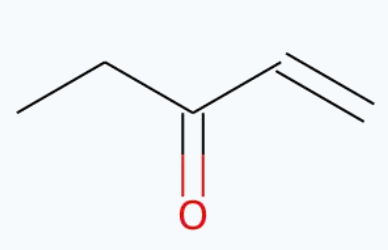
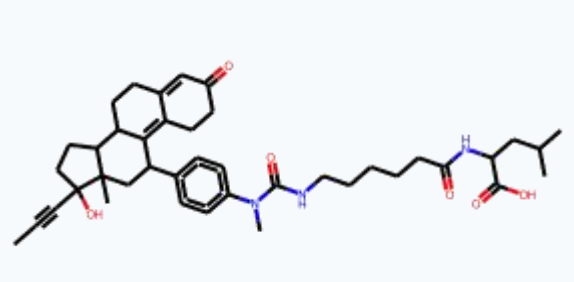
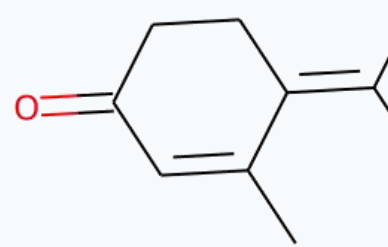
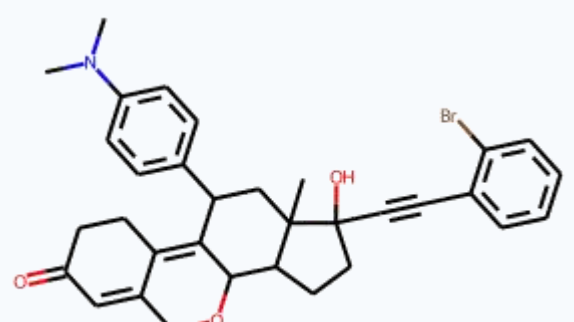
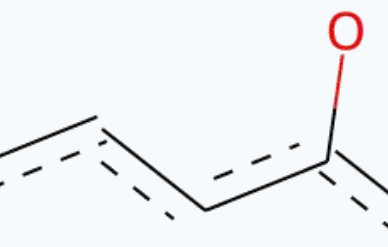
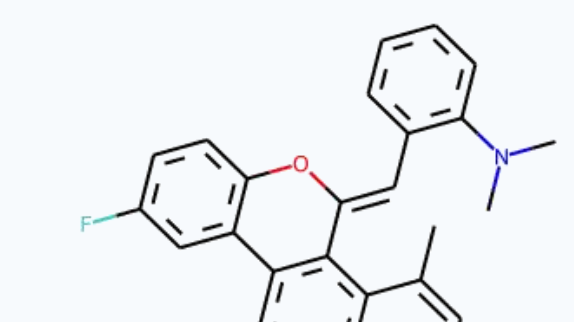
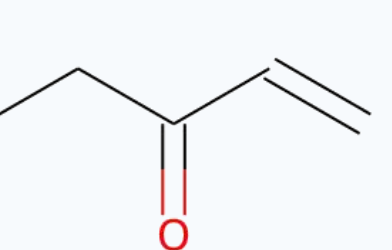
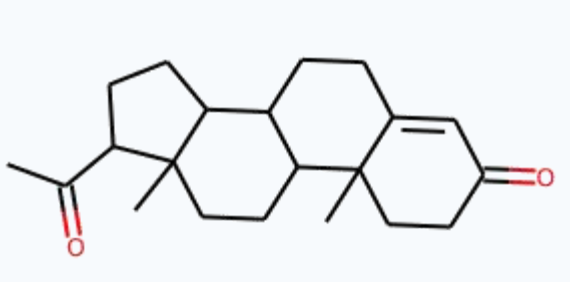
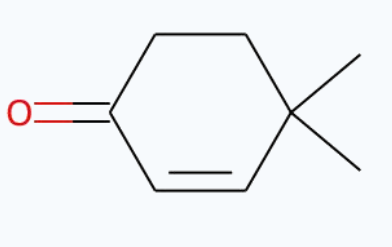
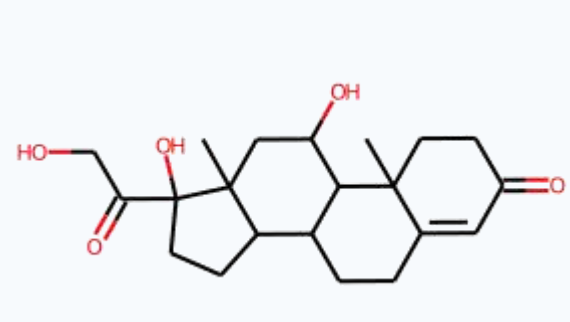
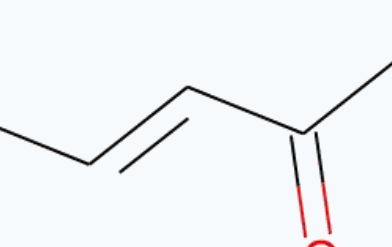
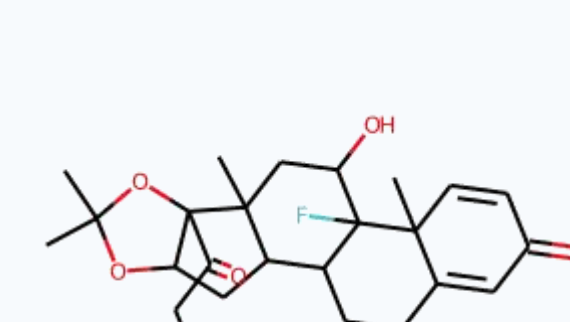


