

## Fractint for Windows (Winfract)

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### Help Topics

[New Features in this Version](#)

[Bugs and Limitations in this Version](#)

[File Menu](#)

[Fractal Formula Selection](#)

[Image Sizing and Color selection](#)

[Coordinate Box options](#)

[Zooming in on an Image, Mandelbrot/Julia Toggling](#)

[Options and Doodads](#)

[Color-Cycling](#)

[Getting a Status Report](#)

[Distribution Policy and Procedures, Contacting the Authors](#)

[A list of Fractint Authors](#)

## **New Features in this Version**

The internal Fractal engine in this release is from Fractint 16.0. (Note that many of the other options introduced in Fractint 16.0, such as its new "demo mode", are not included in this release of Winfract.)

When generated using a Windows video driver capable of 256 simultaneous colors, Plasma Clouds now use only 236 of those colors, leaving 20 colors untouched to avoid conflict with the 20 palette entries reserved by Windows. No more "black holes" in your plasma clouds.

The Color-Cycling option is grayed out and non-selectable if your Windows video driver does not support palettes (and most 16-color Windows video drivers don't).

The Status! option has been moved from the top of the main menu bar into the "Options" menu.

The following New Features are all from Mark Peterson:

**Window Sizing** option - when selected, causes Winfract's window to automatically resize to the size of the fractal image.

**Coordinate Box** option - when selected, causes a Coordinate box to display, showing the position of the cursor in Rectangular, Polar, or Pixel coordinates.

**Zoom Bar** - when selected, causes Winfract to switch to an alternate method of generating a Zoom-box, using a "Zoom-bar" to select the size of the Zoom-box (the Zoom-box is moved around by moving the cursor to a position inside the Zoom-box, holding down the left mouse-button, and moving the cursor and the Zoom-box together).

**SSTOOLS.INI** updating - Winfract now automatically updates your SSTOOLS.INI file (creating a new one if none currently exists) with a special [winfract] section that contains the status of Winfract when you exited the program. When first firing up, Winfract reads this file to determine its start-up options.

When **Running Iconized**, the Winfract ICON is now replaced by an iconized version of the current fractal image. A quick glance at the Icon will now let you know approximately what Winfract is doing.

## Bugs and Limitations in this Version

Fractint for Windows is (and will continue to be) a "port" of Fractint for DOS to the Windows environment, retaining the Fractint for DOS fractal and GIF engines, but replacing its front-end and graphics interface with a Windows engine. As such, its development will always "trail" that of Fractint for DOS, except where the Windows version adds functionality simply by virtue of the Windows interface.

There are four main causes behind the various bugs and limitations in this program. For future brevity, these causes are tagged FFD (problems related to the fact that much of the code is and will continue to be from Fractint For Dos), WLIM (limitations caused by Windows, or some Windows video drivers that aren't limitations under MS-DOS), NWP (a Novice Windows Programmer write the original program, and he often had no idea what he was doing), and NYI (Not Yet Implemented - hey, we're working on it!). FFD and WLIM limitations are probably permanent - hopefully NWP and NYI problems are less so.

- You can run only one "instance" of Fractint for Windows. FFD: Fractint for DOS is riddled with initialized FAR data ("char far myvalue = 0;"), and the Windows SDK silently but firmly tags any program containing code like that as a single-instance program.

- Palette support is limited. WLIM: The "stock" VGA driver supplied with Win 3.0 doesn't support palette-modification by applications at all (most third-party 256-color SuperVGA drivers do, though). "Plasma Clouds" look Godawful using anything less than a 256-color Windows driver. Color-cycling is limited to video drivers which support palette-manipulation (and, for 16-color drivers, color-cycling affects the background windows as well). Note that 16-bit and 24-bit "true color" Windows video drivers don't \*have\* a palette to manipulate, so we can't color-cycle them.

- Fractint for Windows is not as "background process" friendly as it should be when it updates the screen image - if it is updating a large image in its entirety, it can grab the machine for seconds at a time. NWP - we're making attempts to update the screen in periodic "chunks", but we need to improve some more in this area.

