

1 Application of fractal models to delineate mineralized zones in
2 the Pulang porphyry copper deposit, Yunnan, Southwest China

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9 Abstract

10 The aim of this study is to delineate and identify various mineralized zones and
11 barren host rocks based on surface and subsurface lithogeochemical data from the
12 Pulang porphyry copper deposit, Southwest China, utilizing the number-size (N-S),
13 concentration-volume (C-V) and power spectrum-volume (S-V) fractal models. The
14 N-S model reveals three mineralized zones characterized by Cu thresholds of 0.28%
15 and 1.45%: <0.28% Cu represents weakly mineralized zones and barren host rocks,
16 0.28%-1.45% Cu represents moderately mineralized zones, and >1.45% Cu represents
17 highly mineralized zones. The results obtained by the C-V model depict four
18 geochemical zones defined by Cu thresholds of 0.25%, 1.48% and 1.88%,
19 representing nonmineralized wall rocks (Cu<0.25%), weakly mineralized zones
20 (0.25%-1.48%), moderately mineralized zones (1.48%-1.88%), and highly
21 mineralized zones (Cu>1.88%). The S-V model is used by performing a 3D fast
22 Fourier transformation of assay data in the frequency domain. The S-V model reveals
23 three mineralized zones characterized by Cu thresholds of 0.23% and 1.33%: <0.23%
24 Cu represents leached zones and barren host rocks, 0.23%-1.33% Cu represents
25 hypogene zones, and >1.33% Cu represents supergene enrichment zones. All the
26 multifractal models indicate that high-grade mineralization occurs in the central and
27 southern parts of the ore deposit. Their results are compared with the alteration and

28 mineralogical models resulting from the 3D geological model using a logratio matrix.
29 The results show that the S-V model is best at identifying highly mineralized zones in
30 the deposit. However, the results of the C-V model for moderately and weakly
31 mineralized zones are more accurate than those obtained from the N-S and S-V
32 models.

33 Keywords: Fractal; Concentration-volume (C-V) model; Number-size (N-S) model;
34 Power spectrum-volume (S-V) model; Mineralized zone; the Pulang porphyry copper
35 deposit

36 **1. Introduction**

37 The definition and delineation of different mineralized zones and non-mineralized
38 wall rocks are the main goal in economic geology and mineral exploration.
39 Investigation of ore mineralogy and paragenetic sequence provides useful data on
40 ore-forming processes in deposits because typical characteristics of various types of
41 ore deposits are reflected by their mineral assemblages (Craig and Vaughan, 1994;
42 White and Hedenquist, 1995). Common methods generally use mineralography,
43 petrography and alteration mineral assemblage analysis to delineate various
44 mineralized zones in porphyry deposits (Beane, 1982; Schwartz, 1947; Sillitoe, 1997;
45 Berger et al., 2008). Lowell (1968) first proposed a conceptual model of the lateral
46 and vertical variations in mineralogy within alteration zones. Some similar models
47 were developed for potassic alteration, which is usually situated in the center and deep
48 parts of porphyry ore deposits, based on this conceptual model (Sillitoe and Gappe,
49 1984; Cox and Singer, 1986; Melfos et al., 2002). Fluid inclusion and stable isotope
50 studies are other methods used to outline different mineralization phases based on
51 thermometric and isotope element parameters and other geological parameters (e.g.,
52 Boyce et al., 2007; Faure et al., 2002; Wilson et al., 2007). Drillhole data and logging
53 information, including mineralogical information, host rock changes and
54 alterations are helpful in delineating mineralization zones. Different geological
55 interpretations could be used to detect zone boundaries, which may also lead to
56 different results because the elemental grade distribution may not be taken into

57 consideration.

58 Non-Euclidean fractal geometry (Mandelbrot, 1983) is an important branch of
59 nonlinear mathematical sciences and has been applied in various research fields of the
60 geosciences since the 1980s. The relationships between geology, geochemistry and
61 mineralogical settings and spatial information can be researched by methods based on
62 fractal geometry (Afzal et al., 2011; Carranza, 2008, 2009). Bolviken et al. (1992)
63 and Cheng et al. (1994) have shown that geochemical patterns of various elements
64 have fractal dimensions. The concentration-area (C-A) model was proposed by
65 Cheng et al. (1994) to recognize geochemical anomalies from background
66 concentrations and calculate elemental thresholds of different geochemical data.
67 Furthermore, many other fractal models have been proposed and applied in
68 geochemical exploration work, including the number-size (N-S) fractal model
69 proposed by Mandelbrot (1983) and Agterberg (1995), the power spectrum-area (S-A)
70 fractal model proposed by Cheng et al. (1999), the concentration-distance (C-D)
71 fractal model proposed by Li et al. (2003), the concentration-volume (C-V) fractal
72 model proposed by Afzal et al. (2011) and the power spectrum-volume (S-V) fractal
73 model proposed by Afzal et al. (2012).

74 Methods of fractal analysis also illustrate the relationships between geological,
75 geochemical and mineralogical settings and spatial information derived from the
76 analysis of mineral deposit occurrence data (Carranza, 2008; Carranza et al., 2009;
77 Goncalves et al., 2001). Various geochemical processes can be described based on the
78 differences in fractal dimensions obtained from the analysis of relevant geochemical
79 data. Afzal et al. (2011) considered that the log-log plots obtained by fractal methods
80 are useful tools to delineate different geological populations of geochemical data, and
81 the thresholds could be determined as some breakpoints in those plots.

82 The application of fractal models to delineate various grade mineralization zones
83 was dependent on the relationships between the metal grades and volumes (Afzal et
84 al., 2011; Agterberg et al., 1993; Cheng, 2007; Sim et al., 1999; Turcotte, 1986). Afzal
85 et al. (2011 and 2012) proposed a concentration-volume (C-V) and power
86 spectrum-volume (S-V) fractal model to delineate different porphyry-Cu mineralized

87 zones and barren host rocks. In this paper, N-S, C-V and S-V fractal models were
88 applied to delineate various mineralized zones and barren host rocks in the Pulang
89 porphyry copper deposit, Yunnan, Southwest China.

90 **2. Fractal models**

91 2.1. Number-size (N-S) fractal model

92 The number-size (N-S) method proposed by Mandelbrot (1983) can be utilized
93 to describe the distribution of geochemical populations (Sadeghi et al., 2012). In this
94 method, geochemical data does not undergo any preprocessing (Mao et al., 2004).
95 This model shows a relationship between desirable attributes (e.g., Cu concentration
96 in this study) and their cumulative number of samples (Sadeghi et al., 2012). A
97 power-law frequency model has been proposed to explain the N-S relationship
98 according to the frequency distribution of elemental concentrations and cumulative
99 number of samples with those attributes (e.g., Li et al., 1994; Sadeghi et al., 2012;
100 Sanderson et al., 1994; Shi and Wang, 1998; Turcotte, 1996; Zuo et al., 2009a).

101 The N-S model proposed by Mandelbrot (1983) can be expressed as follows:

$$102 \quad N(\geq\rho)=F\rho^{-D} \quad (1)$$

103 where ρ denotes the element concentration, $N(\geq\rho)$ denotes the cumulative number of
104 samples with concentrations greater than or equal to ρ , F is a constant and D is the
105 scaling exponent or fractal dimension of the distribution of element concentrations.
106 According to Mandelbrot (1983), log-log plots of $N(\geq\rho)$ versus ρ show linear
107 segments with different slopes $-D$ corresponding to different concentration intervals.

108 2.2. Concentration-volume (C-V) fractal model

109 Afzal et al. (2011) proposed a concentration-volume (C-V) fractal model based
110 on the same principle of the concentration-area (C-A) model (Cheng et al., 1994) to
111 analyze the relationship between the concentration of ore elements and accumulative
112 volume with concentrations greater than or equal to a given value (Afzal et al., 2011;
113 Zuo et al., 2016; Lin et al., 2013; Sadeghi et al., 2012; Soltani et al., 2014; Sun and
114 Liu, 2014; Wang, G. et al., 2012). This model can be expressed as follows:

115
$$V(\rho \leq v) \propto \rho^{-a_1}; V(\rho \geq v) \propto \rho^{-a_2} \quad (2)$$

116 $V(\rho \geq v)$ and $V(\rho \leq v)$ represent the occupied volumes with concentrations above or
117 equal to and less than or equal to the contour value v ; v indicates the threshold value
118 of a zone; and a_1 and a_2 are the characteristic indexes. The thresholds obtained by this
119 method indicate the boundaries between the different grade mineralization zones and
120 barren host rocks of ore deposits. The drillhole data of the elemental concentrations
121 were interpolated by using geostatistical estimation to compute $V(\rho \geq v)$ and $V(\rho \leq v)$,
122 which are the volume values enclosed by a contour level ρ in a 3D model.

123 2.3. Power spectrum-volume (S-V) fractal model

124 Different geochemical patterns in the spatial domain could be seen as layered
125 signals of various frequencies. Cheng et al. (1999) proposed the power spectrum-area
126 (S-A) fractal model to recognize geochemical anomalies from backgrounds utilizing
127 the method of spectrum analysis in the frequency domain according to this argument.
128 This model is combined with a concentration-area (C-A) model (Cheng et al. 1994),
129 offering a useful tool to determine an optimum threshold value between various
130 patterns based on the scaling property.

131 Afzal et al. (2012) proposed the power spectrum-volume (S-V) fractal model to
132 delineate different grade mineralization zones based on the same principle as the S-A
133 model proposed by Cheng et al. (1999). The S-V model was utilized in the frequency
134 domain by applying a fast Fourier transformation to the assay data. The straight lines
135 obtained by log-log plotting indicate the relationships between the power spectra and
136 relevant volumes of ore elements. These relationships were utilized to recognize the
137 hypogene zones and supergene enrichment zones from the barren host rocks and the
138 leached zone of the deposit. The recognition of various mineralization zones is based
139 on the power-law relationships between the power spectra and occupied volumes. The
140 formula is as follows:

141
$$V(\geq S) \propto S^{-2/\beta} \quad (3)$$

142 where the power-law relationships between the power spectra ($S = \|F(W_x, W_y,$
143 $W_z)\|$) and occupied volumes with power spectra greater than or equal to S can be

144 indicated by this form; F represents the fast Fourier transformation of the
145 measurement $\mu(x, y, z)$; and W_x , W_y and W_z indicate wave numbers or angular
146 frequencies in the X, Y and Z directions in a 3D model. The range of index β is $0 < \beta \leq 2$
147 or $1 \leq 2/\beta$ with the special cases of $\beta=2$ and $2/\beta=1$ corresponding to nonfractal and
148 monofractal expressions, and $1 < 2/\beta$ corresponding to multifractals (Cheng, 2006).