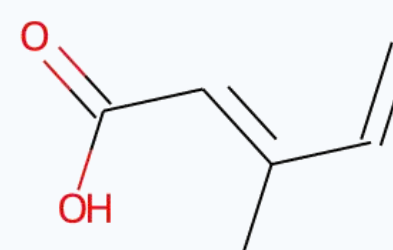
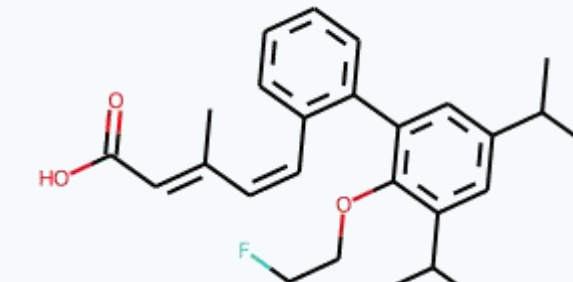
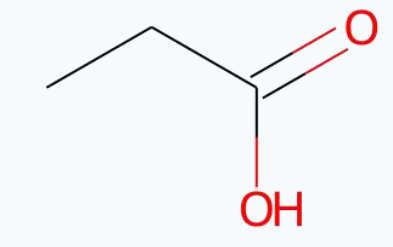
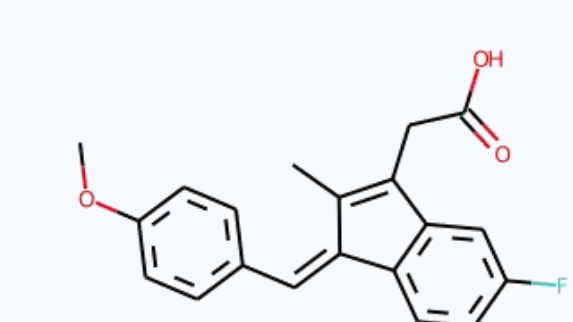
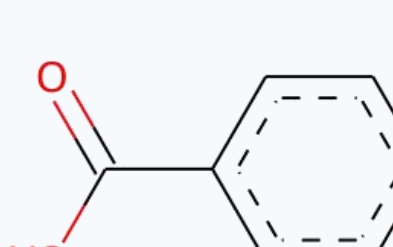
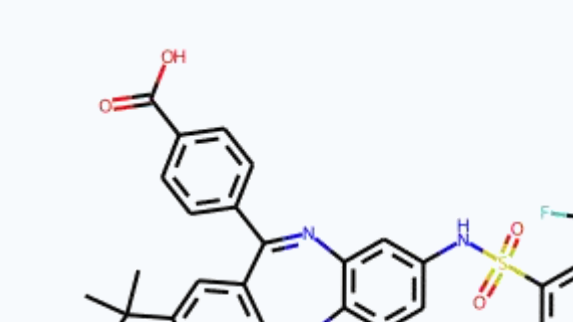
PR

