

Chuchua
896228

Sample name	Ba in ppm	Ti in ppm	Sc in ppm	Ca	Cs in ppm
22369	58 +/- 7	9498 +/- 9710000	44 +/- 2	44	12 +/- 2
22375	63210 +/- 2310	10859 +/- 787500	51 +/- 2	2	<10
22382	335 +/- 8	10067 +/- 879500	46 +/- 2	24.1	12 +/- 2
22386	45820 +/- 2310	10321 +/- 788500	45 +/- 2	237.4	<10
22391	50459 +/- 2540	11239 +/- 919600	49 +/- 2	243.1	<10
22395	84424 +/- 4260	10611 +/- 757900	47 +/- 2	234.1	<10
22400	1796 +/- 15	12151 +/- 9413900	44 +/- 2	246.7	13 +/- 2
22403	2314 +/- 18	8822 +/- 778900	40 +/- 2	240.7	10 +/- 2
22417	94 +/- 8	11111 +/- 10011000	47 +/- 2	248.7	17 +/- 2

Sample name	Cr in ppm	Co in ppm	Ni in ppm	Rb in ppm	Sr in ppm	Y in ppm	Zr in ppm
22369	250 200 +/- 10	40 39 +/- 2	79 +/- 2	<20 5 +/- 6	190 236 +/- 3	32 37 +/- 2	111 +/- 2
22375	200 235 +/- 10	35 35 +/- 2	77 +/- 2	<20 < 7	64 117 +/- 2	18 35 +/- 2	109 +/- 2
22382	220 221 +/- 10	35 34 +/- 2	76 +/- 2	<20 2 +/- 6	76 86 +/- 2	32 33 +/- 2	111 +/- 2
22386	200 211 +/- 10	32 36 +/- 2	65 +/- 2	40 1 +/- 5	30 46 +/- 2	18 36 +/- 2	109 +/- 2
22391	200 230 +/- 10	35 36 +/- 2	74 +/- 2	20 < 7	36 57 +/- 2	20 42 +/- 2	118 +/- 2
22395	190 216 +/- 10	31 36 +/- 2	63 +/- 2	<20 < 7	44 84 +/- 2	14 39 +/- 2	104 +/- 2
22400	250 246 +/- 10	37 36 +/- 2	95 +/- 2	<20 2 +/- 6	74 87 +/- 2	34 37 +/- 2	130 +/- 2
22403	210 219 +/- 10	37 35 +/- 2	67 +/- 2	<30 < 7	110 127 +/- 2	28 30 +/- 2	100 +/- 2
22417	270 237 +/- 10	47 45 +/- 2	88 +/- 2	<20 8 +/- 6	90 108 +/- 2	32 39 +/- 2	123 +/- 2

use these values

Sample name	Nb in ppm	Mo in ppm	Th in ppm	U in ppm
22369	6 < 5	< 3	5 +/- 5	10 +/- 4
22375	< 2 < 5	5 +/- 3	4 +/- 5	< 4
22382	6 1 +/- 3	< 3	3 +/- 5	< 4
22386	2 < 5	< 3	4 +/- 5	0 +/- 4
22391	2 < 5	< 3	< 5	2 +/- 4
22395	< 2 < 5	< 3	< 5	< 4
22400	6 1 +/- 3	< 3	< 5	< 4
22403	6 1 +/- 3	< 3	9 +/- 5	2 +/- 3
22417	6 5 +/- 3	< 3	4 +/- 5	8 +/- 4

