# **A Visual Encoding System for Comparative Exploration** of Magnetic Resonance Spectroscopy Data

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INTRODUCTION

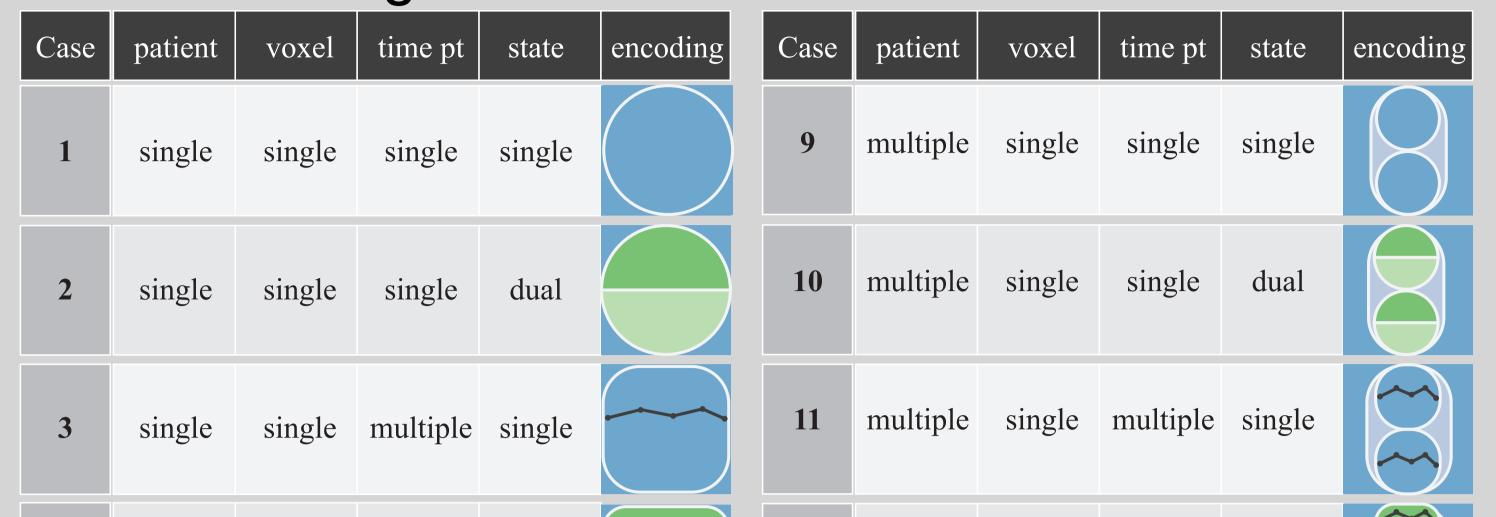
What is MRS? Magnetic resonance spectroscopy (MRS) is an in vivo non-invasive biochemical imaging technique utilized for tissue metabolite characterization used often for early detection and treatment of tumors and neuropsychiatric conditions. <sup>2,3</sup>

Presenting a novel visualization system for the exploratory analysis of MRS data

Why is this data difficult to visualize? MRS data output is inherently abstract, with a steep learning curve to successful interpretation. Data often have poor signal to noise ratio, making comparison difficult. 2,3

## RESULTS

### Visual Encodings

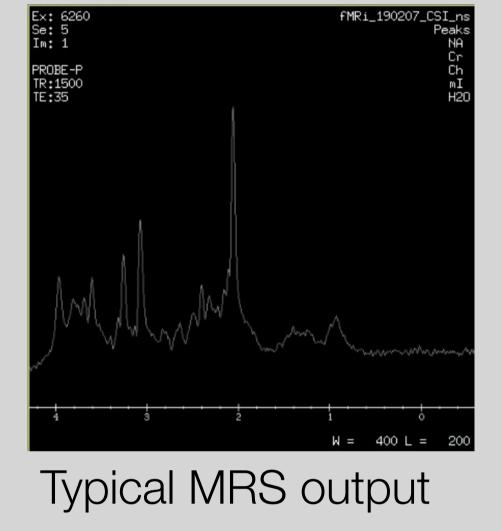


What are the current toolset limitations?