149 By using the method of geostatistical estimation, the drillhole data of elemental
150 concentration values were interpolated to construct a block model of ore element
151 distribution. The power spectrum values can be obtained by 3D fast Fourier
152 transformation of the ore element grades. The logarithm of all the power spectrum
153 values and accumulative volume values were calculated. Additionally, the log-log plot
154 between power spectrum and volume was drawn according to previously determined
155 values. Then, the filters were constructed on the basis of threshold values obtained by
156 the log-log plot of S-V. Finally, the power spectra were converted back to the space
157 domain by utilizing inverse fast Fourier transformation.

158 **3. Geological setting of the Pulang porphyry copper deposit**

159 The Pulang porphyry copper deposit is situated in the southern end of the Yidun
160 continental arc, Southwest China (Fig. 1). The continental arc was produced due to the
161 westward subduction of Garze–Litang oceanic crust (Deng et al., 2014b, 2015; Wang
162 et al., 2014). The Pulang ore deposit, one of the largest porphyry copper deposits in
163 China (Deng et al., 2012, 2014a; Mao et al., 2012, 2014), is characterized by a typical
164 porphyry-type alteration zone. The geological characteristics of the deposit, including
165 the alteration types and their zonation, the geometry of the orebody, the metallogenic
166 time and the geodynamic settings, have been systematically researched (Leng et al.,
167 2012; Li et al., 2011, 2013). The deposit consists of five ore-bearing porphyry bodies
168 covering an area of approximately 9 km², and the explored ore tonnage of Cu is
169 estimated to be 6.50 Mt (Liu et al., 2013).

170 The outcrop strata of the Pulang deposit are dominated by Upper Triassic
171 Tumugou Formation clastic rocks and andesite and Quaternary sediments (Fig. 1c).
172 The Triassic porphyry intrusions primarily comprise quartz diorite porphyry, quartz

173 monzonite porphyry, quartz diorite porphyrite and granodiorite porphyry. The
174 Tumugou Formation strata were intruded by the quartz diorite porphyry with an age
175 of 219.6 ± 3.5 Ma (zircon U-Pb dating) (Pang et al., 2009). Then, quartz monzonite
176 porphyry with an age of 212.8 ± 1.9 Ma and granodiorite porphyry with an age of
177 206.3 ± 0.7 Ma (zircon U-Pb dating) (Liu et al., 2013) crosscut the quartz diorite
178 porphyry. The quartz monzonite porphyry is considered to be associated with
179 mineralization because its age is similar to the molybdenite Re-Os isochron age of
180 213 ± 3.8 Ma from the orebody (Zeng et al., 2004). Moreover, the Cu concentrations
181 of the quartz monzonite porphyry are higher than those of the other porphyries.

182 The porphyry-type alteration zones transition from early potassium-silicate
183 alteration through quartz-sericite alteration to propylitization, upward and outward
184 from the core of the quartz monzonite porphyry (Fig. 4). The wall rocks near the
185 porphyries were mostly changed into hornfels. Systematic drilling has demonstrated
186 that the potassium-silicate and quartz-sericite zones host the main orebodies,
187 constituting the core of mineralized zones. The propylitic zones and hornfels only
188 develop weak mineralization. The orebodies occur mainly in potassium-silicate and
189 quartz-sericite and occur as veins in the propylitic zones and hornfels. The major rock
190 types in the deposit are quartz monzonite porphyry, quartz diorite porphyrite, granite
191 diorite porphyry, quartz diorite porphyry and hornfels (Fig. 2). Metallic minerals
192 mainly include pyrite, chalcopyrite with a small amount of molybdenite and pyrrhotite
193 (Fig. 3).

194 **4. Fractal modeling**

195 Based on the geological data (which include the collar coordinates of each
196 drillhole, azimuth and dip (orientation), lithology and mineralogy) recorded from 130
197 drillholes in the Pulang deposit, 20492 lithogeochemical samples were collected at 2
198 m intervals. The laboratory of the 3rd Geological Team of the Yunnan Bureau of
199 Geology and Mineral Resources utilized the iodine-fluorine and oscillo polarographic
200 method to analyze the concentrations of Cu and associated paragenetic elements, and
201 its analytical uncertainty is less than 7% (Yunnan Diqing Nonferrous Metal Co. Ltd.,

202 2009). Only Cu concentrations were studied in this study. The histogram and Q-Q plot
203 of the log-transformed Cu data indicate that the distribution of Cu data is log-normal
204 (Fig. 5). The experimental semivariogram of the Cu data of the Pulang deposit
205 indicates a range and nugget effect of 320.0 m and 0.25, respectively (Fig. 6). The
206 spherical model is fitted with regard to the experimental semivariogram. The 3D
207 model of the Cu concentration distribution of the Pulang deposit was produced with
208 the ordinary kriging method using Geovia Surpac software on the basis of the
209 semivariogram and anisotropic ellipsoid. Fundamentally, the accuracy of the
210 interpolation results mainly depends on whether the interpolation model accurately
211 fits the spatial distribution characteristics of the deposit. Ordinary kriging was used
212 because it is compatible with a stationary model; it only requires a variogram, and it is
213 the most commonly used form of kriging (Chilès and Delfiner, 1999). Goovaerts
214 (1997) showed that the values in unsampled locations are estimated by the ordinary
215 kriging method according to the moving average of the interest variables, satisfying
216 various distribution forms of data. Ordinary kriging is a spatial estimation method in
217 which the error variance is minimized. This error variance is based on the
218 configuration of the data and its variogram (Yamamoto, 2005). The correct variogram
219 in kriging interpolation can guarantee the accuracy of the interpolation results.

220 The accuracy of the spatial interpolation analysis is verified by comparing the
221 difference between the measured values and the predicted values to select the best
222 variogram model. To test the variogram model, the cross-validation method was used
223 to determine whether the parameters of the variogram model were correct. The
224 distribution of the residual is normal (Fig. 7), and the mean error between the actual
225 and estimated Cu grades is equal to 0 (Table 1). This result indicates that this model is
226 reasonable and that the variogram parameters used for estimating the Cu grade are
227 unbiased.

228 The obtained block models were used as inputs to the fractal models. The Pulang
229 deposit was modeled by 20 m×20 m×5 m voxels, and they were decided by the grid
230 drilling dimensions and geometrical properties of the deposit (David, 1970). The
231 Pulang deposit is totally modeled with 150973 voxels. The terms “highly”,

232 “moderately” and “weakly” have been used to classify mineralized zones based on
233 fractal modeling, in accordance with the classification of the ore grades in the deposit.

234 **4.1. Number-size (N-S) fractal modeling**

235 The N-S model was applied to the Cu data (Fig. 8). The selection of breakpoints
236 as threshold values is an objective decision because geochemical populations are
237 defined by different line segments in the N-S log-log plot. The straight fitted lines
238 were obtained based on least-square regression (Agterberg et al., 1996; Spalla et al.,
239 2010). In other words, the intensity of element enrichment is depicted by each slope
240 of the line segments in the N-S log-log plots (Afzal et al., 2010; Bai et al., 2010).

241 Based on the classification of the 3D model of Cu data and the thresholds obtained
242 from the N-S fractal model (Table 2), highly mineralized zones are situated in the
243 southern and central parts of the Pulang deposit and coincide with the
244 potassium-silicate alterations. However, small and highly mineralized zones are
245 located in the central parts of the Pulang deposit (Fig. 9). Moderately mineralized
246 zones occur along a northwest-southeast trend and correlate with the phyllic zones.
247 Weakly mineralized zones and barren host rocks are situated in the marginal parts of
248 the area.

249 **4.2. Concentration-volume (C-V) fractal modeling**

250 The occupied volumes corresponding to the Cu grades were computed to obtain
251 the concentration-volume model according to the 3D model of the Pulang deposit.
252 Through the obtained C-V log-log plot, the threshold values of the Cu grades were
253 determined (Fig. 10). These results indicate the power-law relationship between Cu
254 grade and volume. According to these results (Table 3), the low-concentration zones
255 exist in many parts of the deposit and occur along a northwest-southeast trend.
256 Moderately and highly mineralized zones are situated in several parts of the central
257 deposit and to the south of the deposit (Fig. 11).

258 **4.3. Power spectrum-volume (S-V) fractal modeling**

259 Based on the geological data (which include the collar coordinates of each
260 drillhole, azimuth and dip (orientation), lithology and mineralogy) recorded from 130

261 drillholes in the deposit, a 3D model and block model of the distribution of Cu in the
262 Pulang deposit were constructed with ordinary kriging using Geovia Surpac software.

263 The power spectrum (S) was calculated for the 3D elemental distribution using
264 3D fast Fourier transformation in MATLAB (R2016a). The logarithmic values of the
265 power spectra and relevant volumes were plotted against each other (Fig. 12). The
266 straight lines fitted in the log-log plot indicate different relationships between the
267 power spectra and occupied volumes. The thresholds of $\log S = 7.81$ and $\log S = 8.70$
268 were determined by the log-log S-V plot. The 3D filters were designed to separate
269 different mineralization zones on the basis of these threshold values. Inverse fast
270 Fourier transformation was used to convert the decomposed components back into the
271 space domain by using MATLAB (R2016a). According to the results, the Cu
272 concentrations of the hypogene zones range from 0.23% to 1.33% (Table 4), and
273 values of $>1.33\%$ Cu correspond to the supergene enrichment zones, whereas values
274 of $<0.23\%$ Cu correspond to the leached zone and barren host rocks (Fig. 13).

