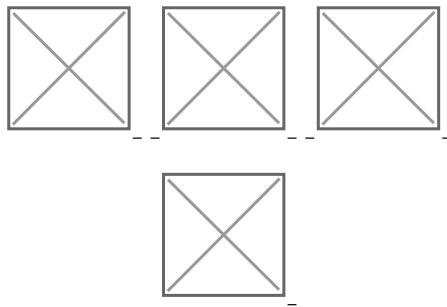


Alpaca Electron



Alpaca Electron



*Alpaca Electron is built from the ground-up to be the easiest way to chat with the alpaca AI models.
No command line or compiling needed!*

☐ Features + to-do

- ☑ Runs locally on your computer, internet connection is not needed except when downloading models
- ☑ Compact and efficient since it uses [llama.cpp](#) as its backend (which supports Alpaca & Vicuna too)
- ☑ Runs on CPU, anyone can run it without an expensive graphics card
- ☑ No external dependencies required, everything is included in the installer
- ☑ "Borrowed" UI from *that* popular chat AI :trollface:
- ☑ Supports Windows, MacOS, and Linux (untested)
- ☑ Docker-ized ☐
- ☑ Context memory

- ☐ Chat history
- ☐ Integration with Stable Diffusion
- ☐ DuckDuckGo integration for web access
- ☐ GPU acceleration (cuBLAS & openBLAS)

Demo

Demonstration

Quick Start Guide

1. Download an Alpaca model (7B native is recommended) and place it somewhere on your computer where it's easy to find.

Note

Download links will not be provided in this repository.

2. Download the latest installer from the [releases page](#) section.
3. Open the installer and wait for it to install.
4. Once done installing, it'll ask for a valid path to a model. Now, go to where you placed the model, hold shift, right click on the file, and then click on "Copy as Path". Then, paste this into that dialog box and click .
5. The program will automatically restart. Now you can begin chatting!

Note

The program will also accept any other 4 bit quantized .bin model files. If you can find other .bin Alpaca model files, you can use them instead of the one recommended in the Quick Start Guide to experiment with different models. As always, be careful about what you download from the internet.

Troubleshooting

General

- If you get an error that says "Invalid file path" when pasting the path to the model file, you probably have some sort of misspelling in there. Try copying the path again or using the file picker.
- If you get an error that says "Couldn't load model", your model is probably corrupted or incompatible. Try downloading the model again.
- If you face other problems or issues not listed here, create an issue in the "Issues" tab at the top of this page. Describe in detail what happens, and include screenshots.

Windows

- If the model has been loaded into RAM but text generation doesn't seem start, check to see if your CPU is compatible with the AVX2 instruction set. If it does not support AVX2, Alpaca Electron will use AVX instead, which is much slower so be patient.
- If you get an error saying "vcruntime140_1.dll is missing" or nothing happens at all and the model was not loaded into RAM, try installing the Microsoft Visual C++ Redistributable.

MacOS

- If you get an error that says "App can't be opened because it is from an unidentified developer.", go to the Applications folder. Then, hold the control key and click on the app. Then click "Open", then click "Open" when it gives you a warning. Your preference will be saved and MacOS will let you open the app normally from now on.
- If the above method does not work, try running the following command in terminal: `xattr -cr /Applications/Alpaca\ Electron.app/`

Linux

- You can either download the prebuilt app (packaged as tar.gz) from the releases page, extract it and execute it with `./"alpaca electron"` or build the application on yourself.
- If you want to build the application yourself:

“ Clone the repository:

```
git clone https://github.com/ltsPi3141/alpaca-electron.git
```

Change your current directory to alpaca-electron:

```
cd alpaca-electron
```

Install application specific dependencies:

```
npm install --save-dev
```

Build the application:

```
npm run linux-x64
```

Change your current directory to the build target:

```
cd release-builds/'Alpaca Electron-linux-x64'
```

Run the application with `./'Alpaca Electron'`

Docker Compose

- You can run this electron application with docker compose. Therefore you need to complete the following steps:

“ Clone the repository:

```
git clone https://github.com/ltsPi3141/alpaca-electron.git
```

Change your current directory to alpaca-electron:

```
cd alpaca-electron
```

Build the container image:

```
docker compose build
```

Run the application container:

```
docker compose up -d
```

- If no window opens up run `docker compose up` (without the -d). If there is an error like `Authorization required, but no authorization protocol specified` run `xhost local:root` on your docker host.

🔧 Building

Prerequisites

- [Node.js](#)
- [Git](#)
- If you're on Windows and are planning on building llama.cpp binaries also, [CMake](#).

(OPTIONAL) Building llama.cpp from source

1. Clone llama.cpp's GitHub repo

```
git clone https://github.com/ggerganov/llama.cpp  
cd llama.cpp
```

2. Build llama.cpp On Windows:

```
mkdir build  
cd build  
cmake ..  
cmake . --config Release
```

On Linux and MacOS:

```
make
```

Running the project from source

1. Clone the GitHub repo

```
git clone https://github.com/ItsPi3141/alpaca-electron  
cd alpaca-electron
```

2. Install node packages

```
npm install
npm run rebuild
```

“ **Info** If you are on Linux, replace `npm run rebuild` with `npm run rebuild-linux`

3. (OPTIONAL) Use your own llama.cpp build

“ **Warning**

This step is not required. Only do it if you had built llama.cpp yourself and you want to use that build. Otherwise, skip to **step 4** If you had built llama.cpp in the previous section, copy the `main` executable file into the `bin` folder inside the alpaca-electron folder.

Make sure the file replaces the correct file. E.g. if you're on Windows, replace chat.exe with your file. If you're on arm64 MacOS, replace chat_mac_arm64. Etc...

4. Start the Electron app

```
npm start
```

Building a release and installer

Run one of the following commands:

- `npm run win`
- `npm run mac-x64`
- `npm run mac-arm64`
- `npm run linux-x64`

You can only build for the OS you are running the build on. E.g. if you are on Windows, you can build for Windows, but not for MacOS and Linux.

Credits

Credits go to [@antimatter15](#) for creating alpaca.cpp and to [@ggerganov](#) for creating llama.cpp, the backbones behind alpaca.cpp. Finally, credits go to Meta and Stanford for creating the LLaMA and Alpaca models, respectively.

Special thanks to [@keldenl](#) for providing arm64 builds for MacOS and [@W48B1T](#) for providing Linux builds

Revision #2

Created 21 June 2023 09:19:03 by tinfoil-hat

Updated 21 June 2023 09:20:01 by tinfoil-hat