- You cannot "abort" the File Print operations (File Save and File Open can be aborted simply by selecting another option). NWP, NYI: We're performing the entire print operation with a single 'StretchDIBits()' call. For now, just fire up a print operation, sit back, and wait for the results. Also, don't expect to run a background process (like a file transfer) while Fractint is printing. Finally, the File Print option only works on images up to 800x600 on my 386 with 6MB of RAM and LaserJet II, for some reason.

- There are lots of places where this program doesn't thoroughly check that Windows-specific procedures that it invokes ran correctly, and it crashes on occasion because of that. NWP, NYI: Basically, if we haven't hit a particular bug yet, we haven't adjusted the program to look for it.

- Lots of little doodads aren't in this version yet. NYI.

- There are \*far\* too many places in this document that say "Read Fractint-for-DOS's FRACTINT.DOC file for an explanation". NYI.

## File Menu

File **Open** loads either Fractint-for-DOS-style or "generic" GIF files. Note that the correct palette isn't completely displayed on the screen. Windows "reserves" the first twenty palettes for its own use, and "adjusts" Fractint's images accordingly - and for 16-color drivers like the VGA driver supplied with Windows that don't support palette-manipulation by applications programs, cheerfully ignores our attempts at palette-manipulation. If the restored image is resumable (a feature added with version 14.0 of FFD and version 1.2 of FFW), the image will resume calculating as soon as it is loaded.

File **Save** saves your current image as a GIF (version 89a) file, compatible with Fractint-for-DOS, your favorite DOS GIF viewers, and (of course) the "File Open" option. Again, the palette you see on the screen may not be the one that the program is "using" (and gets saved to the disk file). If the saved image is resumable (a feature added with version 14.0 of FFD and version 1.2 of FFW), the image will resume calculating as soon as it is finished saving.

File **3D Open** and File **3D Overlay** load in your image using "3D" transformations. The 3D Open option clears your image first and then loads the new one, while the Overlay option leaves your original image intact and adds the new image over it. For a full explanation of the zillion or so 3D parameters, read Fractint-for-DOS's FRACTINT.DOC file.

File **Print** does the best job it can sending the image to your printer. The program does not "dither" for black-and-white printers - it just "candy stripes" adjacent colors the same way that Fractint-for-DOS does.

**Read Color-Map** and **Write Color-Map** load and save external color-maps in the same manner as Fractint-for-DOS (within Windows' limitations).

File **Copy to Clipboard** copies the currently-displayed image to the Windows clipboard (in "Device-Independent Bitmap" format) where your other Windows programs can collect it. Note that, for 256-color video drivers and 256-color images, the colors in the clipboard won't visibly match the colors in Winfract's window until you change your "focus" to the clipboard's window. That's because Winfract has warned the Clipboard that it is reserving the right to color-cycle its colors, so the Clipboard is just using Windows' default 20 colors to display its colors while Winfract has control of the screen.

GIF and "Graphics Interchange Format" are trademarks of CompuServe Incorporated, an H&R Block Company.

## Fractal Formula Selection

Using the Options **Formula** option, you can select any of over 60 fractal types (every fractal type that is available in version 16.0 of Fractint-for-DOS). After selecting a fractal type, a dialogue box pops up and prompts you for any formula parameters and the screen corners (all with reasonable default values).

A brief list of Fractal types (and their formulas) included in this release (Note: this list is from Fractint 15.0, and doesn't include those fractals new to 16.0):