275 **5. Comparison of the fractal models and geological model of the** 276 **deposit**

277 Alteration models have a key role in zone delineation and in presenting
278 geological models, as described by Lowell and Guilbert (1970). The potassic and
279 phyllic alterations control major mineralization within supergene enrichment and
280 hypogene zones according to these models. Models of Cu mineralization zones
281 derived via fractal models can be compared with geological data to validate the results
282 of analysis in different porphyry Cu deposits. The results of the fractal modeling of
283 the Pulang deposit were compared with the 3D geological model of the deposit
284 constructed by using Geovia Surpac and drillhole data (Fig. 2). Moreover, the results
285 obtained from these fractal models were controlled by mineralogical investigations.

286 Carranza (2011) has illustrated an analysis for the calculation of spatial
287 correlations between two binary datasets, especially mathematical and geological
288 models. An intersection operation between the mineralization zones obtained from
289 fractal models and the different alteration zones in the geological model was

290 performed to derive the amount of voxels corresponding to each of the classes of
291 overlap zones (Table 5). Using the obtained numbers of voxels, the Type I error (T1E),
292 Type II error (T2E), and overall accuracy (OA) of the fractal model were estimated
293 with respect to different alteration zones and the geological data (Carranza, 2011).
294 The OAs of the fractal models of the mineralized zones were compared as follows.

295 A comparison between highly mineralized zones based on the fractal models and
296 potassic alteration zones resulting from the 3D geological model shows that there is a
297 similarity among these fractal models. The overall accuracies for the C-V, N-S and
298 S-V models are 0.50, 0.51 and 0.52, respectively (Table 6), which indicate that the
299 S-V model gives better results for identifying highly mineralized zones in the deposit.
300 The number of overlapped voxels (A) in the S-V model is higher than those in the
301 N-S and C-V models. The correlation (from OA results) between highly mineralized
302 zones obtained from S-V modeling and potassic alteration zones is better than that of
303 the N-S and C-V model because of a strong proportional relationship between the
304 extension and positions of voxels in the S-V model and the potassic alteration zones
305 in the 3D geological model.

306 A comparison between phyllic alteration zones resulting from the 3D geological
307 model and moderately and weakly mineralized zones from the fractal modeling shows
308 that the overall accuracies of the C-V, N-S and S-V fractal models with respect to
309 phyllic alteration zones of the geological model are 0.59, 0.54 and 0.56, respectively.
310 The overall accuracy of moderately and weakly mineralized zones obtained from C-V
311 modeling is higher than that of mineralized zones obtained from N-S and S-V
312 modeling (Table 7). On the other hand, moderately mineralized zones defined by C-V
313 modeling overlap with phyllic zones defined by the 3D geological model. However,
314 the results of the C-V model are more accurate than those of the N-S and S-V models
315 with respect to the phyllic zones defined by the 3D geological model.

316 It could be considered that there are spatial correlations between different
317 modeled Cu zones and geological features such as alterations and mineralogy. Several
318 samples were collected from different drillholes in different grade mineralization
319 zones of the Pulang deposit to validate the results of the fractal models. These

320 samples were analyzed by microscopic identification and XRF (X-ray fluorescence
321 spectrometry). The PL-B82 sample was collected from the drillhole situated in a
322 high-grade mineralization zone and includes a high chalcopyrite content and some
323 molybdenite (Fig. 16a). The PL-B62 sample was collected from the drillhole situated
324 in a moderate-grade mineralization zone and includes a low chalcopyrite content and
325 some pyrrhotite in the polished section (Fig. 16b). The PL-B74 sample was collected
326 from the drillhole located in a weakly mineralized zone with lower chalcopyrite
327 content and some pyrrhotite (Fig. 16c and Fig. 16d). The results obtained from the
328 mineralogy, microscopic identification and drillhole scanning by XRF of these
329 samples indicate that the Cu concentrations are 1.80%, 1.32% and 0.41% in the
330 PL-B82, PL-B62 and PL-B74 samples, respectively (Table 8).

331 **6. Conclusions**

332 In many cases, drillhole logging is dealing with the lack of proper diagnosis of
333 geological phenomena, which can undermine the delineation of mineralized zones
334 because it depends on the subjective interpretation of individual loggers, and no two
335 loggers provide the same interpretations. However, conventional geological modeling
336 based on drillhole data is fundamentally important for understanding the orebody
337 spatial structure. Grades of ore elements are not determined by conventional methods
338 of geological ore modeling, while the variation in ore grades in a mineral deposit is an
339 obvious and salient feature. Given the problems mentioned above, using a series of
340 newly established methods based on mathematical analyses such as fractal modeling
341 seems to be inevitable.

342 In this paper, the number-size (N-S), concentration-volume (C-V) and power
343 spectrum-volume (S-V) fractal models were used to delineate and recognize various
344 Cu mineralized zones of the Pulang porphyry copper deposit in the southern end of
345 the Yidun continental arc, Southwest China. All these fractal models reveal that
346 high-grade Cu mineralized zones are situated in the central and southern parts of the
347 deposit. The Cu threshold values of highly mineralized zones are 1.45% and 1.88%
348 based on the N-S and C-V fractal models. The Cu threshold of supergene enrichment

349 zones is 1.33% based on the S-V fractal model. The models of moderately
350 mineralized zones contain 0.28-1.45% Cu according to the N-S model and 1.48-1.88%
351 Cu according to the C-V model. The hypogene zones contain 0.23-1.33% Cu
352 according to the S-V model. The N-S model reveals weakly mineralized zones and
353 barren host rocks containing <0.28% Cu. In contrast, the C-V model reveals that the
354 barren host rocks contain <0.25% and that the weakly mineralized zones contain
355 0.25-1.48% Cu. The S-V model reveals that the barren host rock and leached zone
356 contain <0.23% Cu.

357 The comparison between highly mineralized zones based on the fractal models
358 and potassic zones resulting from the 3D geological model illustrates that the S-V
359 fractal model is better than the N-S and C-V model because the number of overlapped
360 voxels (A) in the S-V model is higher than those in the N-S and C-V model. The
361 overall accuracies for the C-V, N-S and S-V models are 0.50, 0.51 and 0.52,
362 respectively (Table 6), which indicates that the S-V model gives the best results for
363 identifying highly mineralized zones in the deposit. On the other hand, the correlation
364 (from OA results) between the highly mineralized zones obtained from S-V modeling
365 and the potassic alteration zones is better than those of the N-S and C-V models
366 because of a strong proportional relationship between the extension and positions of
367 the voxels in the S-V model and potassic alteration zones in the 3D geological model.

368 A comparison between phyllic alteration zones obtained from the 3D geological
369 model and moderate grade mineralization zones obtained from the fractal models
370 indicates that the OA values of the C-V, N-S and S-V fractal methods in reference to
371 the phyllic alteration zones of the geological model are 0.59, 0.54 and 0.56,
372 respectively. The overall accuracy of the moderately and weakly mineralized zones
373 obtained from C-V modeling is higher than the mineralized zones obtained from N-S
374 and S-V modeling (Table 7).

375 According to the correlation between the results driven by fractal modeling and
376 geological logging from drillholes in the Pulang porphyry copper deposit, high-grade
377 mineralization zones generated by fractal models, especially the S-V model, have a
378 better correlation with potassic alteration zones resulting from the 3D geological

379 model than from the N-S and C-V models. The highly and moderately mineralized
380 zones obtained from the fractal models are both situated in the southern and central
381 parts of the Pulang deposit and coincide with potassic and phyllic alteration zones.
382 There is a better relationship between the moderately and weakly mineralized zones
383 derived by the C-V model and the phyllic alteration zones from the 3D geological
384 model than those derived by the N-S and S-V models.

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617 **Fig.1.** Geological map of the Pulang porphyry copper deposit, SW China. Modified
618 after Yunnan Diqing Nonferrous Metal Co. Ltd., 2009.

619 **Fig.2.** Geological 3D models including lithology, alteration and 3D drill hole plot
620 with the legend of each in the Pulang porphyry copper deposit. (Scale is in m³.)

621 **Fig.3.** Photographs of alteration and mineralization in the Pulang porphyry copper
622 deposit, SW China. (a) Quartz monzonite porphyry with potassium-silicate alteration;
623 (b) Quartz diorite porphyrite with quartz-sericite alteration; (c) Quartz diorite
624 porphyrite with propylitic alteration; (d) Hornfels. Qtz=quartz; Pl=plagioclase;
625 Kfs=K-feldspar; Bt=biotite; Ser=sericite; Chl=chlorite; Ep=epidote; Py=pyrite;
626 Ccp=chalcopyrite; Mo=molybdenite; Po=pyrrhotite.

627 **Fig.4.** Cross section along exploration line 0 in the Pulang porphyry copper deposit,
628 SW China. Modified after Wang et al., 2012.

629 **Fig.5.** Histograms of (a) the Cu raw and (b) logarithmic transformation data and (c)
630 Q-Q plot of the log-transformed Cu data in the Pulang deposit.

631 **Fig.6.** The experimental semivariogram (omni-directional) of Cu data in Pulang
632 deposit.

633 **Fig.7.** The cross-validation results: (a) residual VS Cu grade; (b) the residual d
634 istribution histogram.

635 **Fig.8.** N–S log–log plot for Cu concentrations in the Pulang deposit.

636 **Fig.9.** Zones in the Pulang deposit based on thresholds defined from the N–S fractal
637 model of Cu data: (a) highly mineralized zones; (b) moderately mineralized zones; (c)
638 weakly mineralized zones and barren host rocks. (Scale is in m³.)

639 **Fig.10.** C–V log–log plot for Cu concentrations in the Pulang deposit.

640 **Fig.11.** Zones in the Pulang deposit based on thresholds defined from the C–V fractal
641 model of Cu data: (a) highly mineralized zones; (b) moderately mineralized zones; (c)
642 weakly mineralized zones; (d) barren host rock.(Scale is in m³.)