- Qualitative comparison between voxel samples is difficult or impossible
- Limited capabilities to compare all permutations of spectral metabolite ratios<sup>1</sup>
- Limited tools for correlating spectral peaks to corresponding metabolites and spatial sample location

# **DATA & TASKS**

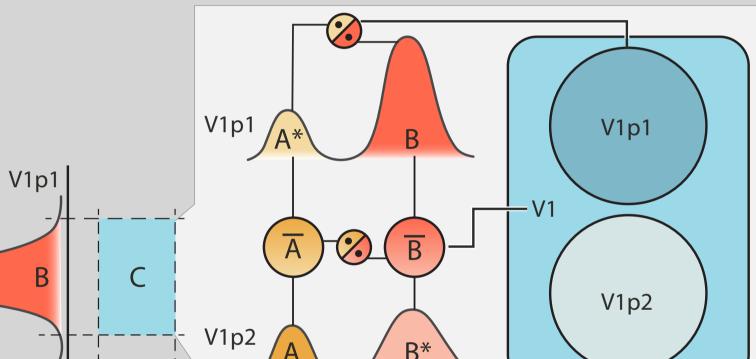
Tier 1	Spectral metabolite samples	Ex: 6260 fMF Se: 5 Im: 1 PROBE-P TR:1500 TE:35
Tier 2	Key attributes: Time, Patient, State, Voxel	MM
Tier 3	Hidden but retained attributes	Typical MRS out



4	single	single	multiple	dual	12	multiple	single	multiple	dual	
5	single	multiple	single	single	13	multiple	multiple	single	single	
6	single	multiple	single	dual	14	multiple	multiple	single	dual	
7	single	multiple	multiple	single	15	multiple	multiple	multiple	single	
8	single	multiple	multiple	dual	16	multiple	multiple	multiple	dual	

Glyphs encoded to each Tier 2 attribute (individual, voxel, time point, state) permute to 16 possible MRS case scenarios.

Calculations We perform 2 different metabolite ratio calculations based on peak integrals, one for the overview layer and one for the detail layer.

