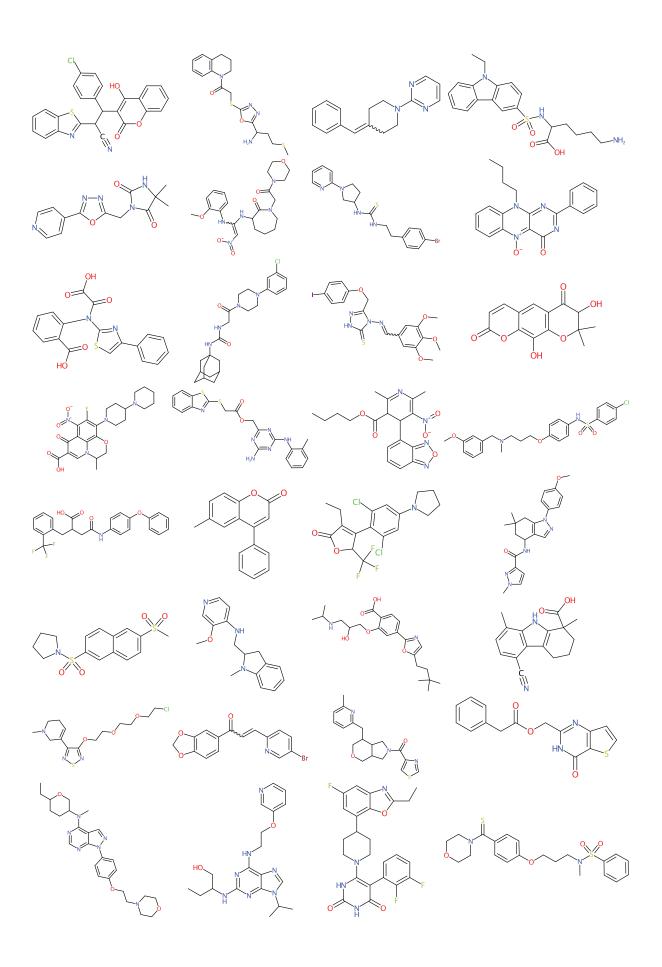# Supporting Information for
Generating Focussed Molecule Libraries for Drug Discovery with Recurrent Neural Networks

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In the following, a few randomly selected molecules produced with the general model trained on ChEMBL are shown.