643 **Fig.12.** S–V log–log plot for Cu concentrations in the Pulang deposit.

644 **Fig.13.** Zones in the Pulang deposit based on thresholds defined from the S–V fractal
645 model of Cu data: (a) the supergene enrichment zones; (b) the hypogene zones; (c) the
646 leached zone and barren host rock (Scale is in m³.)

647 **Fig.14.** Highly mineralized zones in the Pulang deposit: (a) potassium-silicate zone
648 resulted from the 3D geological model from drillhole geological data; (b) N–S
649 modeling of Cu data; and (c) C–V modeling of Cu data; (d) S–V modeling of Cu data
650 (Scale is in m³.)

651 **Fig.15.** Moderately mineralized zones in the Pulang deposit:(a) quartz–sericite zones
652 resulted from the 3D geological model from drillhole geological data; (b) N–S
653 modeling of Cu data; and (c) C–V modeling of Cu data; (d) S–V modeling of Cu data
654 (Scale is in m³.)

655 **Fig.16.** Chalcopyrite content in several samples based on mineralographical study: (a)
656 PL-B82 sample was collected from the drillhole situated in the high grade
657 mineralization zones.; (b) PL-B62 sample was collected from the drillhole situated in
658 the moderately grade mineralization zones.; (c) and (d) PL-B74 sample was collected
659 from the drillhole located at the weakly mineralized zones.

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661 **Table 1** The results of statistical characteristics of the residual.
662 **Table 2** Thresholds concentrations obtained by using N-S model based on Cu% in
663 Pulang deposit.
664 **Table 3** Thresholds concentrations obtained by using C-V model based on Cu% in
665 Pulang deposit.
666 **Table 4** Ranges of power spectrum (S) for different mineralization zones in Pulang
667 deposit.
668 **Table 5** Matrix for comparing performance of fractal modeling results with geological
669 model. A, B, C, and D represent number of voxels in overlaps between classes in the
670 binary geological model and the binary results of fractal models (Carranza, 2011).
671 **Table 6** Overall accuracy (OA), Type I and Type II errors (T1E and T2E, respectively)
672 with respect to potassic alteration zone resulted from geological model and threshold
673 values of Cu obtained through C–V , N–S and S–V fractal modeling.
674 **Table 7** Overall accuracy (OA), Type I and Type II errors (T1E and T2E, respectively)
675 with respect to phyllic alteration zone resulted from geological model and threshold
676 values of Cu obtained through C–V, N–S and S–V fractal modeling.
677 **Table 8** Results of XRF analysis of samples collected from different mineralized
678 zones in the Pulang porphyry copper deposit.

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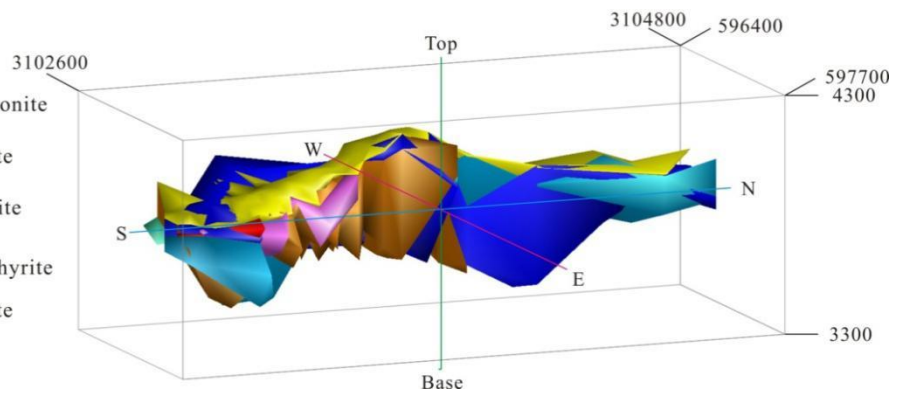
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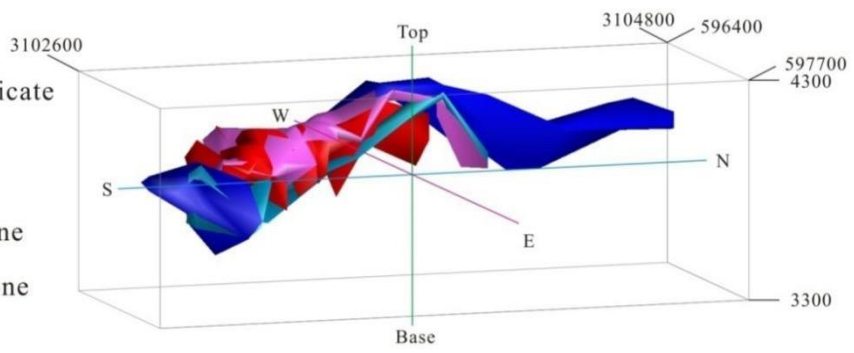
- Qesl
- Quartz monzonite porphyry
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- Diorite porphyrite
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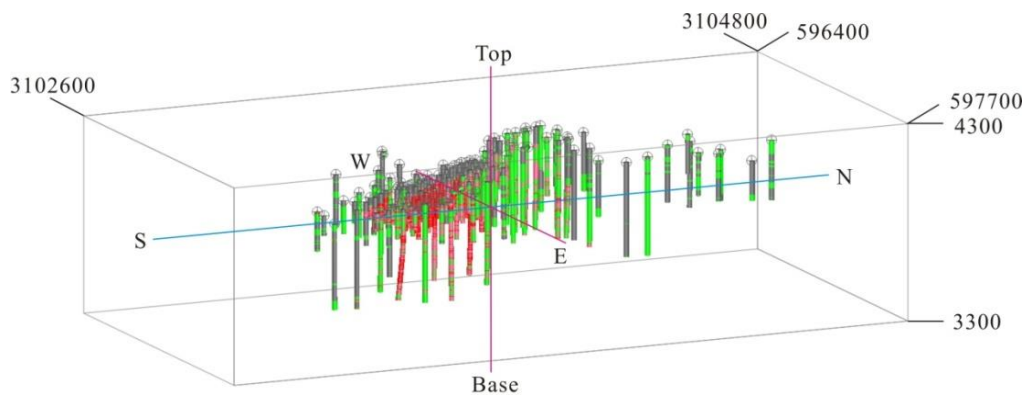
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Alteration

- Potassium-silicate Zone
- Phyllic Zone
- Propylitic Zone
- Hornstone Zone



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Fig. 2.

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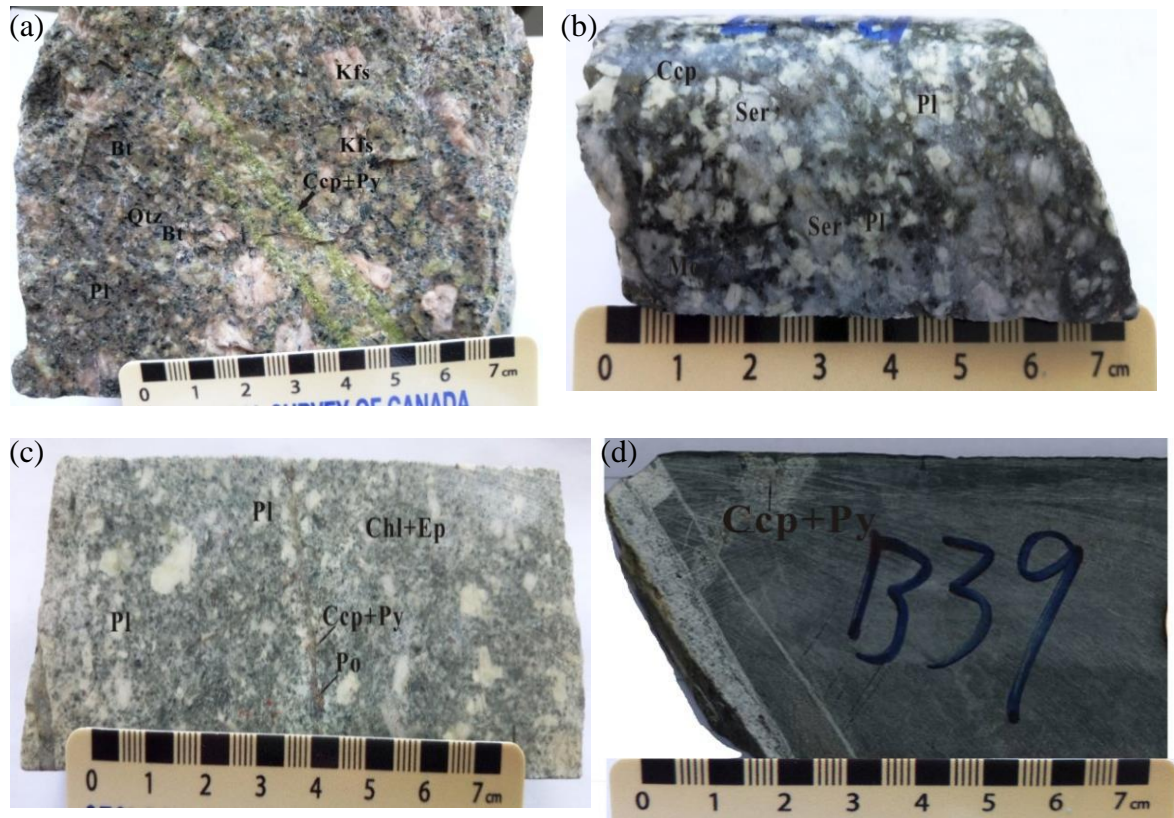


Fig. 3.

