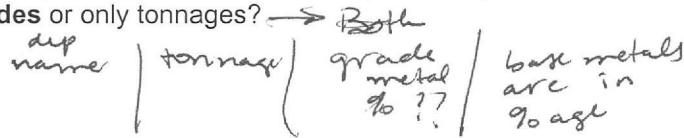


1. We are in the process of refining Yukon-specific deposit models/lists. Do I need to enter **metal grades** or only tonnages? → Both



Results from simulator are in tonnes

2. In the **coder file**, the first column is labeled **Fss**. What does that mean?

↳ don't remember perhaps has to be there

3. In the **model file**, are there grades or tonnages under the **metal columns**, and what units are being used?

grades possibly all in %

4. If estimators chose not to **estimate for 0 deposits**, leave cells blank, fill in 0, or remove columns?

check on min\_code  
- probably throws in 100

5. Does the .dat file (**coder file**) run through another program before Raw2Mark?

probably

↳ Raw 2 re-formats it into format the simulator wants

6. Does **coder data** have to go into **Min\_code** or can it be directly digitized in a **spreadsheet**?

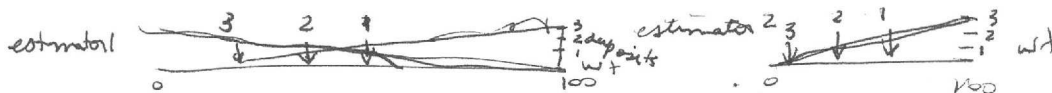
↳ not necessary

↳ comma delimited OK

7. Does Raw2Mark create a probability curve from the estimation?

8. In the **summary file**, what do the # of deposits at 50%, 10%, etc. confidence represent?

output of Raw 2 Mark



9. What is the **command file**, and where is it used?

- tells simulator to start, etc

→ call up command file, and it will give the simulator

10. Where can I **specify which tonnage curve** file names to use?

There is a model #

It may not be in the simulator

The assembler uses the .out files

Data folder  
• .out files

# .dat  
↳ model #

Phase1.exe makes batch files for each file  
(.bat)

There is a program that prepares .txt files

into .dat files

input: ~~it produces~~ a

.dat files

~~then~~ it makes .ben

e.g. 78.ben



Mark 3b

Start up menu in blue screen

Sec 4.1.2 Digital support mode prepared

Phase 1

enter model #: e.g. 30

enter file name for output

never change the intalok.dat

Must have .dat, .dat + intalok.dat in same directory

as phase1.exe

- it makes the .ben

the output file according to the name

ranking

Subject: Ranking  
Date: Fri, 14 May 1999 01:04:14 -0400  
From: Caleen E Kilby <ckilby@compuserve.com>  
To: Al Doherty <aurum@yknet.yk.ca>

Al,

The steps in ranking are outlined below, they can be done either in access or excel, Access is a safer place to sort but it is your choice. ALL SORTING STEPS MUST BE DONE ON THE WHOLE FILE>!!!!

1) Put the data in the program - each record should have (tract id, area, resource value, est90%, est50%, est10%, est5% and est1%) values from previous work.

2) Add columns - two columns after each of resource value, and the 5 estimate columns. These should be named something like resource\_index and resource\_rank, est90\_index and est90\_rank etc.

3) Populate the index columns by placing the corresponding value divided by the tract area in them. Insert the formula in the cell and drag it down the whole file. Do this for the resource\_index and the 5 estimate index columns.

4) now ready to start sorting and adding ranks. Start by sorting the whole file by the resource\_index column. Once sorted put rank numbers in the resource\_rank column. They should run from 1 to the number of tracts in the assessment area. 1 for the lowest rank and the largest number for the highest rank. If there was zero resources the rank number should be zero.

Once the ranks have been added to the resource\_index column move to the est90\_index column and sort the whole file by it. Then enter the rank numbers in the est90\_rank column. Again remember if there is not est90 value then there should be a zero value in the rank column. Rankings may start at 21 with everything less than that having a zero value.

Continue doing this for the rest of the estimate columns.

5) add two columns at far right of the file, a SCORE column and a RANK column.