Cr corrected for Fe absorption

DBASE
CHUATRCE - DBF

Bill

Set up DBase file to be created

Structure	Ba	Ti	Sc	Cs	Nb	Mo	Th	U	Cu	Zn	Yb
	7	7	4	4	3	3	3	3	4	4	

Cr	Co	Ni	Rb	Sr	Y	Zr
4	3	4	3	4	3	4

X Hf
X Ta
2

other lab	Na%	Fe%	As	Se	Br	Sb	La	Ce	Nd	Sm	Eu	Yb	Lu
	5	5	3	3	4	5	5	5	5	5	4	4	4

Sample name	Ba in ppm	Ti in ppm	Sc in ppm	Cs in ppm
22369	58 +/- 7	9498 +/- 97	44 +/- 2	12 +/- 2
22375	63210 +/- 2310	10859 +/- 78	51 +/- 2	<10
22382	335 +/- 8	10067 +/- 87	46 +/- 2	12 +/- 2
22386	45820 +/- 2310	10321 +/- 78	45 +/- 2	<10
22391	50459 +/- 2540	11239 +/- 91	49 +/- 2	<10
22395	84424 +/- 4260	10611 +/- 75	47 +/- 2	<10
22400	1796 +/- 15	12151 +/- 94	44 +/- 2	13 +/- 2
22403	2314 +/- 18	8822 +/- 77	40 +/- 2	10 +/- 2
22417	94 +/- 8	11111 +/- 100	47 +/- 2	17 +/- 2

Sample name	Cr in ppm	Co in ppm	Ni in ppm	Rb in ppm	Sr in ppm	Y in ppm	Zr in ppm
22369	186 +/- 20	39 +/- 2	79 +/- 2	5 +/- 6	236 +/- 3	37 +/- 2	111 +/- 2
22375	213 +/- 20	35 +/- 2	77 +/- 2	< 7	117 +/- 2	35 +/- 2	109 +/- 2
22382	206 +/- 20	34 +/- 2	76 +/- 2	2 +/- 6	86 +/- 2	33 +/- 2	111 +/- 2
22386	192 +/- 20	36 +/- 2	65 +/- 2	1 +/- 5	46 +/- 2	36 +/- 2	109 +/- 2
22391	206 +/- 20	36 +/- 2	74 +/- 2	< 7	57 +/- 2	42 +/- 2	118 +/- 2
22395	192 +/- 20	36 +/- 2	63 +/- 2	< 7	84 +/- 2	39 +/- 2	104 +/- 2
22400	271 +/- 20	36 +/- 2	95 +/- 2	2 +/- 6	87 +/- 2	37 +/- 2	130 +/- 2
22403	204 +/- 20	35 +/- 2	67 +/- 2	< 7	127 +/- 2	30 +/- 2	100 +/- 2
22417	221 +/- 20	45 +/- 2	88 +/- 2	8 +/- 6	108 +/- 2	39 +/- 2	123 +/- 2

do not use Cr

Sample name	Nb in ppm	Mo in ppm	Th in ppm	U in ppm
22369	< 5	< 3	5 +/- 5	10 +/- 4
22375	< 5	5 +/- 3	4 +/- 5	< 4
22382	1 +/- 3	< 3	3 +/- 5	< 4
22386	< 5	< 3	4 +/- 5	0 +/- 4
22391	< 5	< 3	< 5	2 +/- 4
22395	< 5	< 3	< 5	< 4
22400	1 +/- 3	< 3	< 5	< 4
22403	1 +/- 3	< 3	9 +/- 5	2 +/- 3
22417	5 +/- 3	< 3	4 +/- 5	8 +/- 4

Bill.

(Cr) Cr. results need correcting for Fe absorption edge interference - ready in a couple of days
 meanwhile results are good to about 20% relative - worse with high iron.
 Your Fe are much the same \therefore precision will be good (1-2%) but accuracy will be worse

Paul.

Chu Chua Results

Sample name	Nb in ppm	Mo in ppm	Th in ppm	U in ppm	Cu in ppm	Zn in ppm	
22369-0.1726	< 5	< 3	5 +/- 5	< 1	10 +/- 4	75 +/- 3	88 +/- 1
22375-0.1761	< 5	5 +/- 3	4 +/- 5	.4	< 4		
22382-0.1879	1 +/- 3	< 3	3 +/- 5	.3	< 4	64 +/- 3	87 +/- 1
22386-0.1847	< 5	< 3	4 +/- 5	< 1	0 +/- 4	53 +/- 2	76 +/- 1
22391-0.1753	< 5	< 3	< 5	.6	2 +/- 4	75 +/- 3	128 +/- 1
22395-0.1891	< 5	< 3	< 5	< 3	< 4	113 +/- 3	152 +/- 2
22400-0.1826	1 +/- 3	< 3	< 5	.4	< 4	33 +/- 4	192 +/- 2
22403-0.1846	1 +/- 3	< 3	9 +/- 5	.1	2 +/- 3	66 +/- 3	102 +/- 2
22417-0.1651	5 +/- 3	< 3	4 +/- 5	.1	8 +/- 4	60 +/- 3	86 +/- 2