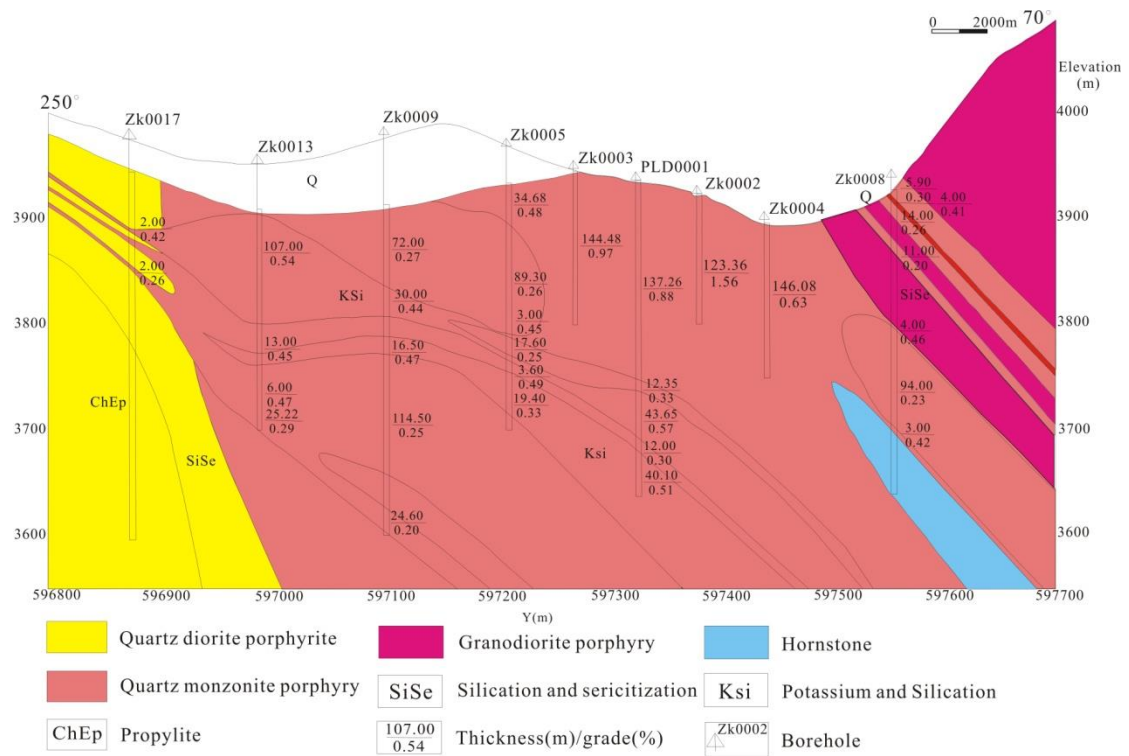
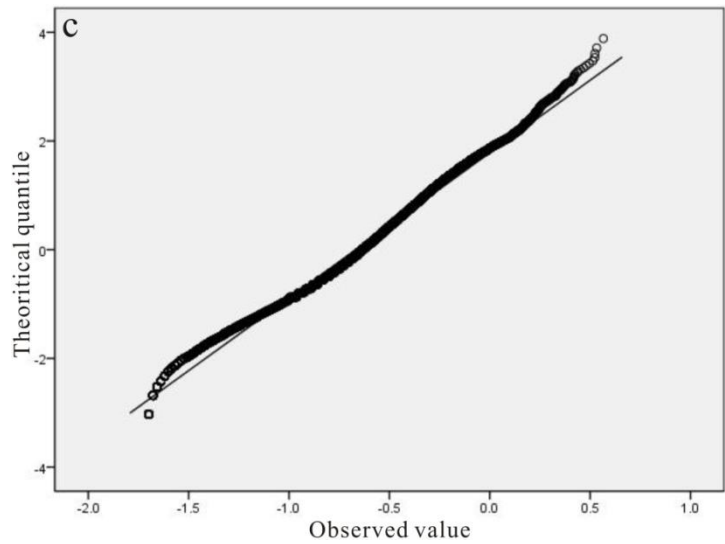
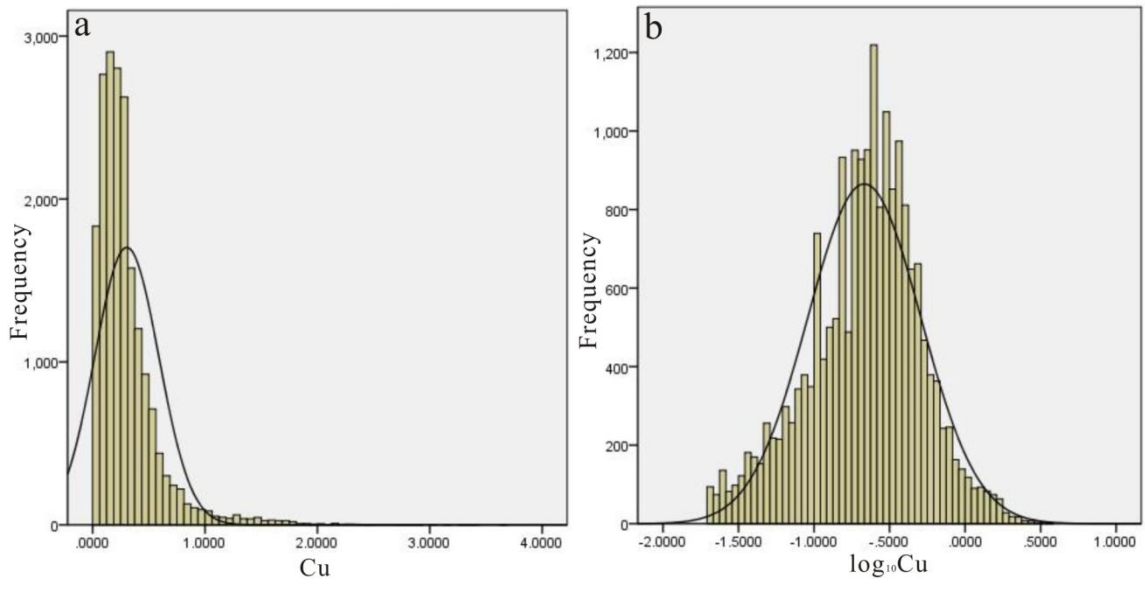


Fig. 4.

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Fig. 5.

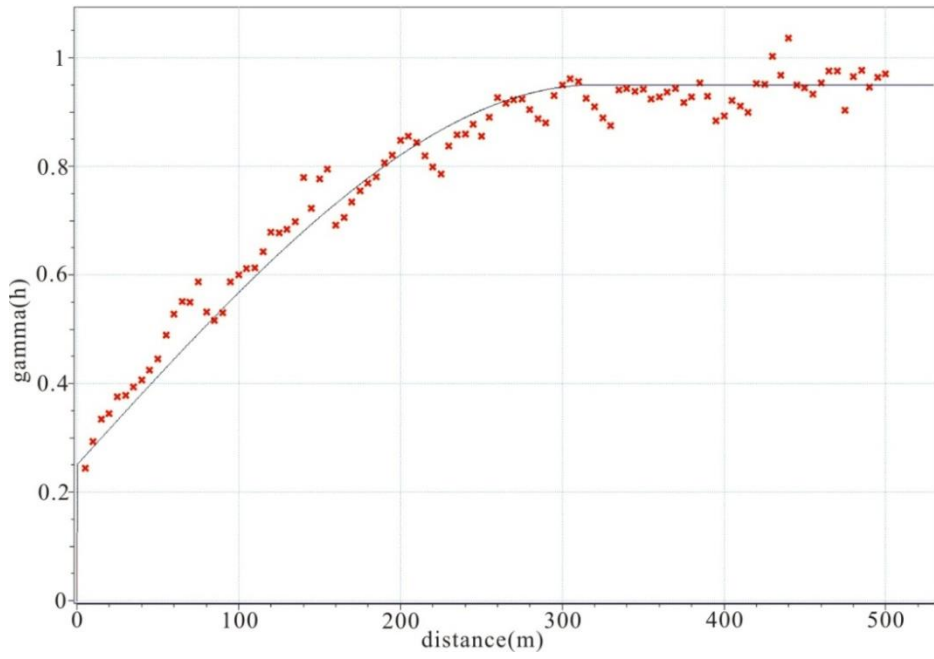


Fig. 6.

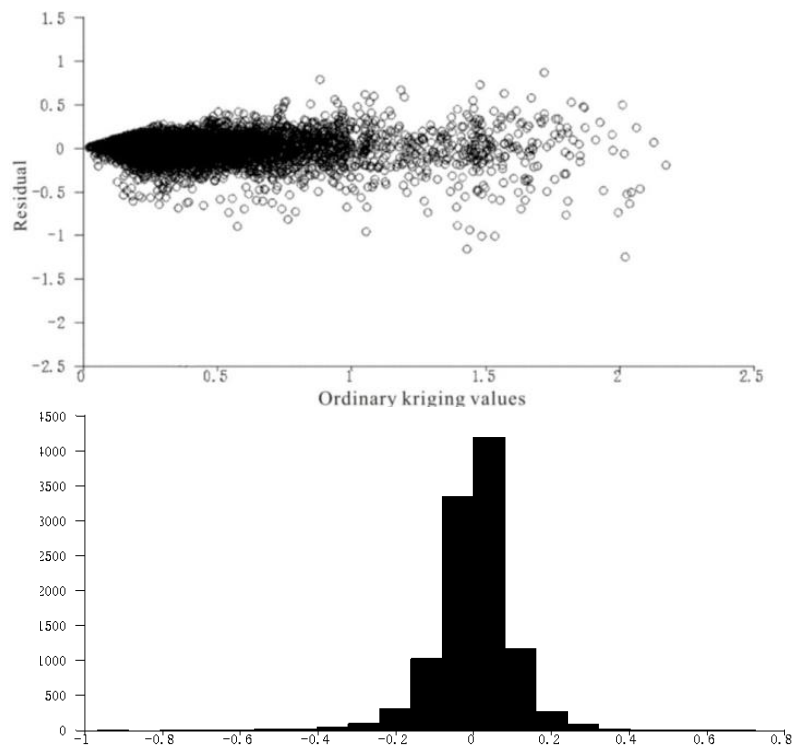


Fig. 7.

