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**APPLICATION  
FOR QUANTIFYING THE LITHOLOGY  
OF A ROCK FORMATION  
USING GEOPHYSICAL WELL LOGS  
AND CONSTRAINED OPTIMIZATION METHODS**

## **1. INTRODUCTION**

Porosity and water saturation are the key reservoir parameters for hydrocarbon and water prospecting. To determine the mentioned parameters the set of equations that combine geological parameters with well logs is constructed [1, 3]. The goal of this paper is to present a comprehensive application for processing and interpreting these well logs.

## **2. PROBLEM**

Porosity and lithology are calculated from the set of equations describing model of a rock formation (1). These equations combine porosity and percentage volume of rock components with corresponding logs through the given geological factors and parameters of formation water. Water saturation is obtained from resistivity log (RT or RX0) [3].

$$P_i \pm \delta_i w_i = \sum_{j=1}^N (a_{ij} \times V_j) \quad (1)$$

where:

- $P_i$  – geophysical measurement,
- $\delta_i$  – measurement error,
- $w_i$  – weight of measurement,
- $a_{ij}$  – geological factor,
- $V_j$  – percentage volume of component (mineral, porosity, water saturation),
- $N$  – total number of components.

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Optimal solution to above-mentioned set of equations is obtained by minimizing the error function, defined as a difference between theoretical value and observed value [1, 3]. The objective function is given by formula (2):

$$f(\phi, V_j, S_w) = \sum_{i=1}^M \left( \frac{P_i^{theor} - P_i^{obs}}{P_i^{obs}} \right)^2 \quad (2)$$

where:

- $\phi$  – porosity,
- $V_j$  – percentage volume of mineral component,
- $S_w$  – water saturation,
- $P_i^{theor}$  – theoretical value,
- $P_i^{obs}$  – observed value,
- $M$  – total number of measurements.

Assumptions:

- 1) percentage volume of each component is between 0 and 1;
- 2) total percentage volume of mineral components and porosity is equal 1.

### 3. SOLUTION

Application for well log processing and interpretation was written in C# programming language, using integrated development environment Microsoft Visual Studio 2008 Professional and ALGLIB – a cross-platform numerical analysis and data processing library. For finding a proper solution to (2), *minbleic* subpackage was used – a part of ALGLIB library which supports boundary and linear equality/inequality constraints [4]. This subpackage allows to use BLEIC (Boundary, Linear Equality – Inequality Constraints) algorithm, based on nonlinear conjugate gradient method, to solve following optimization problems (3):

$$\begin{aligned} & \min f(x) \text{ s.t.} \\ & l_i \leq x \leq u_i, \quad l, x, u \in R^N \\ & Ax \circ b, \quad A \in R^{K \times N}, \text{ where } \circ \text{ is any combination of } \leq, =, \geq \end{aligned} \quad (3)$$

For every depth *minbleicoptimize* was performed using the optimal solution from previous depth as starting points. For the first depth starting points was obtained by the Monte Carlo method [2]. Gradient was calculated by ALGLIB through numerical differentiation provided by *minbleiccreatef* method.

## Optimization methods

Nonlinear conjugate gradient method (NCGM) is generalized form of the conjugate gradient method adapted to nonlinear optimization. In general, the NCGM is a deterministic optimization algorithm that finds the local minimum of a nonlinear function using its gradient. Monte Carlo method (MCM) is used for finding a global minimum of the objective function by random sampling. The MCM is a time-consuming method, therefore number of iteration was limited to  $10^5$  iterations.

