backscattered electrons) is less than 0.1μ m. The backscattered electron images make allow obtaining in certain cases, depending on the state of the surface, a good crystallographic contrast. The instrument also allows the analysis of elements up to uranium with detection limits as low as 1%.

2.4. Mine planning

The database is intended to be processed by ArcGIS software. It is based on data from 108 well sections and other recognition documents developed by the mining geology department of Benguerir [14]. The preparation of the database which contains the necessary information on the parameters studied by the geoprocessing on ArcGIS software is carried out according to several stages: Identification of phosphate levels and sterile levels:

- Determination of the powers of phosphate levels (PM:the thicknesses of the phosphate layers) and sterile levels (sterile P);
- Calculation of the physico-chemical parameters (BPL: Bone Phosphate of Lime (BPL = 2, 18 * % Phosphorus pentoxide P2O5)), CO2:Carbon dioxide) of each phosphate level;
- Elimination of outliers of the studied parameters.
- Notation of the coordinates (X, Y and Z of each well).

3. Results and discussion

3.1. Benchmark level

Although the SEM study showed the existence of phyllitous minerals, the results of the X-ray diffraction do not show evidence of clay mineral abundance [15]. This mineralogical study is fragmentary for it only concerned a single sample, not representative of the entire thickness of the layer and its possible lithological variations. Moreover, the resolution chosen for X-ray diffraction may not be able to reveal the peaks of these minerals. It is appropriate for an objective study to proceed with a sedimentological step of phases separation before an objective X-ray diffraction study. At this stage, this study is not considered conclusive to the mineralogy and the origin of this yellow landmark layer.

However, calcite and dolomite appear in the analysed sample (Benchmark level) as the most abundant minerals (Table 1). These minerals are often associated with, however, a clear predominance of dolomite over calcite (Figure 2), defining a dolomite or limestone dolomite for this yellow level.

Compound Name	Chemical Formula	PDF Index Name
Dolomite	Ca Mg (C O3)2	Calcium Magnesium Carbonate
Calcite	Ca (C O3)	Calcium Carbonate

Table1: The abundant minerals in the analyzed sample.

The dolomite is characterized in this case by a broad and intense peak at 3.91 Å (22.87 °2Th), while calcite is characterized by peaks with lower intensities. From a mineralogical point of view, the line (interplanar distance) of calcite is generally between 1.03 Å and 3.86 Å (23.05 °2Th), and for dolomite, its main line range between 2.89 Å and 2.90 Å. The yellow color, often attributed to iron oxides in clay sedimentary rocks, is in this case. The presence of these two components is not obvious at first sight; this may be related to phosphate rock leaching and carbonate impregnation. The significance of these benchmark levels remains therefore posed.

3.2. Spacer C4 - C3

X-ray diffraction analyzes show two dominant crystalline forms (Table 2), which are calcium carbonate and hydroxyalpatite with lines that appear respectively at 3.02 Å (forcalcium carbonate) and 3.16 Å (forhydroxyalpatite) (Figure 3).

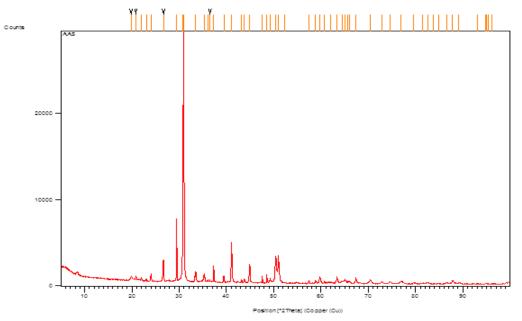
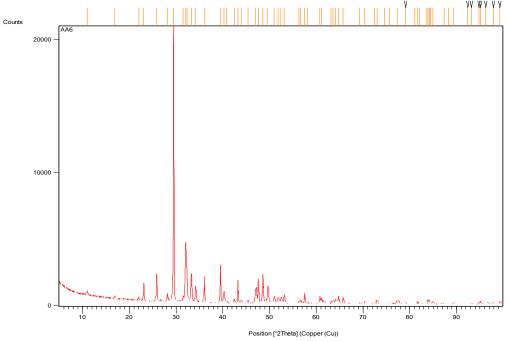


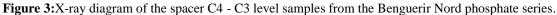
Figure 2: X-ray diagram of clay level samples from the Benguerir North phosphate series

Compound Name	Chemical Formula	PDF Index Name	Semi Quant [%]
Calcium Carbonate	Ca C O3	Calcium Carbonate	66
Hydroxylapatite	Ca5 (P O4)3 (O H)	Calcium Phosphate Hydroxide	34

Table2: The abundant minerals in the analyzed sample.

Calcium carbonate forms 60% of the rock (Table 2); it is therefore a phosphate limestone. Petrographic and SEM studies show that these carbonates are essentially oolithic.