- mandel** = 'Classic' Mandelbrot fractals using 32-bit integer math for speed.  
 $z(0) = 0; z(n+1) = z(n)^2 + C$ , where  $C = Xcoord + i * Ycoord$ .  
Two optional params: real and imaginary perturbations of  $z(0)$ .
- julia** = 'Classic' Julia set fractals using 32-bit integer math for speed.  
 $z(0) = Xcoord + i * Ycoord; z(n+1) = z(n)^2 + C$ .  
Two params required: real and imaginary parts of  $C$ .
- newton**, = Newton Domains-of-attraction (only the coloring schemes are  
**newtbasin** different). First param: the power (from 3 to 10) of the eqn.  
If param=4, the eqn is  $z(n+1) = (3*z(n)^4+1)/(4*z(n)^3)$ .  
Second param is flag to turn on alternating stripes in basin mode.
- complexnewton**, = Newton's fractal type extended to complex numbers.
- complexbasin** Calculates roots of  $z^a + b$ , where both 'a' and 'b' complex.
- lambda** = Classic Lambda fractal.  $z(0) = Xcoord + i * Ycoord$ ;  
 $z(n+1) = Lambda*z(n)*(1 - z(n)^2)$ . Two params: real, imaginary  
parts of Lambda. 'Julia' variant of Mandellambda
- mandellambda**= 'Mandelbrot-Equivalent' for the lambda fractal.  $z(0) = .5$ ;  
 $lambda = Xcoord + i * Ycoord; z(n+1) = lambda*z(n)*(1 - z(n)^2)$ .  
Parameters are perturbations of  $z(0)$ .
- barnsley1** = Alternative 'Mandelbrot'.  $z(0) = Xcoord + i * Ycoord$ ;  
 $z(n+1) = (z-1)*C$  if  $Real(z) \geq 0$ , else  $= (z+1)*modulus(C)/C$ ,  
where  $C = Xcoord + i * Ycoord$ . Params are perturbations of  $z(0)$ .
- barnsleyj1** = 'Julia' corresponding to barnsley1.  $z(0) = Xcoord + i * Ycoord$ ;  
 $z(n+1) = (z-1)*C$  if  $Real(z) \geq 0$ , else  $= (z+1)*modulus(C)/C$ .  
Two params required: real and imaginary parts of  $C$ .
- barnsley2** = Another alternative 'Mandelbrot'.  $z(0) = 0; z(n+1) = (z-1)*C$   
if  $Real(z)*Imag(C) + real(C)*imag(z) \geq 0$ , else  $z(n+1) =$   
 $(z+1)*C$ , where  $C = Xcoord + i * Ycoord$ . Parameters are  
perturbations of  $z(0)$ .
- barnsleyj2** = 'Julia' corresponding to barnsley2.  $z(0) = Xcoord + i * Ycoord$ ;  
 $z(n+1) = (z-1)*C$  if  $Real(z)*Imag(C) + real(C)*imag(z) \geq 0$ ,  
else  $= z(n+1) = (z+1)*C$ . Two params: real and imaginary parts of  $C$ .
- barnsley3**,= Another Barnsley 'Mandelbrot', 'Julia' pair.
- barnsleyj3** See Fractint-for-DOS's FRACTINT.DOC for details.
- sierpinski** = Sierpinski gasket - a Julia set that produces a 'Swiss cheese  
triangle'.  $z(n+1) = (2*x, 2*y - 1)$  if  $y > .5$ ; else  $(2*x-1, 2*y)$   
if  $(x > .5)$ ; else  $(2*x, 2*y)$ . No parameters.
- ifs** = Barnsley IFS Fractal (a fern unless an alternate IFS map has  
been defined using the 'ifs=' command-line option).
- ifs3d** = Barnsley 3D IFS Fractal (a fern unless an alternate IFS map has  
been defined using the 'ifs3d=' command-line option).
- marksmandel**= A variant of the mandel-lambda fractal.  $z(0) = 0$ ;  
 $z(n+1) = ((Xcoord+i*Ycoord)^exp)*z(n) + (Xcoord+i*Ycoord)$ .  
Parameters are perturbations of  $z(0)$ .
- marksjulia** = A variant of the julia-lambda fractal.  
 $z(0) = Xcoord + i * Ycoord; z(n+1) = (z(0)^exp)*z(n) + z(0)$ .

**cmplxmarksjul**, = The above two types generalized with 'exp' a complex rather than a real number

**cmplxmarksmand**

**julibrot** = 'Julibrot' 4-dimensional fractals. See Fractint-for-DOS's FRACTINT.DOC for an description of these fractals (and a description of the prompts involved in invoking them).

**unity** = Mark Peterson's 'Unity' fractal type. Truly Wierd - See Fractint-for-DOS's FRACTINT.DOC for a full description.