### ALGLIB library – example of use

Minimize:

$$f(x, y) = (1 - x)^2 + 100(y - x^2)^2$$

subject to:

$$0 \leq x \leq 3$$

$$0 \leq y \leq 3$$

$$x + y = 4$$

Code:

```
class Program
{
    static void Main(string[] args)
    {
        gradient();
    }
    static void gradient()
    {
        double[] x = new double[] {0, 0}; //initial values x = x[0], y = x[1]
        double[,] c = {
            { 1, 0, 0}, //x[0] >= 0
            { 1, 0, 3}, //x[0] <= 3
            { 0, 1, 0}, //x[1] >= 0
            { 0, 1, 3}, //x[1] <= 3
            { 1, 1, 4} //x[0] + x[1] = 4
        };
        int[] ct = new int[] { 1, -1, 1, -1, 0 }; // >= <= >= <=
        alglib.minbleicstate state;
        alglib.minbleicreport rep;
```

```

//stopping conditions for the optimizer
double epsg = 0;
double epsf = 0;
double epsx = 0;
int maxits = 0;

//differentiation step
double diffstep = 1.0e-6;

alplib.minbleiccreatef(x, diffstep, out state); //creating optimizer
alplib.minbleicsetlc(state, c, ct); //adding linear constraints
alplib.minbleicsetcond(state, epsg, epsf, epsx, maxits); //tuning stopping conditions
alplib.minbleicoptimize(state, rosenrock, null, null); //optimization
alplib.minbleicresults(state, out x, out rep); //obtaining results

//printing values of x[0] and x[1]
Console.WriteLine("x = " + Math.Round(x[0], 3).ToString());
Console.WriteLine("y = " + Math.Round(x[1], 3).ToString());

//printing the value of f(x[0], x[1])
double rval = 0;
object o = new object();
rosenrock(x, ref rval, o);
Console.WriteLine("f(x, y) = " + Math.Round(rval, 3).ToString());

Console.ReadKey(); }

static void rosenrock(double[] x, ref double func, object obj)
{
    func = (1 - x[0]) * (1 - x[0]) + 100 * (x[1] - x[0] * x[0]) * (x[1] - x[0] * x[0]);
}
}

```

Results:

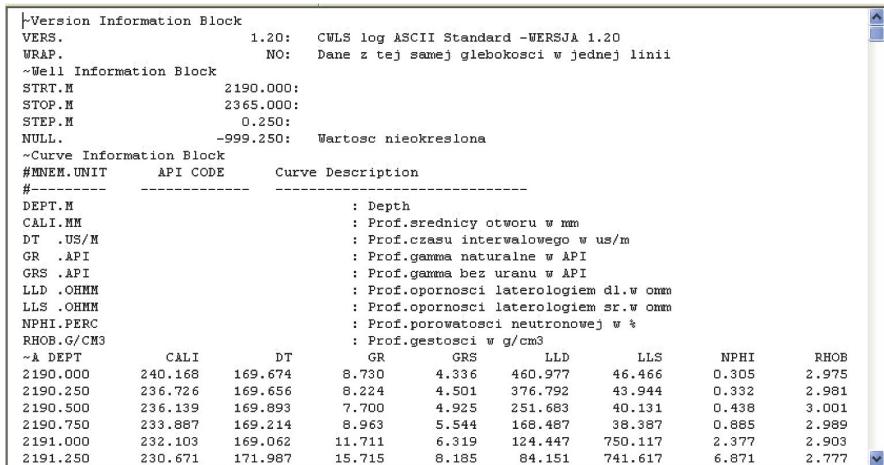
x = 1,561  
y = 2,439  
f(x, y) = 0,315

### **Program description**

For calculations the following set of logs was given:

- interval transit time DT [us/m],
- gamma radiation GR [API],
- neutron porosity NPHI [% in limestone scale],
- bulk density RHOB [g/cm<sup>3</sup>],
- resistivity of the uninvaded formation RT [ $\Omega\cdot m$ ],
- resistivity of the invaded zone RX0 [ $\Omega\cdot m$ ].

