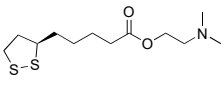
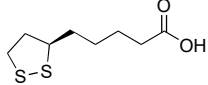
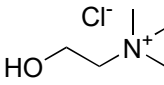
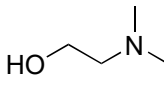
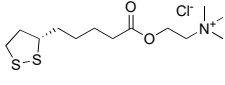
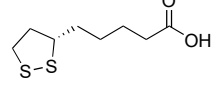
 <chem>CN(C)[CH+]([O-])CCCCC[C@H](CS1SS1)C(=O)O</chem>	 <chem>CN(C)[CH+]([O-])CCCCC[C@H](CS1SS1)C(=O)O</chem>	 <chem>OC(=O)CCCC[C@H](CS1SS1)C(=O)O</chem>	 <chem>CN(C)[CH+]([O-])CO</chem>	 <chem>CN(C)CO</chem>
 <chem>CN(C)[CH+]([O-])CCCCC[C@@H](CS1SS1)C(=O)O</chem>	 <chem>OC(=O)CCCC[C@@H](CS1SS1)C(=O)O</chem>			