

# **Building EMERALD-UI**

## **a protein alignment visualization tool for EMERALD**

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Project: [emerald-ui](#)

# **Presentation Structure**

## **What We'll Cover Today**

- 1. Background & Motivation**
- 2. Limitations of Traditional Methods**
- 3. EMERALD Algorithm**
- 4. EMERALD UI Overview**
- 5. Technical Implementation**
- 6. EMERALD-UI In Action**

# Background & Motivation

## Why Protein Alignment Matters

Proteins are the workhorses of biology, and comparing their sequences helps us understand function, evolution, and disease.

### Example Protein Alignment:

```
A: MKTAYIAKQRQISFVKSHFSRQ  
B: MKTA--AKQRQISFVKSHFSRQ
```

Match: +5    Mismatch: -2    Gap: -6

# How Are Proteins Aligned?

- **Sequence alignment algorithms** compare amino acid sequences to find similarities.
- **Cost matrices** (e.g., BLOSUM, PAM) assign scores for matches, mismatches, and gaps.
- The goal: **maximize total score** by aligning similar regions and penalizing differences.

	C	S	T	A	G	P	D	E	Q	N	H	R	K	M	I	L	V	W	Y	F	
C	9																				C
S	-1	4																			S
T	-1	1	5																		T
A	0	1	0	4																	A
G	-3	0	-2	0	6																G
P	-3	-1	-1	-1	-2	7															P
D	-3	0	-1	-2	-1	-1	6														D
E	-4	0	-1	-1	-2	-1	2	5													E
Q	-3	0	-1	-1	-2	-1	0	2	5												Q
N	-3	1	0	-2	0	-2	1	0	0	6											N
H	-3	-1	-2	-2	-2	-2	-1	0	0	1	8										H
R	-3	-1	-1	-1	-2	-2	-2	0	1	0	0	5									R
K	-3	0	-1	-1	-2	-1	-1	1	1	0	-1	2	5								K
M	-1	-1	-1	-1	-3	-2	-3	-2	0	-2	-2	-1	-1	5							M
I	-1	-2	-1	-1	-4	-3	-3	-3	-3	-3	-3	-3	-3	1	4						I
L	-1	-2	-1	-1	-4	-3	-4	-3	-2	-3	-3	-2	-2	2	2	4					L

# Issues in Protein Alignment

- **Multiple Valid Alignments:** Different alignments may be equally plausible.
- **Gap Placement:** Deciding where to insert gaps affects biological interpretation.
- **Scoring Bias:** Choice of cost matrix influences which regions align.
- **Functional Relevance:** Not all aligned regions are biologically meaningful.

Instead of finding **one "optimal" alignment**, what if we could:

- **Explore the alignment landscape** - Find all good alignments, not just the best
- **Identify trustworthy regions** - Discover where alignments consistently agree
- **Reduce scoring bias** - See robust patterns across different parameters
- **Focus on functional relevance** - Highlight biologically meaningful regions

## The EMERALD Approach:

“ **If multiple high-scoring alignments agree on a region,  
that region is probably functionally important!**

”