The expected input was – beside parameters defined by user – the file with logs encrypted in LAS format (Fig. 1).



```

Version Information Block
VERS.          1.20:  CWLS log ASCII Standard -WERSJA 1.20
WRAP.          NO:   Dane z tej samej glebokosci w jednej linii
~Well Information Block
STRT.M        2190.000:
STOP.M        2365.000:
STEP.M        0.250:
NULL.         -999.250:  Wartosc nieokreslona
~Curve Information Block
#MNEM.UNIT    API CODE      Curve Description
#----- -----
DEPT.M          : Depth
CALI.MM         : Prof.srednicy otworu w mm
DT .US/M        : Prof.czasu interwalowego w us/m
GR .API         : Prof.gamma naturalne w API
GRS .API        : Prof.gamma bez uranu w API
LLD .OHMM       : Prof.opornosci laterologiem dl.w cm
LLS .OHMM       : Prof.opornosci laterologiem sr.w cm
NPHI .PERC      : Prof.porowatosc neutronowej w %
RHOB .G/CM3     : Prof.gestosc w g/cm3
#& DEPT       CALI       DT      GR      GRS      LLD      LLS      NPHI      RHOB
2190.000      240.168   169.674  8.730   4.336   460.977  46.466  0.305   2.975
2190.250      236.726   169.656  8.224   4.501   376.792  43.944  0.332   2.981
2190.500      236.139   169.893  7.700   4.925   251.683  40.131  0.438   3.001
2190.750      233.887   169.214  8.963   5.544   168.487  38.387  0.885   2.989
2191.000      232.103   169.062  11.711  6.319   124.447  750.117  2.377   2.903
2191.250      230.671   171.987  15.715  8.185   84.151   741.617  6.871   2.777

```

**Fig. 1.** Example of input file in LAS format showed in external text editor

### LAS format description

LAS (*Log ASCII Standard*) is a standard file-format to store well log information [5]. Each LAS file contains numerous sections and each section begins with a tilde (~) mark. The sections that make up an LAS file are as follows:

- ~V – version and wrap mode information [mandatory],
- ~W – well identification [mandatory],
- ~C – curve information [mandatory],
- ~P – parameters or constants [optional],
- ~O – other information such as comments [optional],
- ~A – ASCII log data [mandatory].

Well Information section defines the following mnemonics:

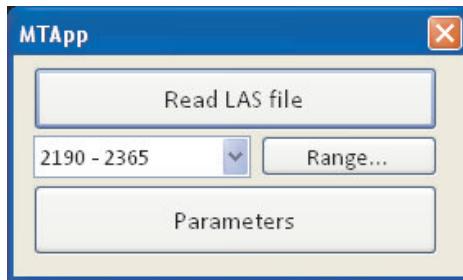
- STRT.M – the first depth in the file;
- STOP.M – the last depth in the file;
- STEP.M – the depth increment used;
- NULL. – the null value.

The hash sign (#) is used to indicate that the line is a comment line.

### Application in use

The main window (Fig. 2) allows to:

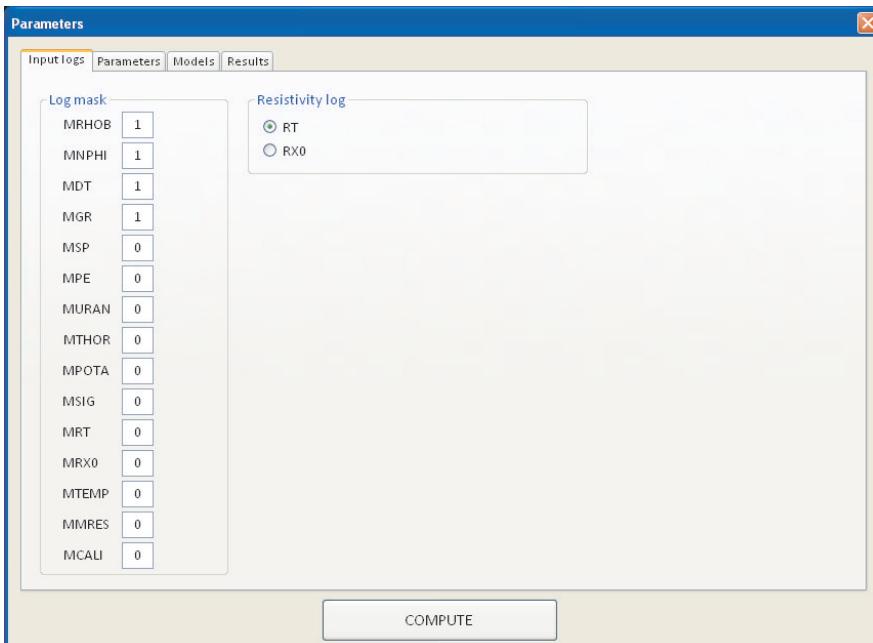
- read from LAS file;
- define depth interval used in interpretation;
- activate *Parameters* window.