## ranking

the score column should contain the weighted sum of all the rankings. OR  
the following formula

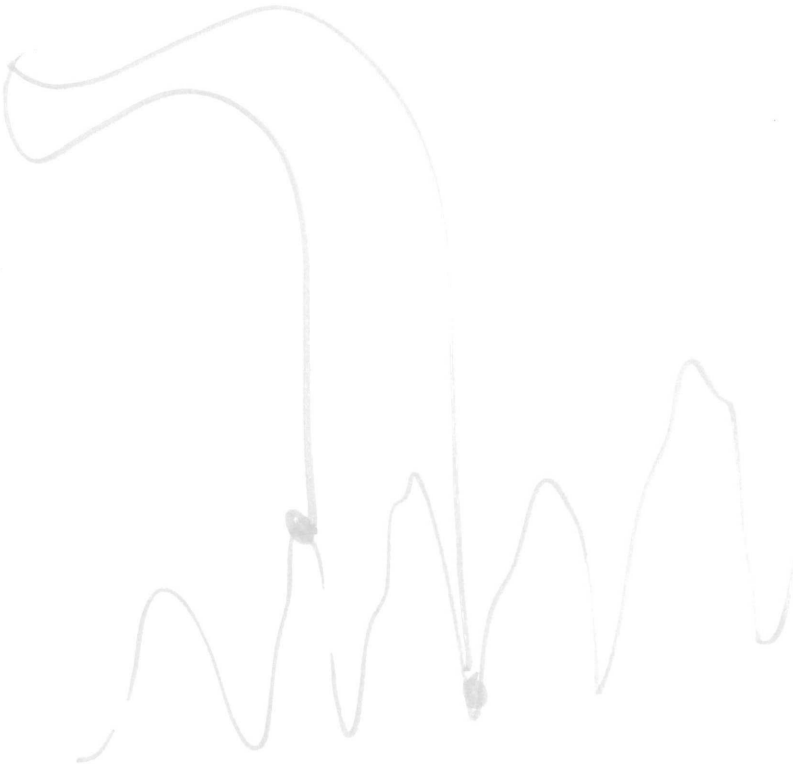
=  
resource\_rank\*1+est90\_rank\*.9+est50\_rank\*.5+est10\_rank\*.1+est5\_rank\*.05  
+est  
1\_rank\*.01

Once all the score values are in place, sort the whole file by the SCORE  
value then and the rank numbers in the rank column. This is then the final  
ranking of the assessment area.

Good Luck.

RS 36 2 52 28

4990



2.5

|



## **Anna.Fonseca**

---

**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Wednesday, February 02, 2000 2:50 PM  
**To:** Anna Fonseca  
**Subject:** mineral deposit simulator help

Hi Anna,

Thanks for your message yesterday. We would be happy to assist you in your mineral deposit simulation. Our rates would be the same as last year at \$60 Cdn / hour. Sometimes I may be able to answer your questions, but most of the time it will be Ward. This email address is fine for queries, and our phone number is below. In addition Ward also has a Telus email address wkilby@telus.net. He checks that regularly as well. If I can't answer something I will pass it to Ward as soon as possible.

With regards to your question about the "indicated reserves" tonnages for Howards Pass - Ward says he does not know where that number came from. Lots of people worked on the deposit models and there may have been some "chicanery" creeping in from time to time. The models should be based on global resource, as you mentioned. I don't know who was involved in that model, but it may have been intentional. It is possible that the Howards Pass deposit has very large resource tonnages, and to facilitate the deposit curve and median tonnage estimate a smaller number was given to it and named "indicated reserves". The numbers are not likely real as the property has never moved to the mining stage. Most importantly, whatever is used....resources or reserves,etc., statistically they must all be the same or the result is meaningless.

Please say hello to Daniele for me (Caleen). I went into the Yukon room at Roundup to say hello, but must have missed her.

Talk to you soon,  
Caleen

Caleen E. Kilby, P.Geol.  
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## **Anna.Fonseca**

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**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Thursday, February 24, 2000 11:30 PM  
**To:** Anna.Fonseca  
**Subject:** RE: mineral deposit simulator help

Anna,

The .def file tells phase1 which column each commodity is in.

You must make up the .def files yourself. This can be done easily with a text editor. Just follow the format of some existing .def files or edit an existing one. Best to put all the model files for an analysis in a completely different directory so things don't get mixed-up.

