

## **PBDAT version 1.24**

*PBDAT* is an obsolete, but possibly still-useful program for transforming raw mass-spectrometric and laboratory data into final, blank, spike, and fractionation-corrected U-Th-Pb isotopic ratios and associated ages, together with all relevant errors and error-correlations.

### **Installation Instructions**

- 1) Download the PBDAT.ZIP file and User's Manual from this website.
- 2) Extract the various files from PBDAT.ZIP.
- 3) Make a new subdirectory on the hard disk (\PBDAT for example).
- 4) Copy all of the files from PBDAT.ZIP to the new folder of the hard disk.
- 5) Run PBDAT from the PBDAT folder.

The first time you run the program, select the type of display that you're using from the "Output" menu (press ALT-O or use the mouse to pull down the "Output" menu). Your choices will be "Color", "Mono- chrome", or "Compaq" (ignore the last unless you have a monochrome Compaq computer from the 1980's). Next, select the default directory for your LabData and Raw-Datafiles with the "Output Format & File Location" option from the "Output" menu. Browse through the pull-down menus and HELP screens to get an idea of how the program works (press F1 for help with the pull-down menus).

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