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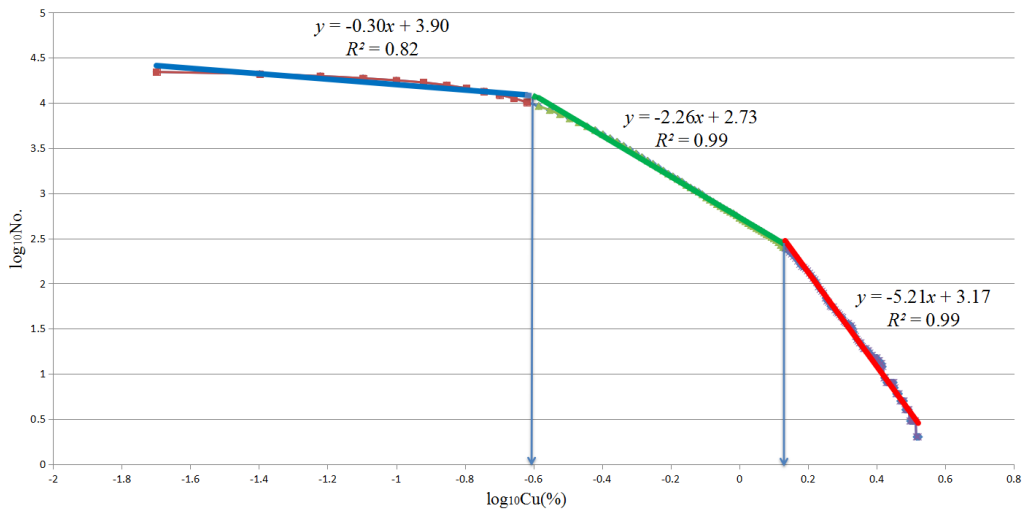


Fig. 8.

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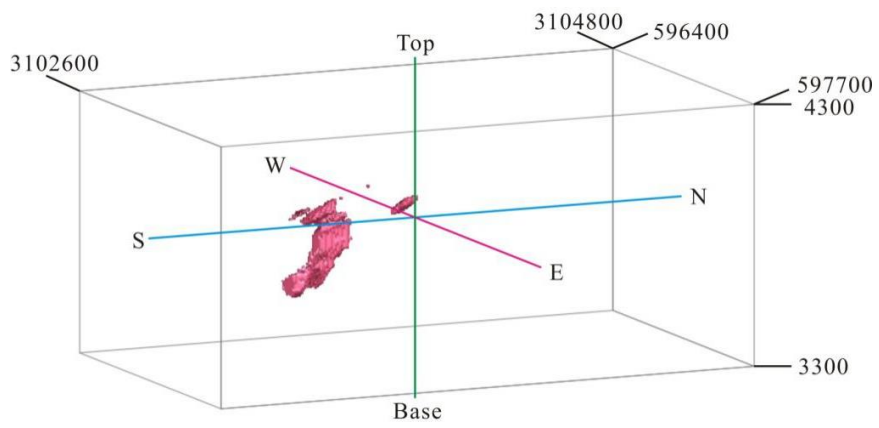
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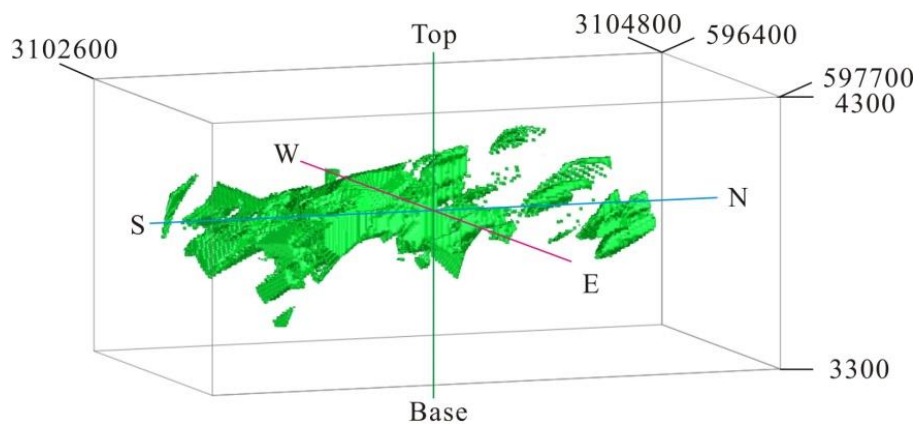
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813 (a)



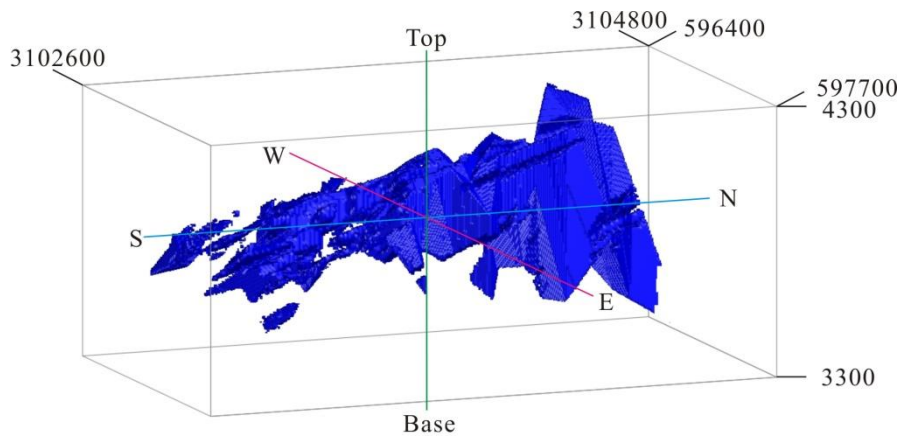
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815 (b)



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817 (c)



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Fig. 9

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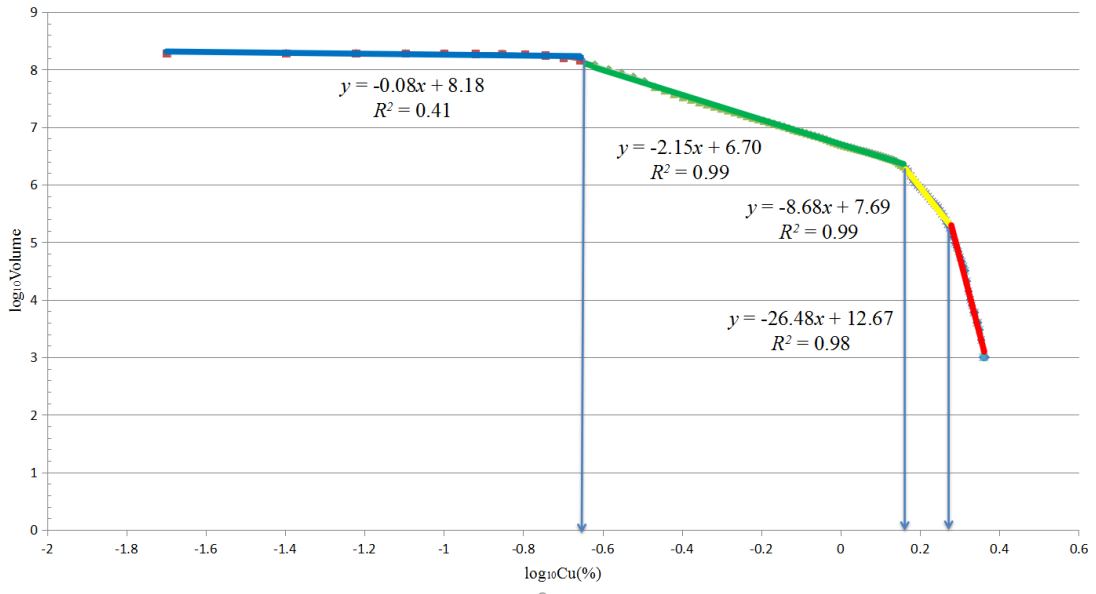


Fig. 10.

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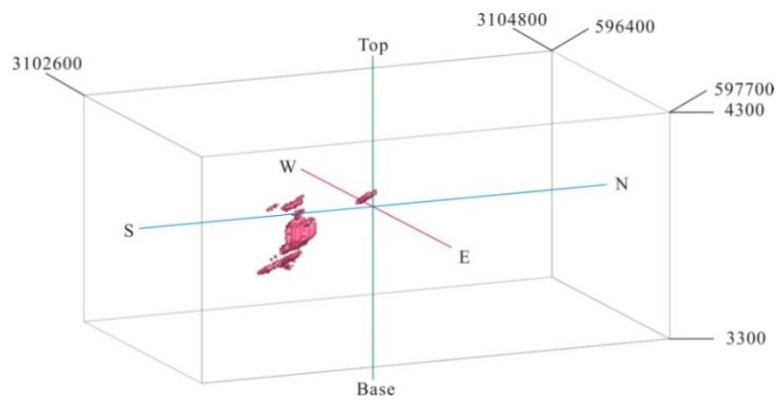
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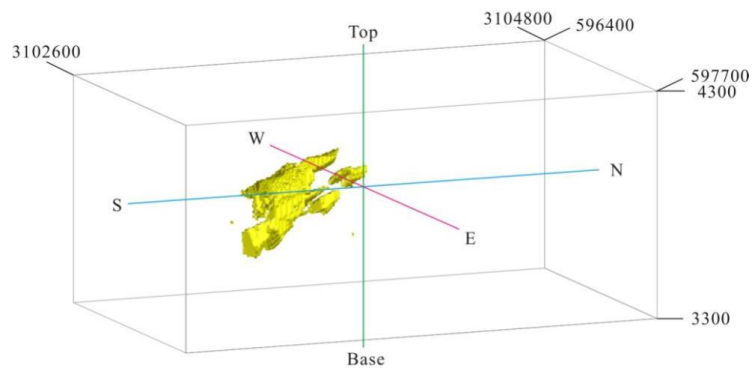
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855 (a)