SA	Representative Structure	SMILES	p-value
		<chem>CC#CC1(O)CCC2C1(C)CC(c1ccc(cc1)N(C(=O)NCCCCC(C(=O)NC(C(=O)O)CC(C)C)C1=C3CCC(=O)C=C3CC21</chem>	2,0367E-67
		<chem>O=C1CCC2=C3C(CC4(C(C3OCC2=C1)CCC4(O)C#Cc1cccc1Br)C)c1ccc(cc1)N(C)C</chem>	3,0253E-135
		<chem>Fc1ccc2c(c1)c1ccc3c(c1C(=Cc1cccc1N(C)O2)C(=CC(N3)C)O)C</chem>	3,1310E-41

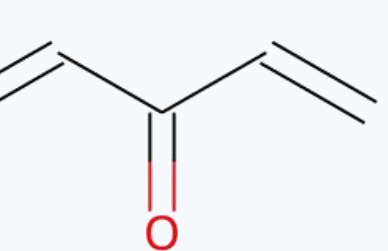
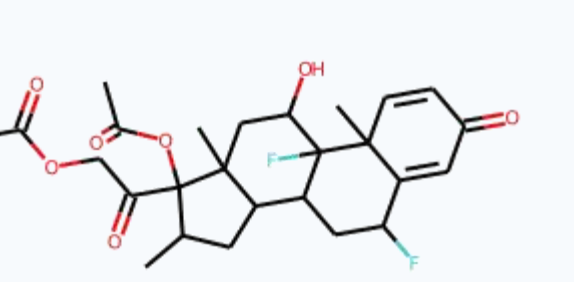
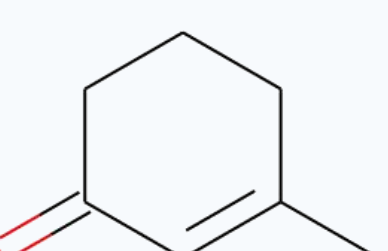
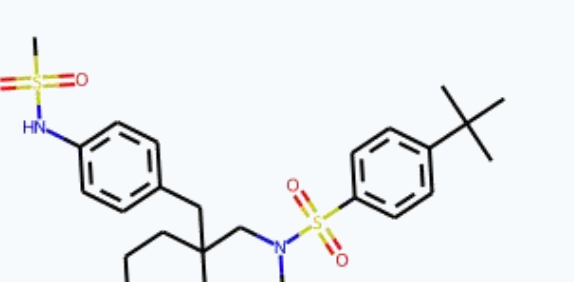

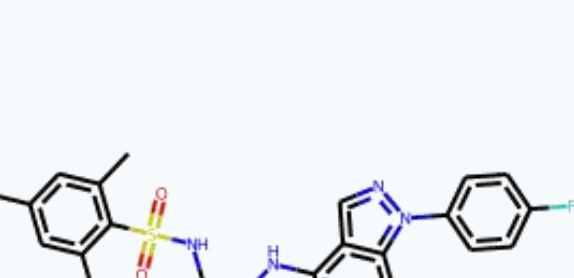
PXR

SA	Representative Structure	SMILES	p-value
		<chem>O=C1CCC2(C(=C1)CCC1C2CCC2(C1CCC2C(=O)C)C)C</chem>	7,5419E-20
		<chem>OCC(=O)C1(O)CCC2C1(C)CC(O)C1C2CCC2=C(C1(=O)CCC12C</chem>	5,9073E-12
		<chem>OCC(=O)C12OC(OC1CC1C2(C)CC(O)C2(C1CCC1=CC(=O)C=CC2(C)F)C)C</chem>	1,1977E-12

