

General Fragments





Fragment-based approach has emerged as a powerful tool in modern drug discovery. Most of the hits from conventional screening would hardly fit the Lipinski rules, not to say the Ro4.

Thus, stricter rules have been proposed – first of all, to increase the potential quality of the obtained hit. At the same time, small library could cover relatively high area of chemical space since the final drug could be a combination of several fragments.

Fragments are widely used in various, sometimes very specific, research projects. Discover Chemspace fragments subsets that can be applicable in NMR-assisted and RSA-assisted screening.



Chemspace **General Fragments** library consists of compounds that

- Comply with Ro3
- No PAINS, "toxic" and reactive moieties*

*Please see also our **Covalent Fragments** set that go beyond this library.

Library size:

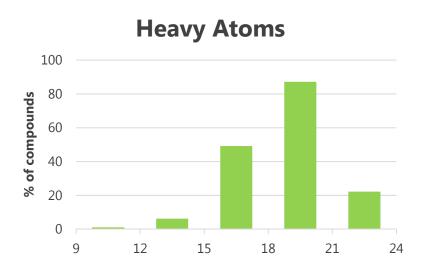
86 837 in-stock compounds

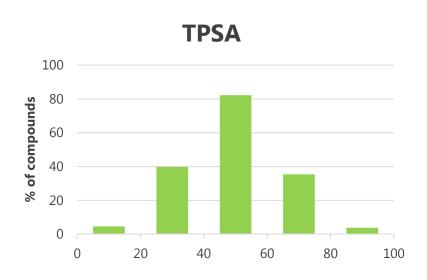
1 244 258 make-on-demand compounds

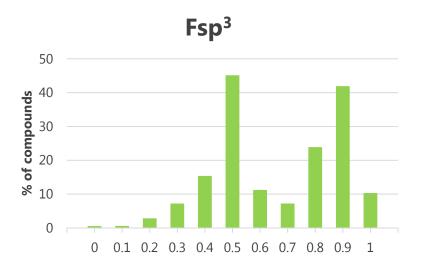
You can order full set or selected subset based on your criteria; all compounds are supplied as powders, solutions, or dry films. Please contact us at sales@chem-space.com for more information.

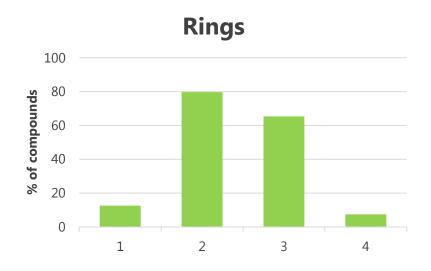


PhysChem properties











Chemspace Compound sets

Discover our **Fragment Libraries**:

- **General** Fragments
- **3D-Shaped** Fragments
- Acid and Amine Fragments
- **Covalent** Fragments

- Fluorine and Heavy Fragments
- **Selected** Fragments
- <u>Singleton</u> Fragments
- **Saturated** and **Spiro** Fragments

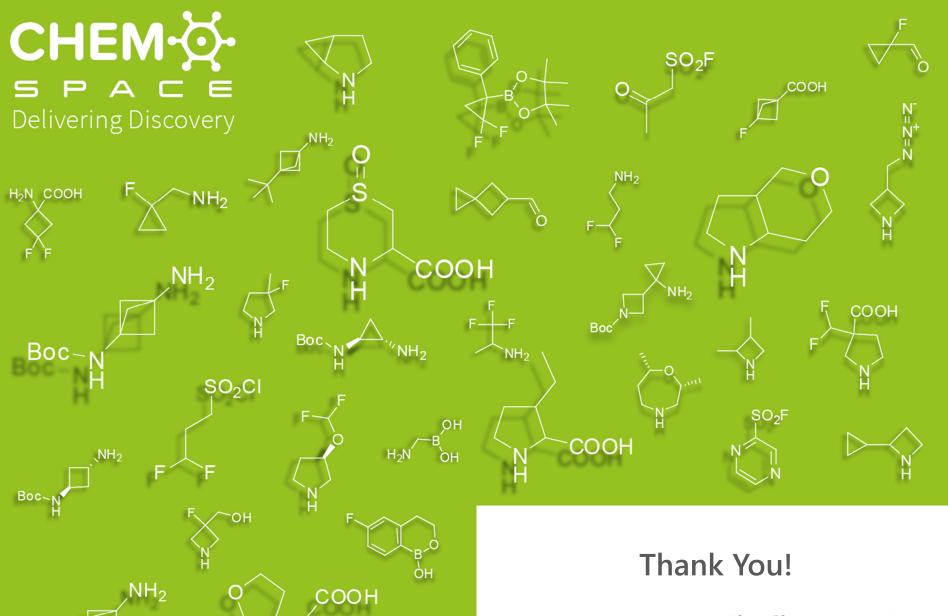


Chemspace Compound sets

Discover our **Screening compounds**:

- ChEMBL analogs
- CNS-Focused library
- Covalent Modifiers
- Drug Impurities
- Drug Repurposing
- Framework-Derived set

- <u>High QED</u> compounds
- Phenotypic Screening set
- PPI Modulators
- <u>Pre-Plated</u> compounds
- RNA-Targeted library
- **Virtual Screening** set



The Chemspace Team

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