**Fig. 2.** Main window

The *Parameters* window (Fig. 3–6) allows to:

- mask logs i.e. to choose logs participating in interpretation;
- select resistivity log for optimization (RT/RX0);
- define: general parameters, mud parameters, mud filtrate parameters, formation water parameters, hydrocarbons parameters, loam parameters;
- select the model parameters;
- mask minerals i.e. to choose minerals participating in interpretation;
- define mineral properties (Fig. 7);
- set output parameters;
- start calculations.



**Fig. 3.** Input logs tab in the *Parameters* window

**Parameters**

Input logs Parameters Models Results

<b>General</b>	<b>Mud</b>	<b>Mud filtrate</b>	<b>Formation water</b>	<b>Hydrocarbons</b>				
TD BHT T0 GG CFF EXPIM EXPSW EXPSX0 TSN	MUTD RM TRM RMC TRMC RHOM PEM URM THM POM GRM SIGM	iflawa 0.001 0.000 0.001 0.000 1.000 0.80000 0.00150 0.00350 0.01500 0.00100 0.00500	RMF TRMF RHOMF DTMF PEMF URMF THMF POMF GRMF SIGMF	0.001 0.000 1.000 650.000 0.8000 0.00150 0.00350 0.01500 0.00100 0.00500	RW TRW RBW RHOW RHOBW DTW PEW URW THW POW GRW SIGW	0.001 0.000 0.000 1.200 1.200 650.000 0.80000 0.01500 0.00350 0.01500 0.00100 0.00500	RH TRH RHOH RHOHR DTH PEH URH GRH SIGH	0.001 0.000 0.800 0.800 769.000 0.80000 0.00150 0.00100 0.00500
					<b>Loam</b>			
					QV PHICL RHODCL	100.000 0.500 2.900		

COMPUTE

**Fig. 4.** Parameters tab in the *Parameters* window

**Parameters**

Input logs Parameters Models Results

<b>Model parameters</b>	<b>Mineral mask</b>	<b>Mineral parametres</b>
DTEQ CLEQ SATEQ PEREQ PHIMN PHIMX SWMN SWMX SWIRMN SWIRMX SX0MN SX0MX	LOAM SANDSTONE LIMESTONE DOLOMITE MIN1 MIN2 MIN3 MIN4 MIN5 MIN6	LOAM SANDSTONE LIMESTONE DOLOMITE MIN1... MIN2... MIN3... MIN4... MIN5... MIN6...