**formula** = Formula interpreter - write your own formulas as text files! See Fractint-for-DOS's FRACTINT.DOC for instructions on this one.

**lambdafn** = lambda-fn fractal.  $z(0) = Xcoord + i * Ycoord$ ;  $z(n+1) = lambda * fn(z(n))$ . Two params: real, imag portions of lambda. 'Julia' variant of Mandelfn. 'fn' is a variable function, and may be one of sin, cos, sinh, cosh, exp, log, or sqr.

**mandelfn** = 'Mandelbrot-Equivalent' for the lambda-fn fractal.  $z(0) = Xcoord + i * Ycoord$ ;  $z(n+1) = z(0)*fn(z(n))$ . Parameters are perturbations of  $z(0)$ . 'fn' is a variable function.

**mandel4** = Fourth-power 'Mandelbrot' fractals using 32-bit integer math.  $z(0) = 0$ ;  $z(n+1) = z(n)^4 + C$ , where  $C = Xcoord + i * Ycoord$ . Two optional params: real and imaginary parts of  $z(0)$  (if not 0).

**julia4** = Fourth-power Julia set fractals using 32-bit integer math.  $z(0) = Xcoord + i * Ycoord$ ;  $z(n+1) = z(n)^4 + C$ . Two params required: real and imaginary parts of C.

**test** = 'test' point letting us (and you!) easily add fractal types. Currently, the 'Distance Estimator' M'brot/Julia Set algorithm. two optional parameters - if none given, uses the M'brot Set"

**plasma** = plasma clouds - random, cloud-like formations. Requires four or more colors. One param: 'graininess' (.5 to 50, default = 2)

**julfn+zsqr**= Julia Biomorph fractal.  $z(0) = Xcoord + i * Ycoord$ ;  $z(n+1) = fn(z(n)) + z(n)^2 + C$ . Two params: real, imag portions of C.

**manfn+zsqr**= 'Mandelbrot-Equivalent' for the Julfn+zsqr fractal.  $z(0) = Xcoord + i * Ycoord$ ;  $z(n+1) = fn(z(n)) + z(n)^2 + (Xcoord + i * Ycoord)$ . Parameters are perturbations of  $z(0)$ .

**manzpower** = 'Mandelbrot-Equivalent' for julzpower.  $z(n+1) = z(n)^m + C$ . Parameters are real and imaginary perturbations, exponent m.

**julzpower** = Juliazpower fractal.  $z(0) = Xcoord + i * Ycoord$ ;  $z(n+1) = z(n)^m + C$ . Three params: real, imag portions of C, exponent m.

**manzppwr** = 'Mandelbrot-Equivalent' for the julzppwr fractal. Use the Space-bar to select julzppwr fractals a/la Mandel/Julia.  $z(n+1) = z(n)^z(n) + z(n)^m + C$ . Parameters are real perturbation, imaginary perturbation, and exponent m.

**julzppwr** =  $z(0) = Xcoord + i * Ycoord$ ;  $z(n+1) = z(n)^z(n) + z(n)^m + C$ . Three params: real, imag portions C, and the exponent m.

**manfn+exp** = 'Mandelbrot-Equivalent' for the julfn+exp fractal. Use the Space-bar to select julfn+exp fractals a/la Mandel/Julia.  $z(n+1) = fn(z(n)) + e^z(n) + C$ . Parameters are real perturbation, and imaginary perturbation of  $z(0)$ .

**julfn+exp** = julia fn+exp fractal.  $z(0) = Xcoord + i * Ycoord$ ;  $z(n+1) = fn(z(n)) + e^z(n) + C$ . Two params: real, imag portions C.

**popcorn** = orbits of  $x(n+1) = x(n) - h*\sin(y(n) + \tan(3*y(n)))$  and  $y(n+1) = y(n) - h*\sin(x(n) + \tan(3*x(n)))$  plotted for EACH screen pixel and superimposed.

