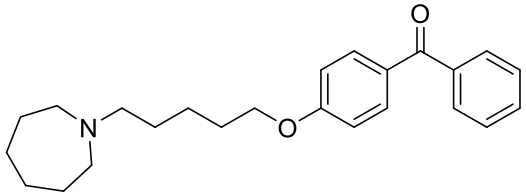
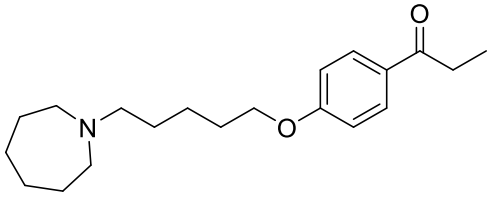
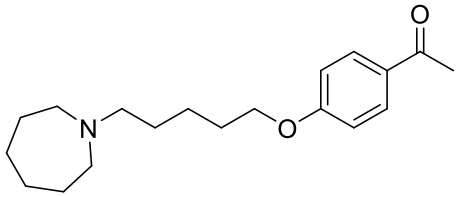
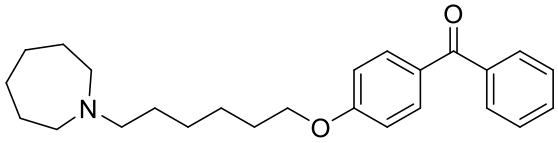
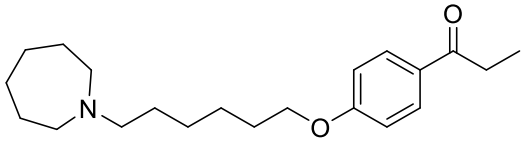
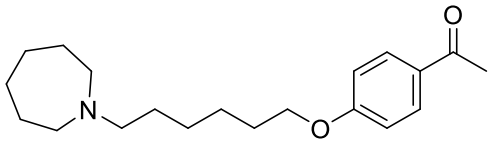
Preliminary calculations done using SwissAdme server (http://www.swissadme.ch/ (Last accessed on 28 December 2022)

name	Compound - structure	Molecular weight	Consensus Log P
Lead 2		365	4.82
15		317	4.16
14		303	3.79
Lead 3		380	5.18
17		331	4.45
16		317	4.15