COMPUTE

**Fig. 5.** Models tab in the *Parameters* window

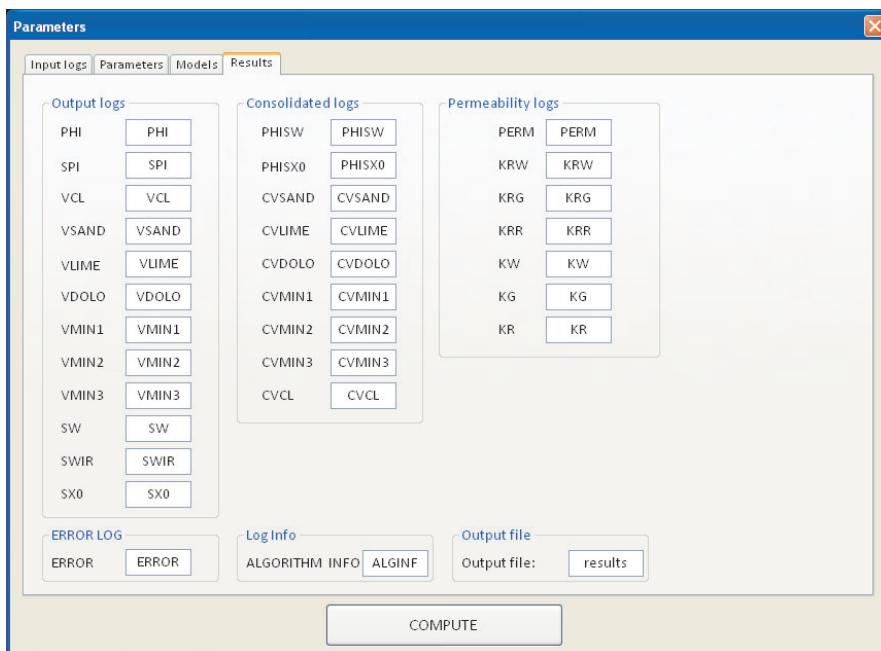


Fig. 6. Results tab in the *Parameters* window

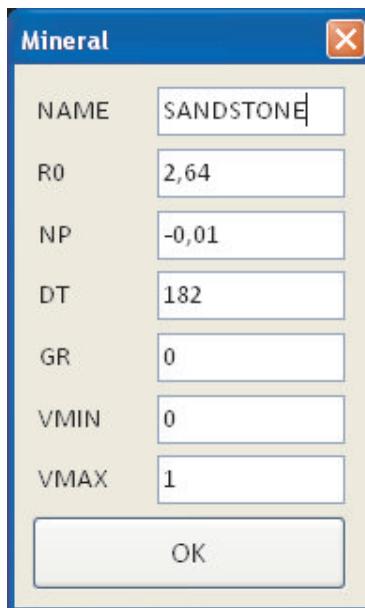
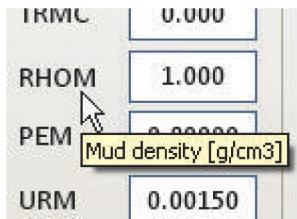


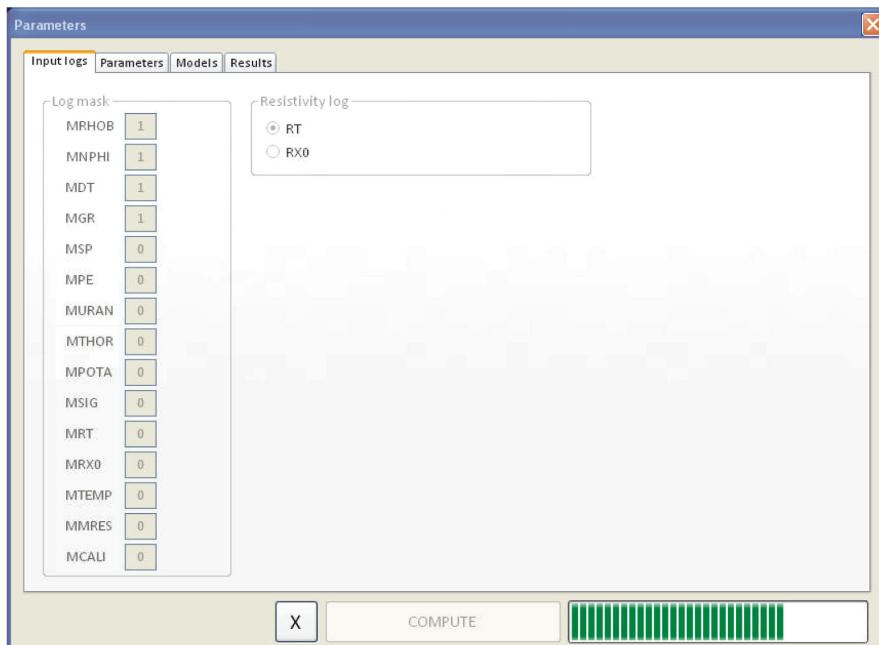
Fig. 7. Mineral window

Every component contained in *Parameters* window can provide short information (hint) when the user places the pointer on its name. This mechanism is known as *tooltip* (Fig. 8).