**popcornjul** = Julia set from popcorn formula.  $z(0) = Xcoord + i * Ycoord$ . Orbit calculated as in popcorn.

**bifurcation** = 'Bifurcation' fractal. Pictoral representation of a population growth model. Let P = new population, p = oldpopulation,

$r = \text{growth rate. The model is: } P = p + r * p * (1 - p).$   
**biflambda** = Bifurcation variation: model is:  $P = p + r * p * (1 - p)P.$   
**bif+sinpi** = Bifurcation variation: model is:  $P = p + r * \sin(\pi * p).$   
**bif=sinpi** = Bifurcation variation: model is:  $P = p + r * \sin(\pi * p).$   
**lorenz** = Lorenz attractor - orbit in three dimensions defined by:  
 $x_{\text{new}} = x + (-a * x * dt) + (a * y * dt)$   
 $y_{\text{new}} = y + (b * x * dt) - (y * dt) - (z * x * dt)$   
 $z_{\text{new}} = z + (-c * z * dt) + (x * y * dt)$   
Parameters are dt, a, b, and c.  
**lorenz3d** = 3D Lorenz attractor with 3D perspective.  
**rossler3D** = Orbit in three dimensions defined by:  
 $x_{\text{new}} = x - y * dt - z * dt$   
 $y_{\text{new}} = y + x * dt + a * y * dt$   
 $z_{\text{new}} = z + b * dt + x * z * dt - c * z * dt$   
Parameters are dt, a, b, and c.  
**henon** = Orbit in two dimensions defined by:  
 $x_{\text{new}} = 1 + y - a * x * x$   
 $y_{\text{new}} = b * x$   
Parameters are a and b  
**pickover** = Orbit in three dimensions defined by:  
 $x_{\text{new}} = \sin(a * y) - z * \cos(b * x)$   
 $y_{\text{new}} = z * \sin(c * x) - \cos(d * y)$   
 $z_{\text{new}} = \sin(x)$   
Parameters are a, b, c, and d.  
**gingerbread** = Orbit in two dimensions defined by:  
 $x <- 1 - y + |x|$   
 $y <- x$   
**diffusion** = Diffusion Limited Aggregation. Randomly moving points accumulate. One param: border width (default 10)  
**fn(z\*z)** =  $z(0) = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = \text{fn}(z(n)) * z(n)$   
**fn\*fn** =  $z(0) = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = \text{fn}(n) * \text{fn}(n)$   
**fn\*z+z** =  $z(0) = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = p1 * \text{fn}(z(n)) + p2 * z(n)$   
Parameters are real and imaginary components of p1 and p2.  
**fn+fn** =  $z(0) = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = p1 * \text{fn1}(z(n)) + p2 * \text{fn2}(z(n))$   
Parameters are real and imaginary components of p1 and p2.  
**sqr(1/fn)** =  $z(0) = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = (1/\text{fn}(z(n))) * (1/\text{fn}(z(n)))$   
**sqr(fn)** =  $z(0) = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = \text{fn}(z(n)) * \text{fn}(z(n))$   
**spider** =  $c(0) = z(0) = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = z(n) * z(n) + c(n);$   
 $c(n+1) = c(n)/2 + z(n+1)$   
**manowar** =  $c = z1(0) = z(0) = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = z(n)^2 + z1(n) + c;$   
 $z1(n+1) = z(n);$   
**tetrate** =  $z(0) = c = X_{\text{coord}} + i * Y_{\text{coord}}; z(n+1) = c^z(n)$   
**kamtorus** = series of orbits superimposed.  $x(0) = y(0) = \text{orbit}/3;$   
 $x(n+1) = x(n) * \cos(a) + (x(n) * x(n) - y(n)) * \sin(a)$   
 $y(n+1) = x(n) * \sin(a) - (x(n) * x(n) - y(n)) * \cos(a)$   
After each orbit, 'orbit' is incremented by a step size. Parameters are a, step size, stop value for 'orbit', and points per orbit.  
**kamtorus3d** = Same as kamtorus but in 3D with 'orbit' the z dimension.  
**magnetj1** =  $z(0) = X_{\text{coord}} + i * Y_{\text{coord}};$   

$$z(n+1) = \left| \frac{z(n)^2 + (c-1)}{2 * z(n) + (c-2)} \right|^2$$
  
Parameters are real and imaginary parts of c.  
**magnetm1** = 'Mandelbrot' variant with  $c = X_{\text{coord}} + i * Y_{\text{coord}}$  and  $z(0) = 0.$   
Orbit formula same as magnetj1

Parameters are perturbations of  $z(0)$ .