3.3. Yellow benchmark level

The images of the sampleobtained by scanning electron microscope (SEM) show four types of crystals (Figure 4 and Figure 5):

- Crystals in rhombohedral form: are the most frequent, with a size varying around 43.3µm. These crystals thus correspond to dolomite and calcite already confirmed by X-ray study.
- 29µm cubic crystals which probably have to correspond to halite crystals (rock salt).
- Crystals in laminated form correspond to phyllosilicates (Clays).
- Irregularly shaped crystals that probably correspond to quartz.
- Thus, these observations, in contrast with the results of the RX, confirm the presence of clays in this reference level.
- The qualitative chemical analysis carried out with the energy dispersive micro-analyzer (EDXA) accounts for the presence of the Ca, Si, Mg, as well as Al and Fe elements (Figure 6) which enter into the compositions of all the raised minerals including clays.

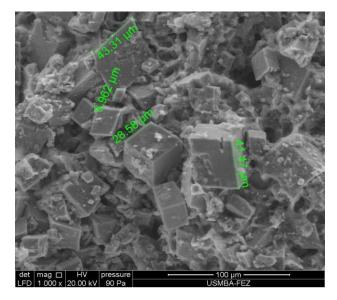


Figure 4: Images of the scanning electron microscopy of the benchmark level.

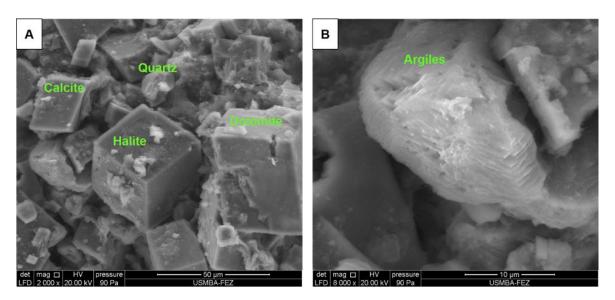


Figure 5: The different shapes of the crystals.A- Rhombohedral and cubic. B- Laminated.

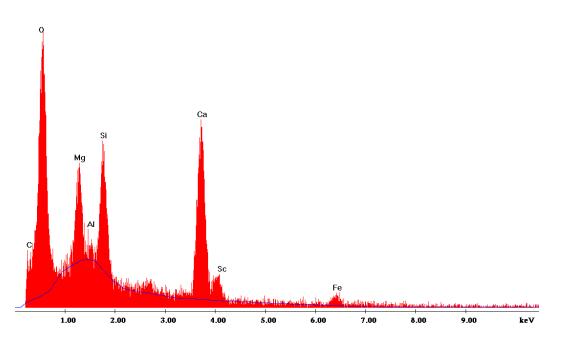


Figure 6: Spectrum of chemical elements of benchmark level (SEM).

3.4. Spacer layer 4 - layer 3

Observation under a scanning electron microscope (SEM) shows that this level consists essentially of oolites of size ranging between $396\mu m$ and $410\mu m$, and of rounded or sometimes irregular shapes, bone debris, as well as quartz grains [16]. X-rays showed the predominance of carbonates and apatite in the composition of this level (Figure 7).

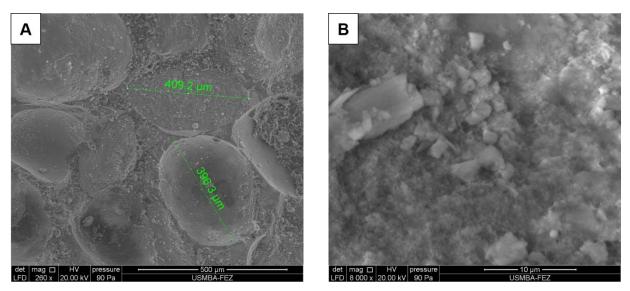


Figure 7: SEM image of the structure of phosphate constituents. A- The ooliths. B- Bone fragments.

Qualitative analyzes using a scanning electron microscope confirm that the level consists of phosphate components associated with carbonates. The presence of clay minerals is not to be excluded in this interlayer as well as organic matter. Sc (Scandium) is also a constituent element (Figure 8).

3.5. Geographic Information System (GIS)

After the data collection phase, ArcGIS software spatial interpolation methods are applied to characterize statistically the characteristics of the studied panel. We are particularly interested in the calculation of some classical statistical parameters such as the mean, the standard deviation, etc ... These statistical parameters are very determining as to the possibility of interpolation by Kriging (Kriging: denotes a set

of methods of interpolation based on mathematical and statistical models that all allow to estimate the value of a property at a point of geographical space from neighboring observations (the autocorrelation, being a function of distance).

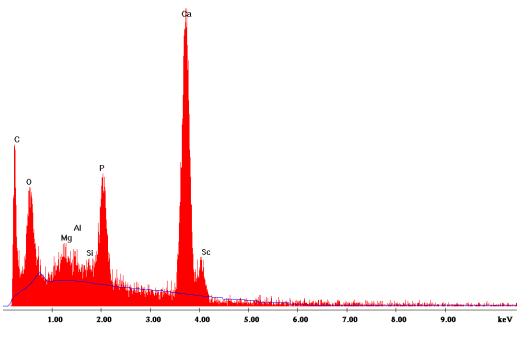


Figure 8: Spectrum of chemical elements of Spacer layer 4 (C4) – layer 3 (C3).