**Fig. 8.** Tooltip example

After clicking the *COMPUTE* button a progress bar (Fig. 9) which visualize how many of depths was already computed became visible. After completion of calculations the *Results* window is shown.



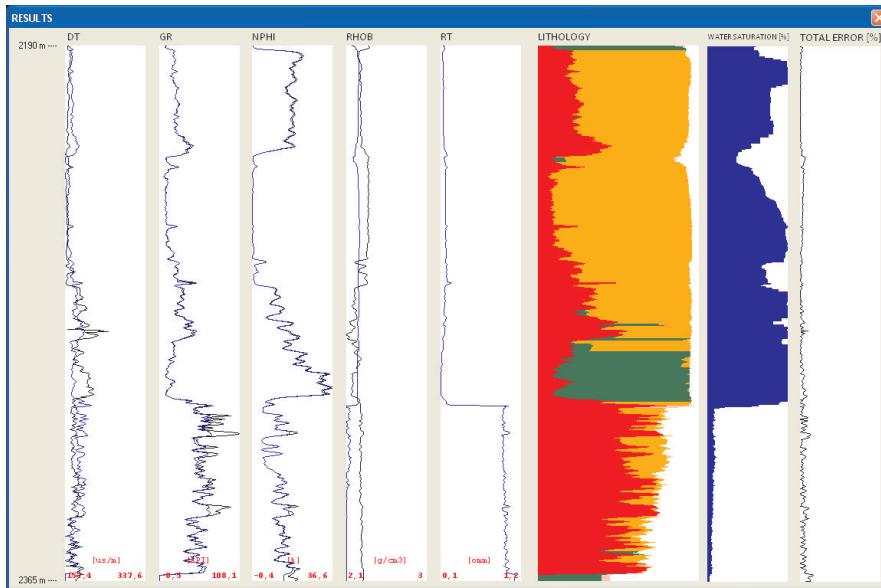
**Fig. 9.** Parameters window during computation

The *Results* window (Fig. 10) consists of:

- graphs of logs participating in interpretation (both experimental and theoretical – Fig. 12),
- RT/RX0 graph,

- lithology graph,
- water saturation graph,
- total error graph.

When user clicked on the lithology graph the *Legend* window appears (Fig. 11).