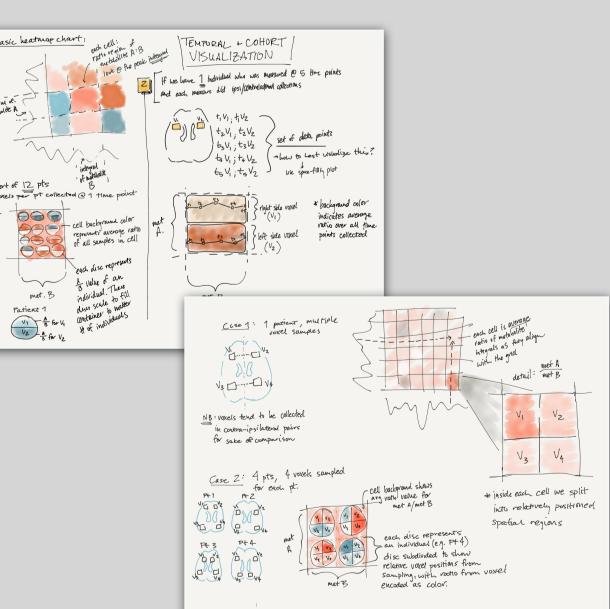


### Core Tasks

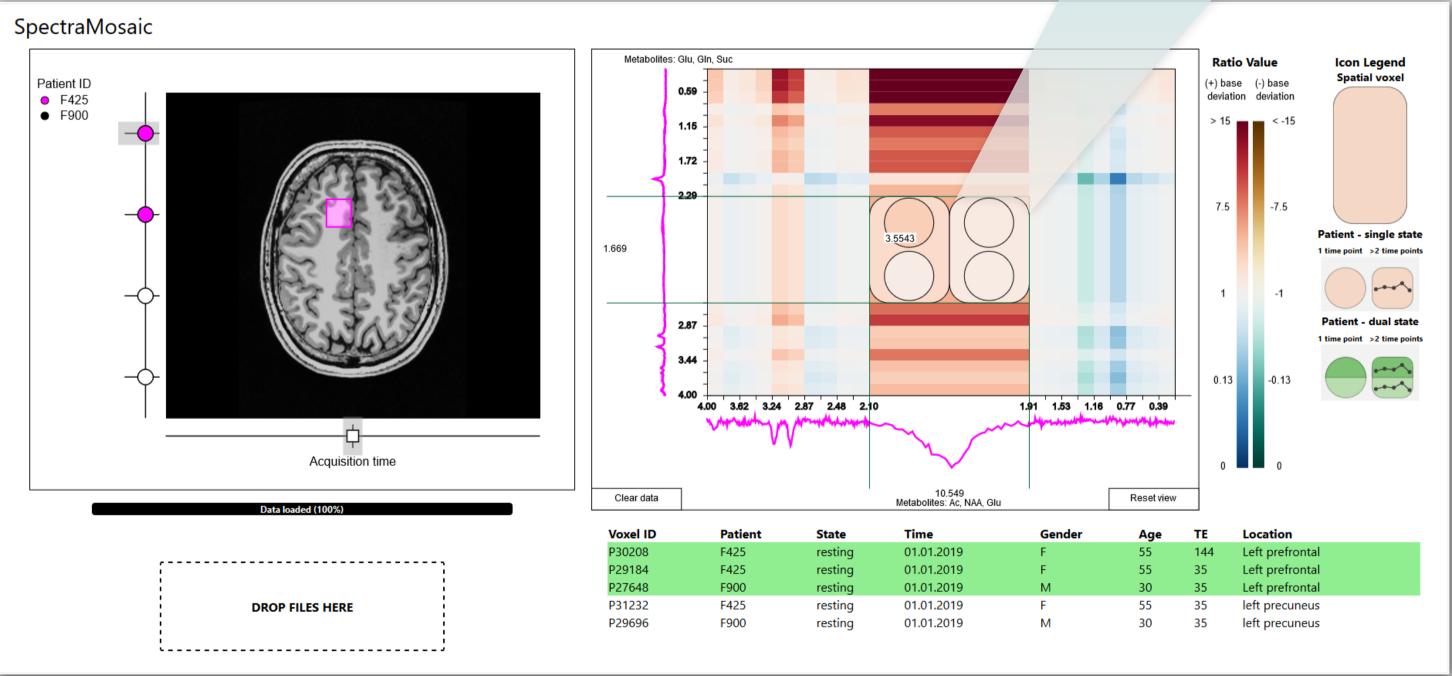
- 1. Correlate spectral peaks with corresponding metabolites
- Match voxel sample to spectral output 2.
- Determine the ratio of metabolites in a sample(s) 3. by calculating the ratio of their peak integrals
- 4. Select a subset of spectral acquisitions from a single group for further analysis
- 5. Discover outliers in spectral output
- Compare spectral metabolite concentrations within and across voxel samples

## Design Requirements

- Web application
- 2. Spectral peak identification
- Linking between data sources 3.
- 4. Support comparison of spectral metabolite concentrations as ratios
- 5. Layered design: overview vs. detail

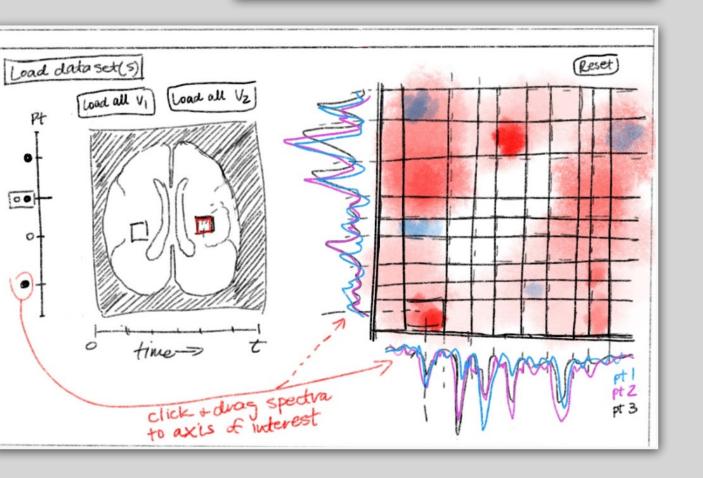


# **SpectraMosaic Application**



We realized this visual encoding system as part of an interactive insight-generation tool for rapid exploration and comparison of metabolite ratio variation for deeper insights to these complex data.

We then developed a taxonomy of visual encodings to represent the range of different metabolite **concentration ratios** at different dimensional tiers.



### CONCLUSIONS

We have developed a novel visual encoding system for selected MRS data attributes using simple glyph shapes with diverging color maps to represent variation. We have implemented this system in a spectral visual analysis tool, and plan to continue working with domain experts to extend our visual system to manage groups and uncertainty.

### REFERENCES

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- 2. POSSE S., OTAZO R., DAGER S. R., ALGER J.: MR spectroscopic imaging: Principles and recent advances. Journal of Magnetic Resonance Imaging 37 (2013), 1301–1325. doi:10.1002/jmri. 23945. 1 3. VAN DER GRAAF M.: In vivo magnetic resonance spectroscopy: Basic methodology and clinical applications. Euro- pean Biophysics Journal 39, 4 (2010), 527-540. doi:10.1002/9780470882221. 1





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