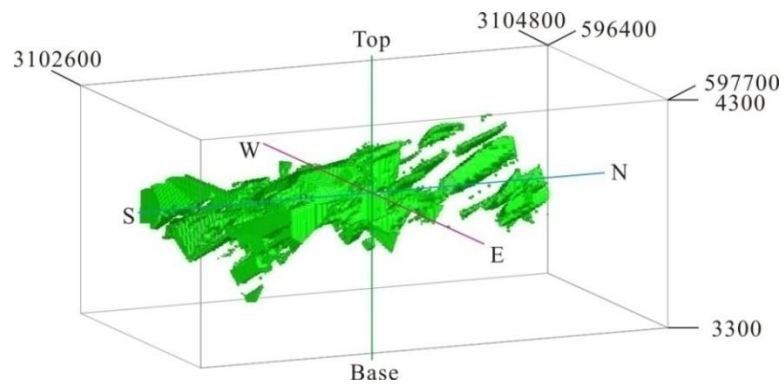
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857 (b)



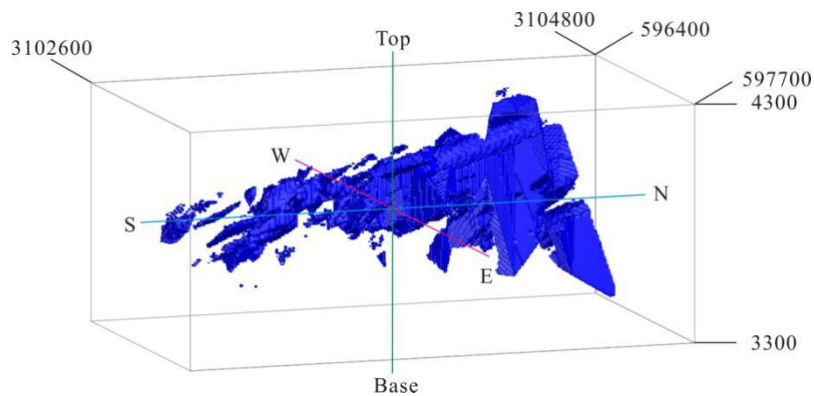
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859 (c)



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861 (d)

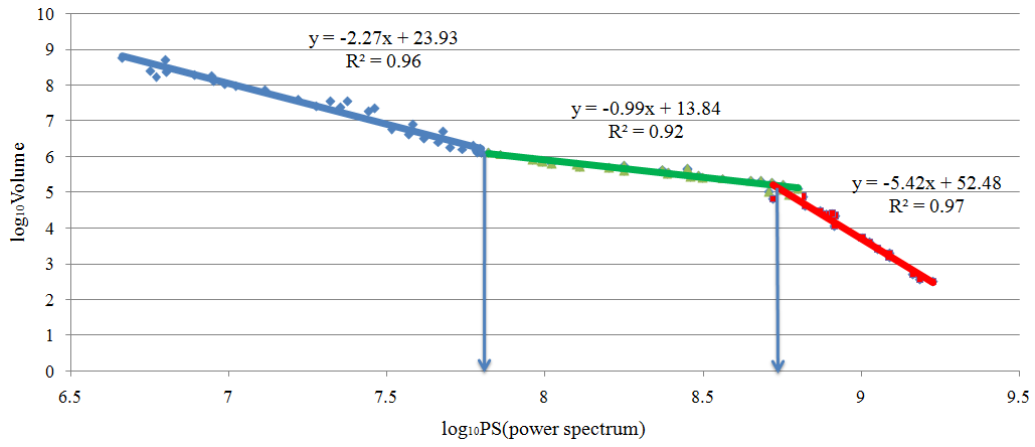


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Fig. 11.

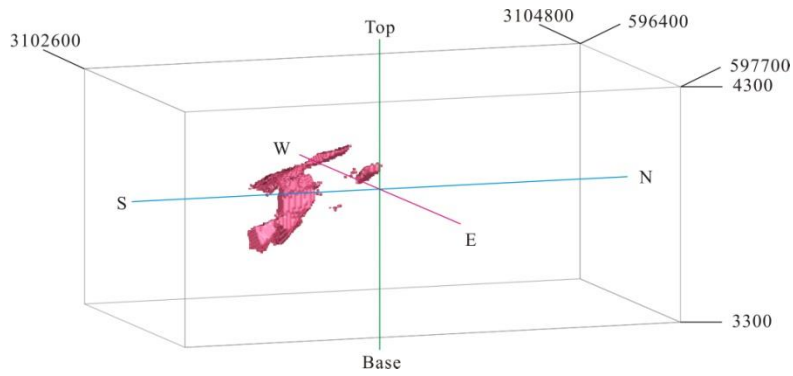


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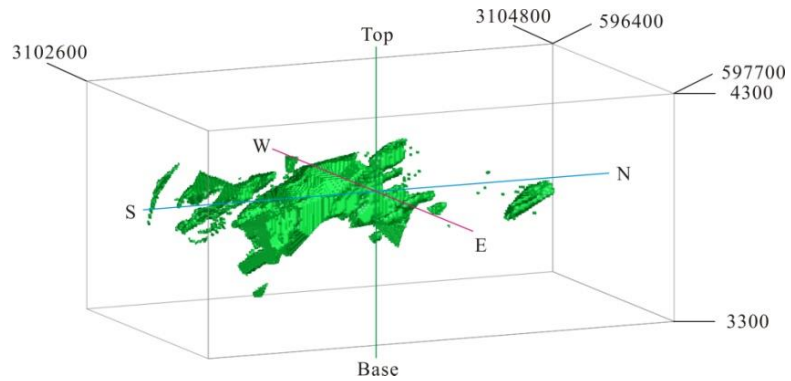
867 (a)

Fig. 12.



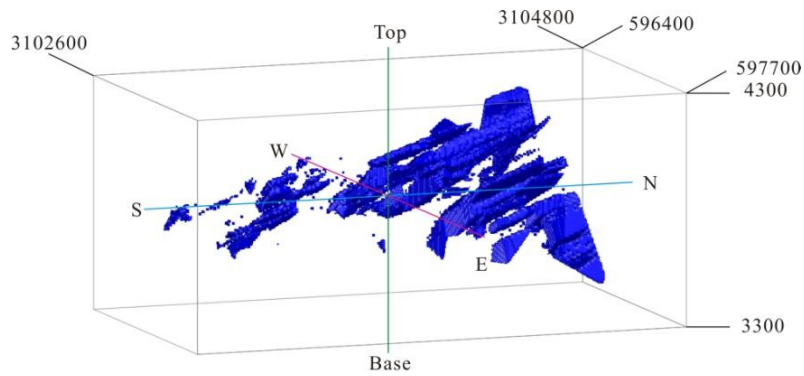
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869 (b)



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871 (c)

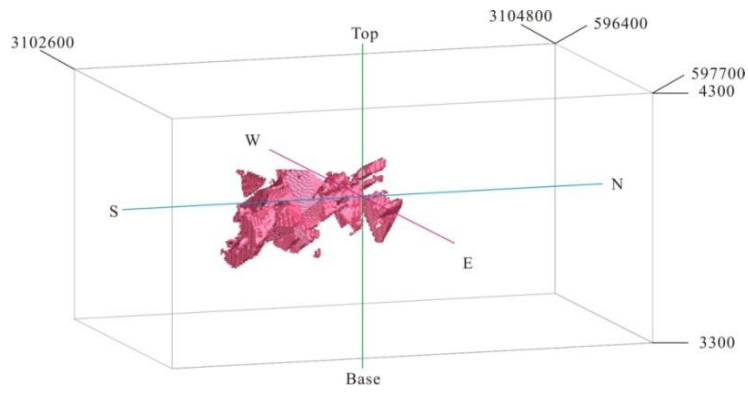


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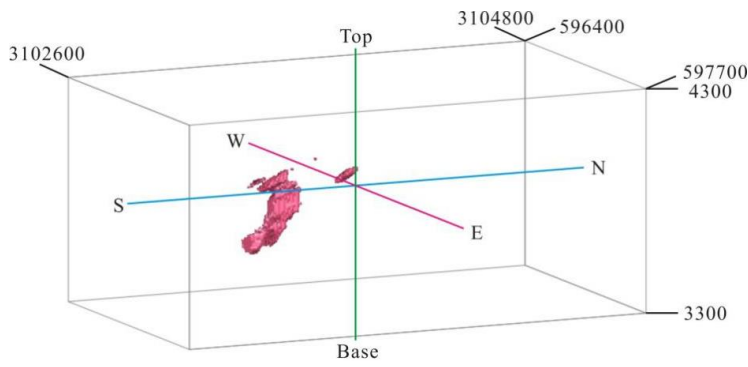
Fig. 13.

874 (a)



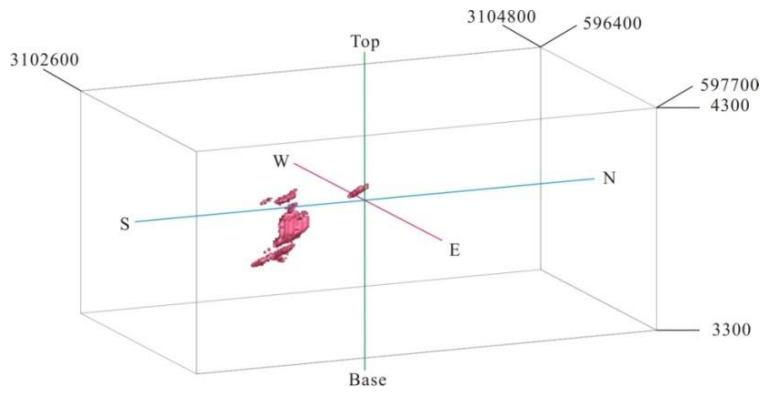
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876 (b)



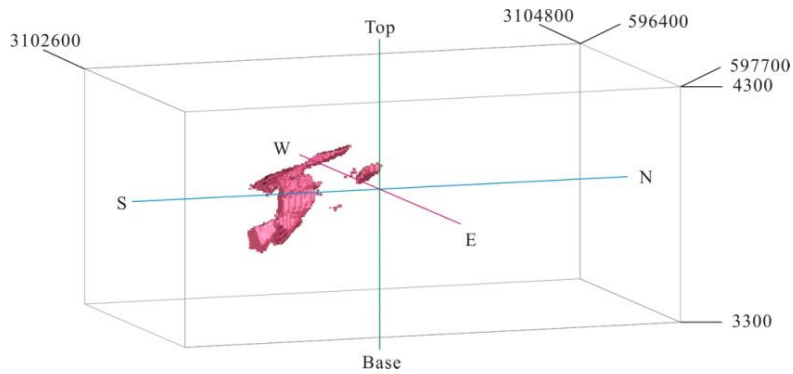
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878 (c)