In the .def file the numbers below the model name and description are:

- the first number is the number of commodities to follow
- the numbers below this are the commodity numbers in column order (right to left) in the .dat file.
- to see what these number codes represent look in the Mark3b.bas program with a text editor. Early on in the program will be a list of commodities and associated numbers. These are the numbers the simulator is expecting to represent a commodity. If required you could change the commodity-number association and then recompile the source program.

It is also worth reading the Read\_me.pgm file with a text editor. It gives a reasonable account of what is going on. (in the execute directory)

Hope that helps.

Ward

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## **Anna.Fonseca**

---

**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Thursday, March 02, 2000 8:09 PM  
**To:** Anna.Fonseca  
**Subject:** .def files

Anna,

Your correct (right). Even though I do confuse left and right (usually in driving tests). I meant left to right.

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## **Anna.Fonseca**

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**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Friday, March 03, 2000 2:16 PM  
**To:** Anna.Fonseca  
**Subject:** Mineral Deposit Simulator - phase 1

Hi Anna,

1. Regarding the .def files: It looks like there need to be 10 values after the initial number. So it may be best to follow the same format and just use a text editor to modify an existing .def file to make a new one.
2. Regarding the models: You can make up your own list but will have to recompile the source code. Or.....just keep your own records of the model numbers and what they mean.
3. Regarding "minor podiform chromite": Singer & Page have a grade and tonnage model for Minor Podiform Chromite in Cox & Singer's Mineral Deposit Models Bulletin (1986). They list deposits & references. Do you have access to this? Or.....I can fax you the pages. Let me know..... (and a fax number).

bye,  
Caleen

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## **Anna.Fonseca**

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**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Wednesday, March 08, 2000 9:47 AM  
**To:** Anna.Fonseca  
**Subject:** phase 1 error messages

Hi Anna,

Ward left this morning for Toronto and was so busy getting his talks ready, it was hard to fit everything else in. But we discussed your two queries last night and I have some replies for you.

With regards to output file error messages:

This is likely due to too few deposits in the model and/or non-consistent relationships between commodities from one deposit to another. There is no real way to correct this if you are happy with the deposits in the model. It is telling you that the grade and tonnage relationships are very dissimilar between deposits. You must then question if the deposits belong to the same model.

Caleen

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## **Anna.Fonseca**

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**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Wednesday, March 08, 2000 9:53 AM  
**To:** Anna.Fonseca  
**Subject:** RE: Raw2Mark

Anna:-

Each run must contain the same number of estimators: i.e.) all the estimates made by 3 people together, then all the estimates by 2 people in another run. If things like the weighting, etc. do not add up then there will be problems. That's why we used the data entry program. The only real way we can sort this out is to see the data that was used for input. The problem is in the data.

Caleen

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## **Anna.Fonseca**

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**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Friday, March 10, 2000 10:00 AM  
**To:** Anna Fonseca  
**Subject:** An answer

Hi Anna,

The data should be sorted by tract number.

bye,  
Caleen

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## **Anna.Fonseca**

---

**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Friday, March 10, 2000 1:48 PM  
**To:** Anna.Fonseca  
**Subject:** Raw2Mark.exe

Hi Anna,

I have received both your emails. I think the problem IS likely with the file created by Excel. Ward has complained about Excel for years - everyone uses it instead of databases (although HE does it as well), and it's a spreadsheet application. I recall last year that we had to hand correct the files. I have sent both your emails on to Ward in Toronto. We don't want him to have too much fun on Friday night. I recall last year that a similar problem occurred and that it was the existence of commas at the END of each line....put there by excel. Try removing the trailing commas. Also, make sure all the numbers are real numbers and not in scientific notation.

If that doesn't work, perhaps Ward will recognize the problem.

bye,  
Caleen

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## **Anna.Fonseca**

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**From:** Caleen E Kilby [ckilby@compuserve.com]  
**Sent:** Thursday, March 16, 2000 11:45 AM  
**To:** Anna.Fonseca  
**Subject:** command file

Hi Anna,

In reply to your attached file with questions, ward sent the following comments. Please let me know if this answers your queries.

The sequence of numbers in the Raw2mrk command output file are the same as would be entered when running the simulator in the interactive mode. If you try running it in interactive mode you will see all the available options, the command file is simply answering the questions in order.

So if you run Mark3b.exe and compare the values in the command file you will see what is being requested and supplied. This way you could then go in with a text editor and change the command file to make better headings etc.

Hope that helps.  
Caleen

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