**magnetj2** =  $z(0) = X_{\text{coord}} + i * Y_{\text{coord}}$ ;  

$$z(n+1) = \sqrt{\frac{z(n)^3 + 3*(C-1)*z(n) + (C-1)*(C-2)}{3*(z(n)^2) + 3*(C-2)*z(n) + (C-1)*(C-2) - 1}}$$

Parameters are real and imaginary parts of  $c$ .

**magnetm2** = 'Mandelbrot' variant with  $c = X_{\text{coord}} + i * Y_{\text{coord}}$  and  $z(0) = 0$ .  
 Orbit formula same as magnetj2  
 Parameters are perturbations of  $z(0)$ .



## **Image Sizing and Color selection**

Using the Options **Image Size** option, you can select either a pre-defined image size, or make up your own (up to an internal limit of 2048x2048) and select a two-color, 16-color, or 256-color image. Those of you with 24-bit color video cards will have to hold yourself back to 256-color images for the moment - sorry.

If you have selected the "Window Sizing" option, your image window will automatically re-size to the image size you have just selected.

If your image size is larger than your window, you can use the scroll bars in the window to scroll around your image.

## **Zooming in on an Image, Mandelbrot/Julia Toggling**

With this release of Winfract, there are two different methods of generating and using "Zoom-Boxes".

If you do not have Mark Peterson's new "Zoom-Box" option enabled, Winfract's zoom box is activated in the same fashion as it was in earlier releases:

To zoom in on a fractal image, just move your mouse pointer to the upper-left corner of the "zoom-box" area, push and hold down the left mouse-button, move the mouse pointer to the lower-right corner of the "zoom-box" area, and let go of the mouse button. The program will immediately begin a new image using the "zoom-box" coordinates you selected.

As you are moving your mouse with the left-button down, the "zoom-box" area will be displayed as an X-ORed rectangle. If you change your mind about zooming as you are moving the mouse around ("No, No, Not \*There\*!"), just make one of the rectangular coordinates less than five pixels across - the program will "forget" about the zoom-box if you let go of the left mouse button and that is the case.

If the aspect ratio of your "zoom-box" is reasonably close to that of the full image size (within 33 percent or so), the program adjusts the "zoom-box" to match the aspect ratio exactly before performing the zoom.

If you have enabled Mark Peterson's new "Zoom-Box" option, a vertical Zoom-Box scroll-bar is displayed in addition to your fractal image. The Zoom-Box defaults to the mid-position, in which your Zoom-Box exactly covers the area of your fractal image. Moving the scroll-button up shrinks your zoom-box (and makes it visible on the fractal image). You can then move the zoom-box around by moving your cursor inside the zoom-box and then holding down your left mouse-button and moving the cursor and zoom-box as a single unit. Double-clicking the left mouse button will cause your image to be redrawn using your current zoom-box coordinates.

You can also "Zoom Out" by moving the zoom-box scroll bar below the midpoint. When you do this, the zoom-box that is displayed is actually showing you where (and how small) your currently displayed image would be if you double-clicked the left mouse button at that point. Moving the zoom-box and double-clicking to perform the zoom is done the same way as when you are "zooming in".

You can switch from a "Mandelbrot Set" image to its "Julia Set" at the location of the mouse cursor by clicking on the right mouse button. When the corresponding "Julia Set" image is on the screen, click the right mouse button again to get the "Mandelbrot Set" image back. The terms "Mandelbrot Set" and "Julia Set" are in quotes because many fractal types (mandel4 and julia4, for instance) have this "Mandelbrot Set"/"Julia Set" relationship.

## **Options and Doodads**

You can select and of a number of options and doodads using the "Options Options-and-Doodads" menu. The list of doodads currently includes various algorithms (single/dual pass, solid-guessing, boundary tracing), biomorphs, decomposition, inside color, outside color, maximum iterations, etc - all flagrantly stolen from Fractint-for-DOS.