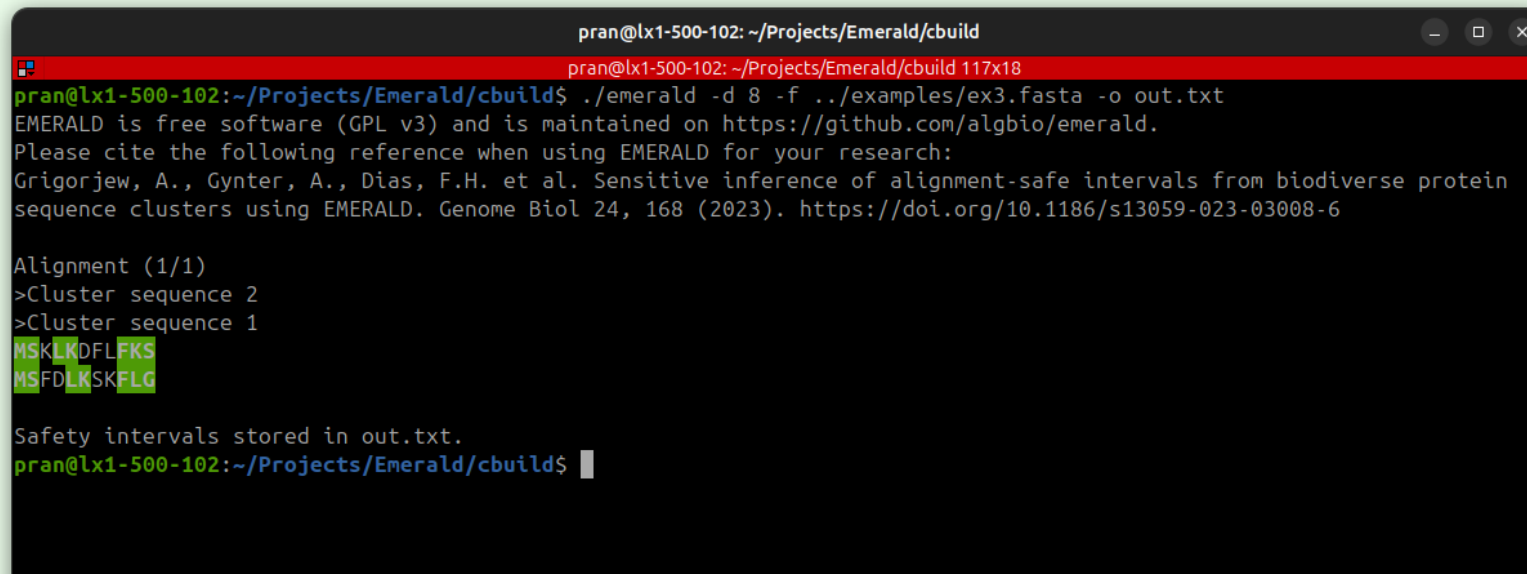
This is exactly what **EMERALD** does - it reveals which parts of your alignment you can trust.

# EMERALD by itself

**EMERALD** is a terminal application for generating protein alignments written in C++.

## Issues

- App only gives raw data, which is hard to interpret
- Only work in the terminal, which is hard to use
- Limited input methods



```
pran@lx1-500-102: ~/Projects/Emerald/cbuild
pran@lx1-500-102: ~/Projects/Emerald/cbuild 117x18
pran@lx1-500-102:~/Projects/Emerald/cbuild$ ./emerald -d 8 -f ../examples/ex3.fasta -o out.txt
EMERALD is free software (GPL v3) and is maintained on https://github.com/algbio/emerald.
Please cite the following reference when using EMERALD for your research:
Grigorjew, A., Gynter, A., Dias, F.H. et al. Sensitive inference of alignment-safe intervals from biodiverse protein
sequence clusters using EMERALD. Genome Biol 24, 168 (2023). https://doi.org/10.1186/s13059-023-03008-6

Alignment (1/1)
>Cluster sequence 2
>Cluster sequence 1
MSKLDLFLFKS
MSFDLKSFLG

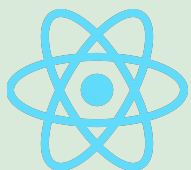
Safety intervals stored in out.txt.
pran@lx1-500-102:~/Projects/Emerald/cbuild$
```

# How EMERALD-UI Solves These Problems

- **Intuitive Visualization:** Draws an interactive, easy-to-use graph from raw EMERALD output.
- **Accessible Interface:** Runs in the browser, no terminal required.
- **Flexible Input:** Paste sequences, upload files, or search directly from UniProt.
- **Alignment Confidence:** Highlights common areas of multiple alignments.
- **Parameter Exploration:** Instantly adjust scoring and see how alignments change.
- **Safety Windows Visualization:** Highlights common regions in 3D structures of proteins.



# Technology Stack



## React

Modern UI framework for building interactive web applications.



## TypeScript

Type-safe code for reliability and maintainability.



## WebAssembly (WASM)

Runs EMERALD executable in-browser.



## D3.js

Interactive, data-driven visualizations.



## UniProt API

Direct protein sequence search.



## Mol\*

3D protein structure visualization.

# **Demo Time**

**Let's see EMERALD-UI in action!**

# **Thank You!**

## **Questions?**

Feel free to ask about EMERALD-UI, protein alignment, or anything from today's presentation.