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#### **CoreMS Framework Overview**

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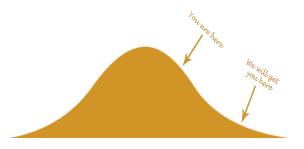
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#### **CoreMS Framework Overview**

- Modern, flexible, stable and sustainable software framework and platform for mass spectra data-processing workflows, including hyphened methods, i.e., LC-MS, IM-MS targeting small molecules analysis of complex mixtures
- Comprehensive molecular formula assignment algorithm merging state of the art approaches and universal mass spectrometry features
- Robust confidence metric for formula assignment enables analysis validation, comparison of algorithms results and ensures analytical reproducibility









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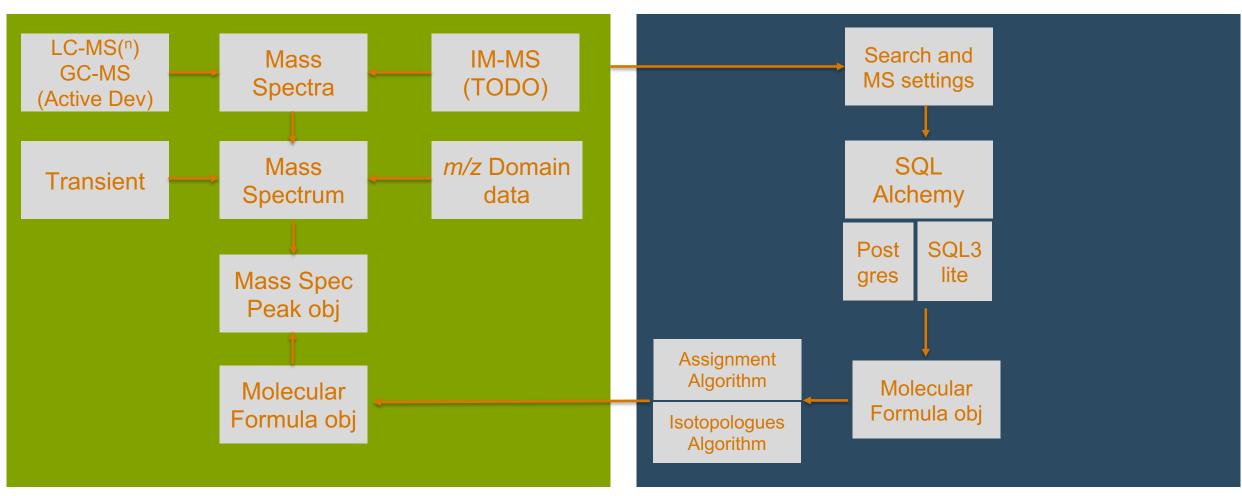
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## **Software Framework**

#### **CoreMS** Hierarchical Data Structure

#### **CoreMS** Molecular Formula Assignment Workflow



For available input, output and data structure types please refer to: <a href="https://github.com/EMSL-Computing/CoreMS/blob/master/README.md">https://github.com/EMSL-Computing/CoreMS/blob/master/README.md</a>



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## **Software Framework**

#### CoreMS Basic Molecular Formula Assignment Script

```
file_name = 'neg_esi_srfa_1ppm_test.d'
                              bruker reader = ReadBrukerSolarix(file location)
                     Input
                               bruker_transient_obj = bruker reader.get transient()
                               T = bruker_transient_obj.transient_time
                               mass_spectrum_obj = bruker_transient.get_mass_spectrum(plot_result=False, auto_process=True)
      Signal processing
Molecular Identification
                              SearchMolecularFormulas(first hit=False).run worker mass spectrum(mass_spectrum_obj)
                               for mspeak in mass_spectrum_obj.sort_by_abundance():
                                    if mspeak:
                                         molecular_formula = mspeak.molecular_formula_lowest_error
                                         pyplot.plot(mspeak.mz exp, mspeak.abundance, 'o', c='g')
           Data analysis
                                         for molecular_formula in mspeak:
                                              if molecular formula.is isotopologue:
       and visualization
                                                 print (molecular_formula.to string())
                                                 print (molecular_formula['13C'])
                                    else:
                                          print(mspeak.mz exp,mspeak.abundance)
                               mass_spectrum_obj
                   Output
```

More examples available at:

https://github.com/EMSL-Computing/CoreMS/tree/master/doc/examples



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## **Software Framework**

**CoreMS** Hierarchical Data Structure CoreMS GC-MS Workflow LC-MS(n) Retention Mass IM-MS GC-MS index (TODO) Spectra (Active Dev) calibration Chrom. Signal Peak Obj **Processing** m/z Domain Mass Reference **Transient** database Spectrum data search SQL3 **Post** Mass Spec gres Peak obj Molecular **Spectral** Compound Obj **Object** Matching

For available input, output and data structure types please refer to: <a href="https://github.com/EMSL-Computing/CoreMS/blob/master/README.md">https://github.com/EMSL-Computing/CoreMS/blob/master/README.md</a>



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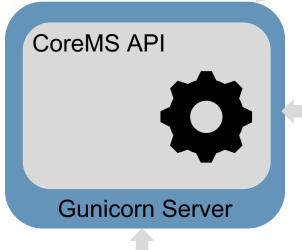
# **Web Application**

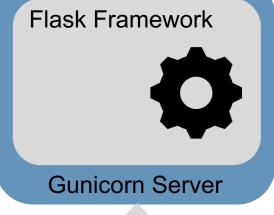
• Graphical User Interface (GUI):

- Back-End:
  - User based temporary storage, and data model for raw, mass-list and reference file data
  - User based data model and storage for results and metadata
  - User based data model and storage of workflow parameters

Redis: Data processing queue management

PostgreSQL: User, Data, Parameters Results







PostgreSQL: Dynamic Molecular Formula Generation





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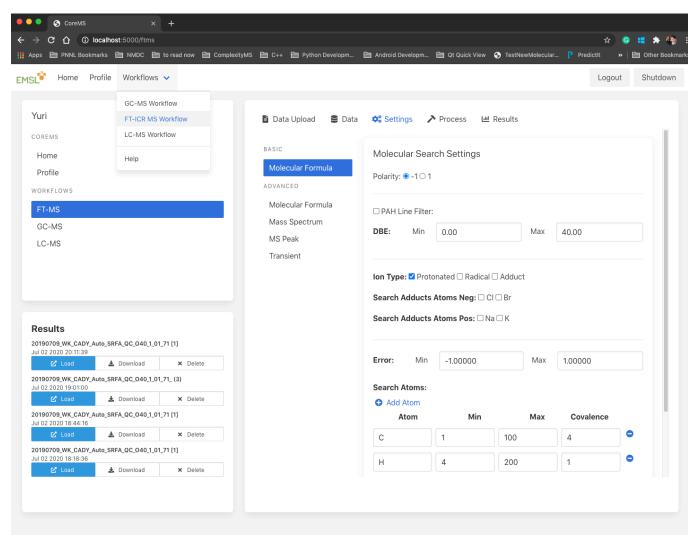
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# **Web Application**

# • Graphical User Interface (GUI):

- Front-End:
  - Web Application framework (HTML, CSS, JavaScript and, jQuery)
  - Data management for temporary raw data storage
  - Data management for workflow parameters, and parameters presets
  - Raw data visualization (Thermo, and Bruker)
  - Workflow result visualization



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# Thank you