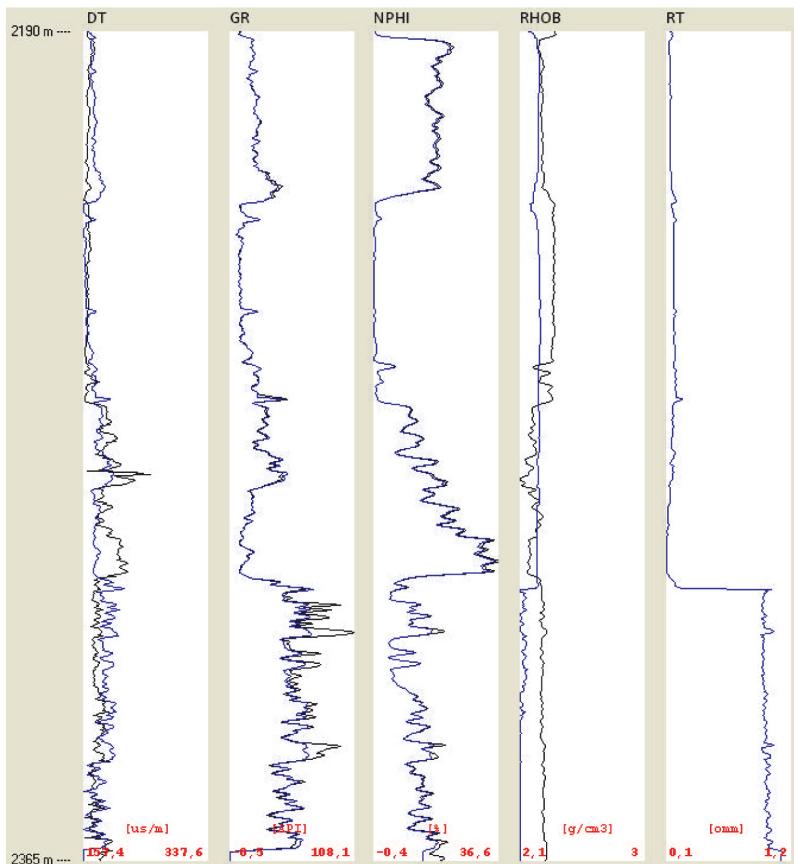


**Fig. 10.** Results window – from the left: DT, GR, NPHI, RHOB, RT, LITHOLOGY, WATER SATURATION [%], TOTAL ERROR [%]



**Fig. 11.** Legend window

Every component contained in the *Results* window can provide basic information such as log name and the range of taken values via *tooltip*. Precise values of percentage volume of selected minerals, pore space and water saturation are available in automatically generated LAS file, placed in folder containing input logs, under the name given by user (Fig. 13).



**Fig. 12.** Experimental logs (black curves) with theoretical logs (blue curves)

~Version Information Block						
VERS.	1.20: CWLS log ASCII Standard -WERSJA 1.20					
WRAP.	NO: Dane z tej samej glebokosci w jednej linii					
~Well Information Block						
STRT.M	2190.000:					
STOP.M	2365.000:					
STEP.M	0.250:					
NULL.	-999.250: Wartosc nieokreslona					
~Curve Information Block						
DEPT.M	: Depth					
PORO.PERC	: Pore space [%]					
LOAM.PERC	: Loam [%]					
SAND.PERC	: Sandstone [%]					
LIME.PERC	: Limestone [%]					
DOLO.PERC	: Dolomite [%]					
WATE.PERC	: Water saturation [%]					
~A DEPT						
	PORO	LOAM	SAND	LIME	DOLO	WATE
2190,00	18,971	9,457	68,866	0,000	2,706	26,808
2190,25	18,405	9,341	72,253	0,000	0,000	26,463
2190,50	17,245	8,639	41,556	32,560	0,000	27,717
2190,75	15,960	10,330	29,324	44,386	0,000	29,523
2191,00	12,654	13,226	0,000	74,121	0,000	36,248
2191,25	9,606	17,801	0,000	72,592	0,000	48,477
2191,50	7,516	21,559	0,000	70,926	0,000	63,991
2191,75	7,733	22,756	0,000	69,511	0,000	64,336
2192,00	7,920	21,186	0,707	70,187	0,000	64,396
2192,25	8,002	20,455	1,015	70,527	0,000	64,524

**Fig. 13.** Results of computation in LAS format showed in external text editor

#### **4. SUMMARY**

The purpose of this paper was to present a comprehensive application for determining the lithology, porosity and water saturation from well logs. Presented application makes it possible to compare experimental logs with computed logs as well as to visualize a total error, what significantly simplify verification of results correctness, allowing to enhance models that was used in interpretation. Moreover, output data encoded in LAS format allow to perform further processing of obtained results in another programs.

#### **REFERENCES**

- [1] Jarzyna J. (ed.), Bała M., Cichy A., Gądek W., Gąsior I., Karczewski J., Marzencki K., Stadtmüller M., Twaróg W., Zorski T.: *Przetwarzanie i interpretacja profilowań geofizyki wiertniczej – system GeoWin*. ARBOR, Kraków, 2002.
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