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880 (d)

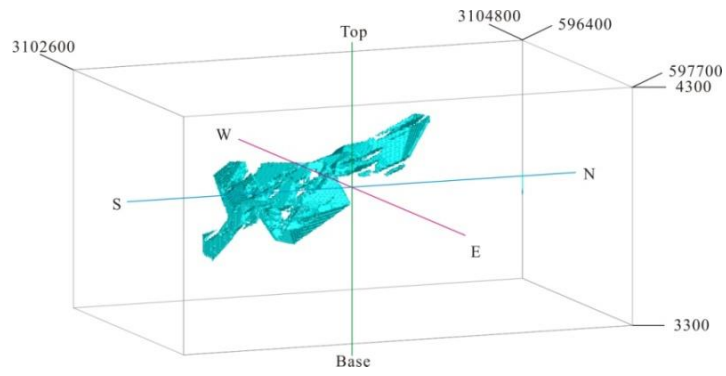


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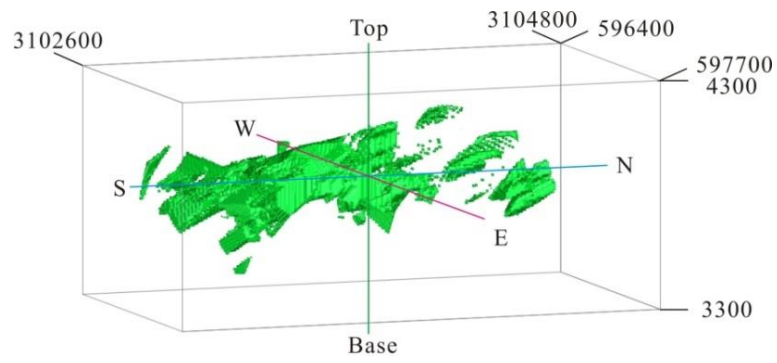
Fig. 14.

883 (a)



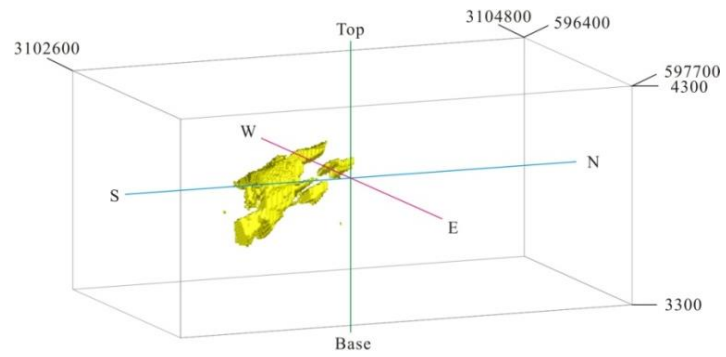
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885 (b)



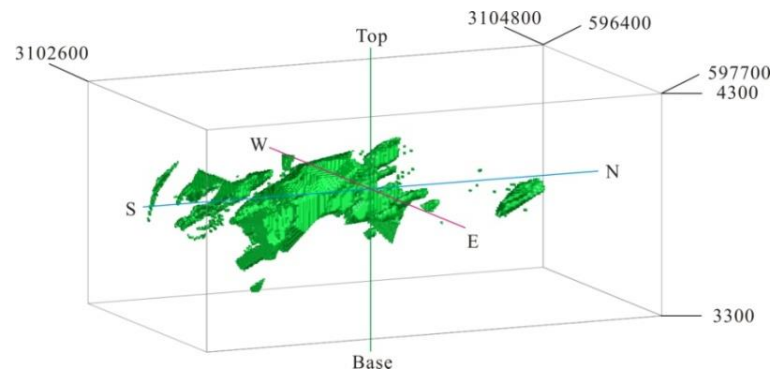
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887 (c)



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889 (d)



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Fig. 15.

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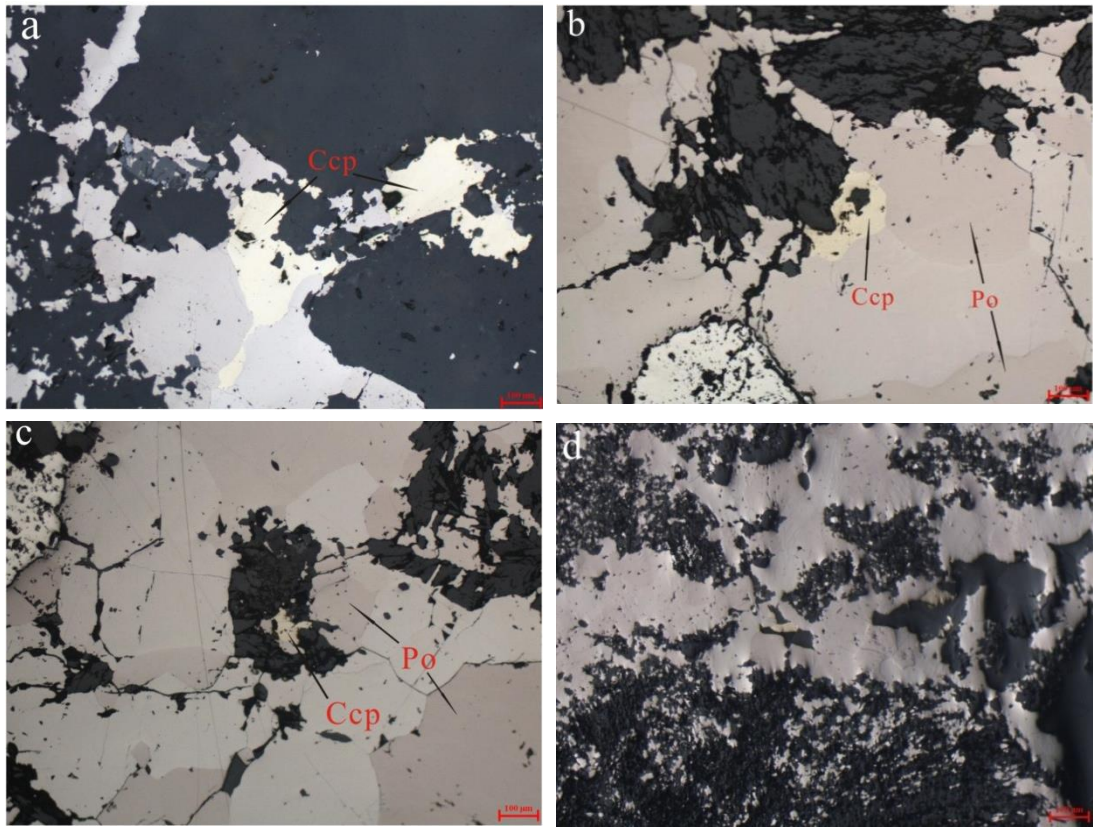


Fig. 16.

937 **Table 1**

Variables	Residual
Mean	0.000
Variance	0.016
Standard Deviation	0.127

938 **Table 2**

Mineralized zones	Thresholds (Cu%)	Range (Cu%)
Barren host rock and weakly mineralized		<0.28
Moderately mineralized	0.28	0.28-1.45
Highly mineralized	1.45	>1.45

939 **Table 3**

Mineralized zones	Thresholds (Cu%)	Range (Cu%)
Barren host rock		<0.25
Weakly mineralized	0.25	0.25–1.48
Moderately mineralized	1.48	1.48–1.88
Highly mineralized	1.88	>1.88

940 **Table 4**

Mineralized zones	PS threshold	Range of PS	Range (Cu%)
leached zone and barren host rock		<7.81	<0.23
hypogene zones	7.81	7.81-8.70	0.23-1.33
supergene enrichment zones	8.70	>8.70	>1.33

941 **Table 5**

		Geological model	
		Inside zone	Outside zone
Fractal model	Inside zone	True positive (A)	False positive (B)
	Outside zone	False negative (C)	True negative (D)
		TypeIerror=C/(A+C)	TypeIIerror=B/(B+D)
		Overallaccuracy=(A+D)/(A+B+C+D)	

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947 **Table 6**

		Potassic alteration of geological model			
		Inside zones		Outside zones	
C–V fractal model of highly mineralized zones	Inside zones	A	2850	B	1360
	Outside zones	C	77927	D	76913
		T1E	0.96	T2E	0.02
		OA		0.50	
N–S fractal model of highly mineralized zones	Inside zones	A	3092	B	1570
	Outside zones	C	75025	D	75473
		T1E	0.96	T2E	0.02
		OA		0.51	
S–V fractal model of supergene enrichment zones	Inside zones	A	4431	B	2318
	Outside zones	C	72985	D	75726
		T1E	0.94	T2E	0.03
		OA		0.52	

948 **Table 7**

		Phyllic alteration of geological model			
		Inside zones		Outside zones	
C–V fractal model of moderately and weakly mineralized zones	Inside zones	A	36518	B	48027
	Outside zones	C	25461	D	69155
		T1E	0.41	T2E	0.40
		OA		0.59	
N–S fractal model of moderately mineralized zones	Inside zones	A	35555	B	46943
	Outside zones	C	23955	D	48223
		T1E	0.40	T2E	0.49
		OA		0.54	
S–V fractal model of the hypogene zones	Inside zones	A	40080	B	44943
	Outside zones	C	26899	D	54239
		T1E	0.40	T2E	0.45
		OA		0.56	

949 **Table 8**

Sample no.	Mineralized zones obtained by fractal models	Cu (%)
PL-B74	Weakly mineralized zones	0.41
PL-B62	Moderately mineralized zones	1.32
PL-B82	Highly mineralized zones	1.80

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