

O=C(O)[C@H]1CCOC[C@H]1c1ccc(I)cc1

The diagram illustrates the decomposition of a chemical SMILES string. A top box contains the full string: O=C(O)[C@H]1CCOC[C@H]1c1ccc(I)cc1. Two arrows point downwards from this box to two separate boxes below. The left box contains the fragment O=C(O)C1CCOCC1, and the right box contains the fragment Ic1ccc(I)cc1.

O=C(O)C1CCOCC1

Ic1ccc(I)cc1