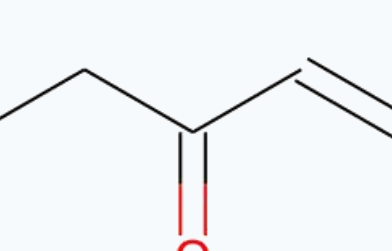
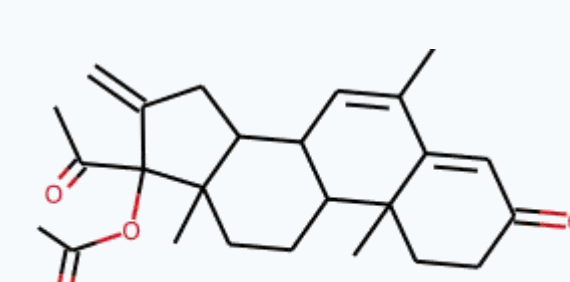
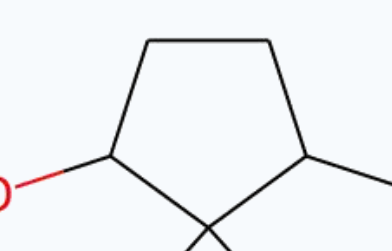
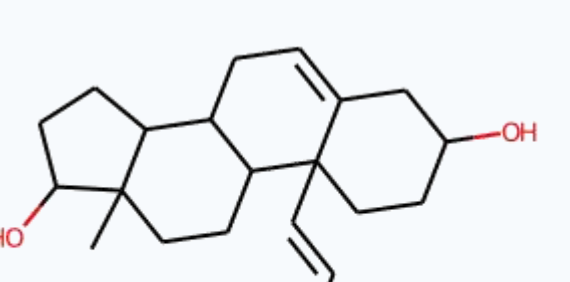
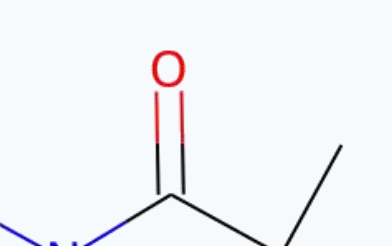
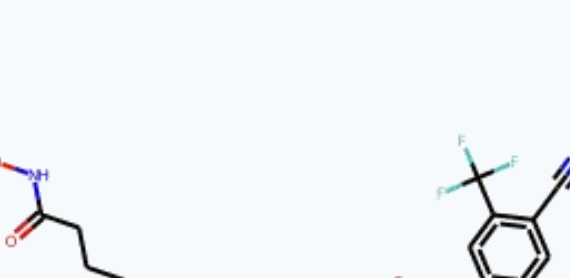
RXR

SA	Representative Structure	SMILES	p-value
		<chem>FC(COc1c(cc(cc1C(C)C)C(C)C)C1=C(CCC1)C=CC(=CC(=O)O)C(F</chem>	2,5303E-30
		<chem>CCOc1ccc(cc1)C=C1c2ccc(cc2C(=C1C)CC(=O)O)F</chem>	1,0458E-08
		<chem>CCCN1c2ccc(cc2N=C(c2c1cc1c(c2)C(C)C)CCC1(C)C)c1ccc(cc1)C(=O)O)NS(=O)(=O)c1cccc(c1)C(F)F)F</chem>	1,2584E-25

