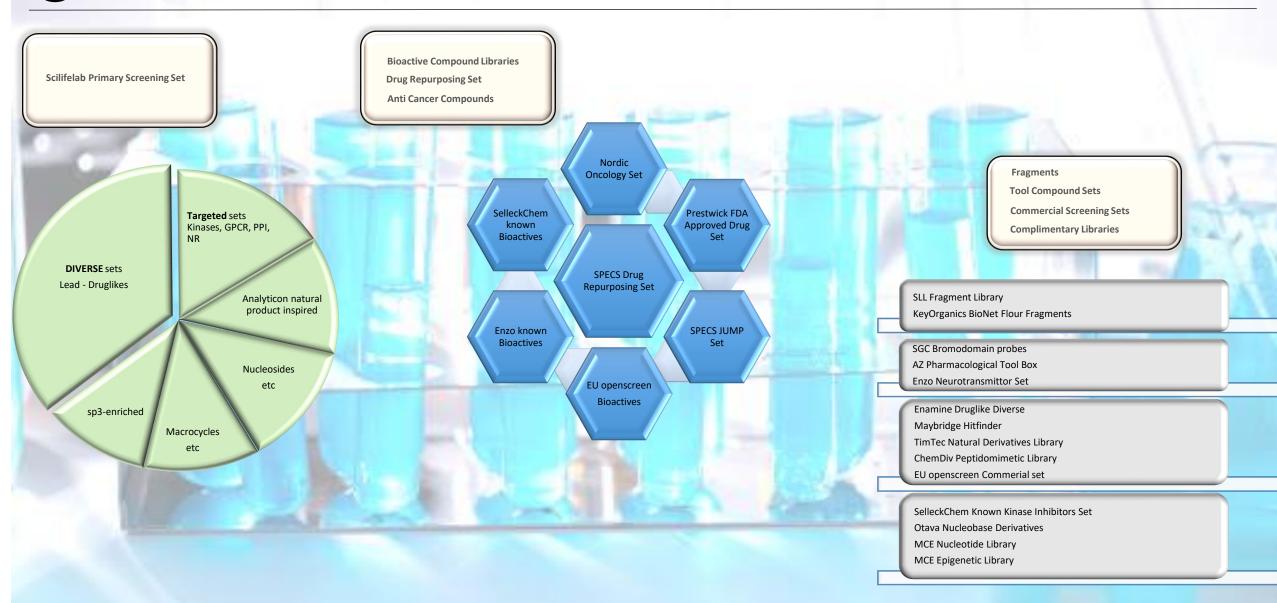


CBCS Primary screening set other available sets









SPECS Drug Repurposing Set

 6000 Drug repurposing set (Phase 1-3, preclinical and launched)built by The Broad Institute of MIT and Harvard, Cambridge, Massachusetts, USA

Prestwick FDA Approved Drug Set

 1200 FDA & EMA approved drugs (current or withdrawn). A diversified marketed drugs library designed for repurposing/repositioning with known bioavailability and safety in humans.

Nordic Oncology Set

• 500 Anti Cancer Compounds shared with the Nordic academia











SelleckChem known Bioactives

• A unique collection of 990 bioactive compounds for screening and high content screening

SPECS JUMP Set(Joint undertaking in morphological profiling)

• 395 compounds Target and MoA, designed by Broad Institute. Suits cellular imaging, image analysis and high dimensional data analytics. Read more: https://jump-cellpainting.broadinstitute.org/









SGC Bromodomain probes

• A library of 40 epigenetic chemical probes.

AZ Pharmacological Toolbox

• 23 compounds with optimized pharmacological properties made available for preclinical research to explore novel disease biology and advance scientific knowledge.

BioMol (Enzo)Neurotrasmittor Set

• 700 CNS receptor ligands. Ideal for screening or identifying recombinant orphan G protein-coupled receptors, target validation, secondary screening, validating new assays, and for routine pharmacological applications. Includes 13 classes of receptor ligands.







SLL Enamine Druglike Diverse

• 28 500 compounds available

SLL Maybridge Hitfinder

- 14 400 compounds available
- Collection designed using a clustering algorithm employing Daylight Fingerprints with the Tanimoto similarity index clustering at 0.71 % similarity.

SLL TimTec Natural Derivatives Library

- 3000 compounds available
- Contains pure natural compounds, synthetic compounds and syntethically modified pure natural compounds: alkaloids, natural phenols, nucleoside analogs, carbohydrates, purines, pyrimidines, flavonoids, steroidal cpds and natural amino acids.

SLL ChemDiv Peptidomimetic Library

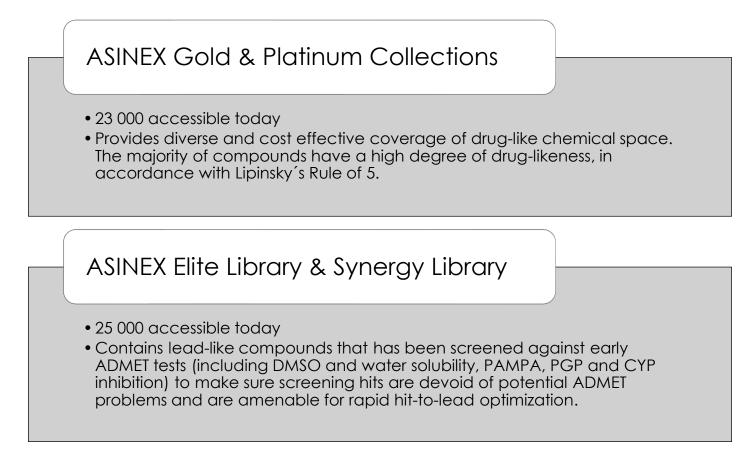
- 15 600 compounds available
- Contain organic molecules that mimic the action of peptides. The molecules may structually resemble petides but are distinctly different in terms of their side chains or their molecular backbones.











PAMPA = Parallel artificial membrane permeability assay PGP = P-glycoprotein inhibition measurement CYP = Cytochrome P450







Fragment screening set 10 - 200mM (5K)

- 5 000 fragments of concentrations between 10 200 mM.
- Designed by Flavio Ballante(CBCS)

KeyOrganics BioNet Flour Fragments

- 462 fragments 200mM available.
- Suitable for NMR screening.

SLL CIMS Fragments

- 1 000 fragments 50mM availble.
- Designed by Anna-Lena Gustavsson(CBCS), Evert Homan(T. Helleday group), Peter Brandt(Kdex)





EU Openscreen Pilot set	
 5 000 cpds Representative and Bioactives 	
EU Openscreen Commercial set	
• 95 000 compounds	
EU Openscreen Fragments	
1000 fragments 200mM available.Suitable for NMR screening.	







SelleckChem Known Kinase

Inhibitors Set

• 378 known kinase inhibitors available

Otava Nucleobase Derivatives

• 958 NIH compounds available

MCE Nucleotide Library

• 177 nucleotides available

MCE Epigenetic Library

• 769 compound available. Includes any process that alters gene activity without changing the DNA sequence and leads to modifications that can be transmitted to daughter cells.

LCBU Chembridge Diverse Part 2

• 7700 diverse compounds available