Once you select any of these options, the program will begin re-drawing your image with the new options in effect.

## Color Cycling

You start and stop Color-cycling mode using the Color-Cycling menu. This menu also lets you determine the direction of the cycling (the designations "in" and "out" were arbitrarily picked based on how the effect looked on a single Mandelbrot image), whether to just rotate the existing colors or generate new ones randomly, and (for random color-generation) whether the colors are to change with a low, medium, or high frequency.

Sorry, but at the current time color-cycling is restricted to display devices whose Windows drivers are capable of palette manipulation - and the "stock" VGA driver distributed with Win 3.0 doesn't have it! Also, color-cycling affects the background windows as well unless your video driver is capable of displaying 256 or more simultaneous colors. If Windows finds itself running with a 16-color video driver, it color-cycles only when its window has the input focus, and returns to the "stock" palette values when it loses this focus.

Winfract can now also use "**Hot-keys**" for color-cycling

- spacebar - toggles color-cycling on and off

- left, right arrows - turn on color-cycling and set the "direction"

- ENTER - initiate random color-cycling with all new colors

(uhh, note that Windows takes a second or so to "fire up" color-cycling, so be patient when you press one of these keys. Breathe in, breathe out, ...)

## Getting a Status Report

You can find out the status of your current image (which formula you are using, its parameters and screen corners, and - most important - whether or not it's still being generated or has finished) by selecting the **Status** option from the Options menu. A "Pup-up" message will tell you about the status of the current image. If the image is still being generated, the program will continue with it when you press the "OK" button.

## **Distribution Policy and Procedures, Contacting the Authors**

The Stone Soup Group is a loosely associated group of fanatic programmers. See Fractint-for-DOS's FRACTINT.DOC for the story behind the 'Stone Soup' name. Fractint for Windows, like Fractint-for-DOS, is freeware, and our "contribution policy" is the same: **Don't want money. Got money. Want admiration.**

### **Conditions on use:**

**Fractint for Windows, like Fractint-for-DOS, may be freely copied and distributed but may not be sold.** It may be used personally or in a business - if you can do your job better by using Fractint for Windows, or use images from it, that's great! It may be given away with commercial products under the following conditions: It must be clearly stated that Fractint for Windows does not belong to the vendor and is included as a free give-away. It must be a complete unmodified release of Fractint, with documentation. There is no warranty of Fractint's suitability for any purpose, nor any acceptance of liability, express or implied.

Virtually all of the Fractint authors can be found on the **Compuserve** network in the **COMART** ('COMputer ART') forum in S 15 ('Fractals'). Fractint development occurs in the COMART forum - most of the authors have never met except there. New members are always welcome! Several of us can also be found on **BIX** in the **GRAPHIC.DISP/FRACTALS** area.

New versions of Fractint-for-DOS and Fractint-for-Windows are uploaded to the Compuserve and BIX networks, and make their way to other systems from those points. The latest version of Fractint-for-Windows can usually be found in the following locations:

**WINFRA.ZIP** - (Executable/Docs) Compuserve: COMART Lib 15 ("Fractals")

BIX: GRAPHIC.DISP/LISTINGS

**WINSRC.ZIP** - (Complete Source) Compuserve: COMART Lib 15 ("Fractals")

BIX: GRAPHIC.DISP/LISTINGS

(What's the latest version? Well, THIS one was, way back when we uploaded it!)

## A list of Fractint Authors

Fractint for Windows was originally ported from Fractint-for-DOS by Bert Tyler. This is the first Windows program that Bert ever wrote, which goes a long way towards explaining a lot of the bugs. Bert's task was made a lot easier by Pieter Branderhorst, who separated the MSDOS-specific code from Fractint-for-DOS's fractal generator modules, making a Windows port of the package possible. Mark Peterson helped track down and eliminate a lot of the bugs that were in the original program, and has since added many of the new options which take advantage of the capabilities of the Windows environment.