Subsequently, the realization of the cards is approached taking into account the objective assigned to the study and the expected results. It consists of:

- Making maps of the physical and chemical characteristics of the phosphate layers by the geostatistical method.
- The connection between the database and the GIS software through a data source that allows a dynamic link to be established so that updates of the data contained in the database are automatically taken into account in the mapping software.

3.6. Maps Creation

After the input database is developed from the well data, it is imported to the ArcMap interface of the ArcGIS geographic information system to be the subject of a geoprocessing aimed at creating iso-value maps of all the studied parameters (Figure 9).

3.7. 3D Model

The structural maps of the other layers, treated in the same way show that the general structure is practically unaffected by major structures. It is generally monoclinal with a slight dip to the northwest[17]. Some irregularities punctuate this structure, particularly in the south-west of the panel as shown by the 3D reconstitution of the panel (Figure 10). At our observation scale, no major structure of the type has been identified or folded. No disturbance has been observed either directly or indirectly in this panel.

3.8. Map listing mining boxes

Listing Cases is the mining database that links each box to its properties [18]. The establishment of the mining listing maps (Figure 11) is based on the compiled data of the physicochemical parameters obtained by the chemical analyzes of the samples taken in the phosphate series from the recognition wells. These data are then recorded on the lithological sections of these different wells for the purposes of correlation and resource evaluation.

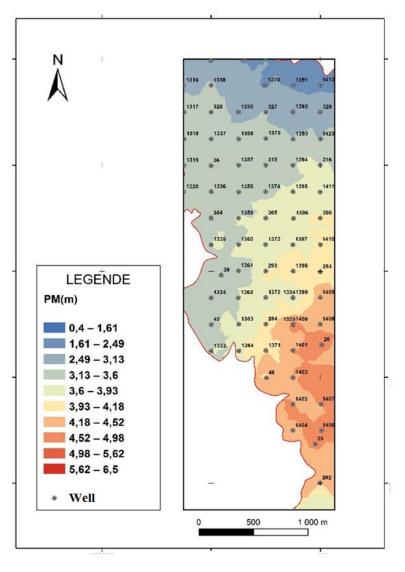


Figure 9: Example of the distribution map of global thickness.

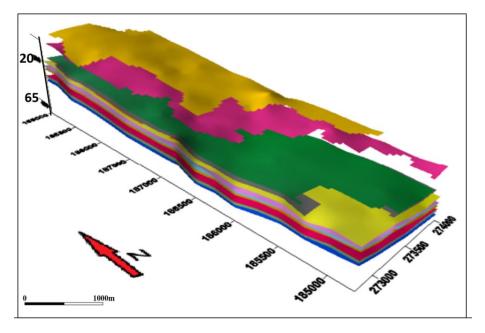


Figure 10: Presentation of phosphate layers in 3D model.

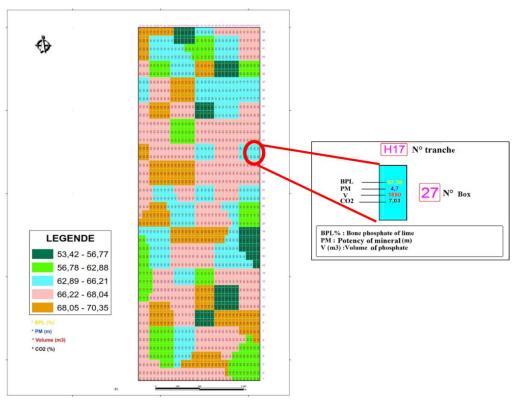


Figure 11: Example of a Map listing cases.

Conclusion

The work presented in this paper deals with a geological and mining study of the panel 7 of the exploration and exploitation zone of North Benguerir belonging to the Ganntour basin. The exploitation of phosphates, although it is done in open-pit, must be preceded by an exploration in order to reconstitute the geometry of the phosphate layers including their extension and their power; the evaluation of their mining quality based on their overall chemical composition and, particularly, their content of P2O5 (Phosphorus pentoxide). These data, in large numbers are obtained through direct or indirect methods based on random sampling or regularly spaced wells on exploration areas. They must be the object of an objective computerized treatment which allows to analyze and discuss the properties of the deposit and mainly to evaluate the volumes of the ore in such a way as to direct the miners towards the best exploitation methods.

In this work, a technique based on the use of ArcGIS software for the processing of data from panel 7 of the North Benguérir deposit is proposed. This technique allowed us to establish different maps showing different properties of the phosphate layers. The main ones are the BPL, CO2; Mg concentration maps, isopach maps, structural maps and finally listing boxes that link each box to all its mining properties. This enabled us reconstruct the geometry of the layers of the entire panel with an objective 3D view of the layers and to quantify the resources of this panel.

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