Fractint for Windows is based heavily on (and uses the fractal generator engines straight out of) Fractint-for-DOS. A partial list of the authors of Fractint-for-DOS includes:

----- Primary Authors (this changes over time) -----  
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----- Contributing Authors -----

Michael Abrash - 360x480x256, 320x400x256 VGA video modes  
Joseph Albrecht - Tandy video, CGA video speedup  
Kevin Allen - Finite attractor and bifurcation engine  
Steve Bennett - restore-from-disk logic  
Rob Beyer - [71021,2074] Barnsley IFS, Lorenz fractals  
Mike Burkey - 376x564x256, 400x564x256, and 832x612x256 VGA video modes  
John Bridges - [73307,606] superVGA support, 360x480x256 mode  
Brian Corbino - [71611,702] Tandy 1000 640x200x16 video mode  
Lee Crocker - [73407,2030] Fast Newton, Inversion, Decomposition..  
Monte Davis - [71450,3542] Documentation  
Chuck Ebbert - [76306,1226] compressed, sqrt log palette, FPU speedups  
Richard Finegold - [76701,153] 8/16/..256-Way Decomposition option  
Mike Gelvin - [73337,520] Mandelbrot speedups  
Lawrence Gozum - [73437,2372] Tseng 640x400x256 Video Mode  
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Norman Hills - [71621,1352] Ranges Option  
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Adrian Mariano - [theorem@blake.acs.washington.edu] Diffusion & L-Systems  
Charles Marslett - [73317,3662] VESA Video and ITT math chip support  
Chris Martin - Paintjet printer support  
Joe McLain - [75066,1257] TARGA Support, color-map files  
Bob Montgomery - [73357,3140] (Author of VPIC) Fast text I/O routines  
Bret Mulvey - plasma clouds  
Ethan Nagel - [70022,2552] Palette editor, integrated Help/Doc system  
Jonathan Osuch - [73227,1432] ITT detect  
Marc Reinig - [72410,77] Lots of 3D options  
Kyle Powell - [76704,12] 8514/A Support  
Matt Saucier - [72371,3101] Printer Support  
Herb Savage - [71640,455] 'inside=bof60', 'inside=bof61' options  
Lee Skinner - Tetrade, Spider, Mandelglass fractal types and more  
Dean Souleles - [75115,1671] Hercules Support  
Kurt Sowa - [73467,2013] Color Printer Support  
Hugh Steele - Cyclorange feature  
Scott Taylor - [72401,410] (DGWM18A) Postscript, KAM Torus, many fn types  
Paul Varner - [73237,411] Floating-point fractal algorithms

Dave Warker - Integer Mandelbrot Fractals concept  
Phil Wilson - [76247,3145] Distance Estimator, Bifurcation fractals  
Richard Wilton - Tweaked VGA Video modes  
...  
Byte Magazine - Tweaked VGA Modes  
MS-Kermit - Keyboard Routines  
PC Magazine - Sound Routines  
PC Tech Journal - CPU, FPU Detectors





## Coordinate Box options

The **Coordinate Box** is courtesy of Mark Peterson. When you have the Coordinate Box checked, a small "coordinate window" constantly displays the current position of your mouse pointer.

From the **Options** menu you can select the coordinates to display in rectangular (default), polar, or pixel coordinates.

**Rectangular** coordinates correspond to Cartesian plane. The coordinates displayed are in absolute units relative to the origin. Fractint uses these coordinates to form a complex number. This complex number initializes one or more variables in the iterative calculation. The  $x$  coordinate is used as the *real* portion of the complex number and the  $y$  coordinate as the *imaginary* portion.

**Pixel** coordinates display the position of a point in terms of the number of pixels, or color dots, relative to the pixel in the upper left-hand corner of the image. For example, if the image size is 200 x 150, then the pixel in the lower right-hand corner of the image is coordinate (199, 149).

**Polar** coordinates display the position of a point in terms of its distance and angle relative to the origin. The angle can be in units of **degrees** (default), **radians**, or **grads**. Most people are familiar with degrees which divide the circle into 360 *degrees*. Grads divide the circle into 400 *grads*. Radians divide the circle into units of  $2\pi$  *radians*.