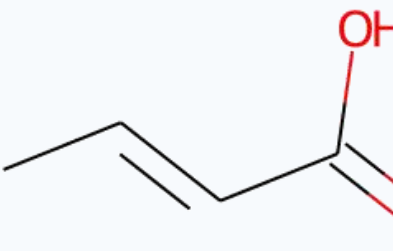
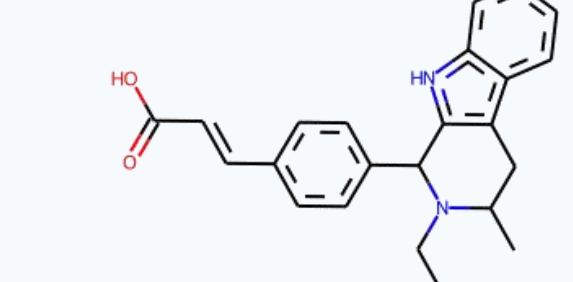
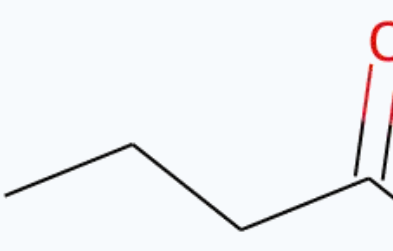
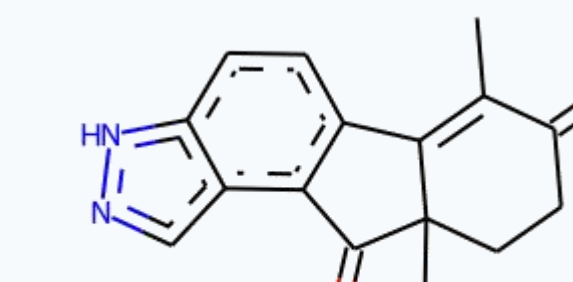
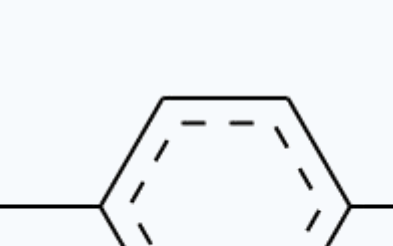
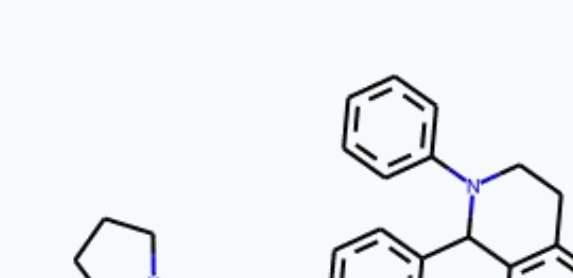
GR

SA	Representative Structure	SMILES	p-value
		<chem>CC(=O)OC1(C(C)CC2C1(C)CC(O)C1(C2CC(C2=CC(=O)C=CC12C)F)C(=O)COC(=O)C</chem>	8,7039E-30
		<chem>CC(C)(C)c1ccc(S(=O)(=O)N2CCCC3=CC(=O)CCC3(Cc3ccc(NS(C)C(=O)=O)cc3)C2)cc1</chem>	1,2213E-18
		<chem>CC(NS(=O)(=O)c1c(C)cc(cc1C)C)CNc1cccc2c1cnn2c1ccc(nc1)F</chem>	8,0567E-27

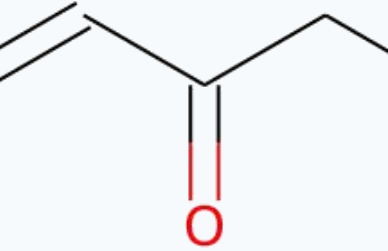
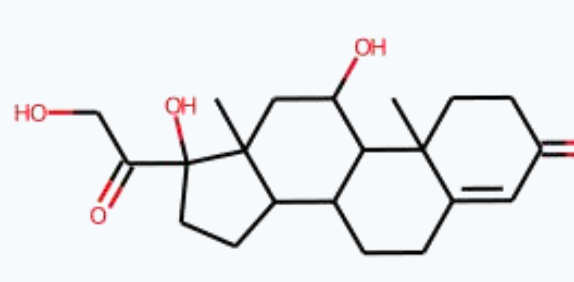
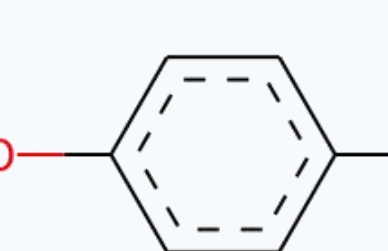
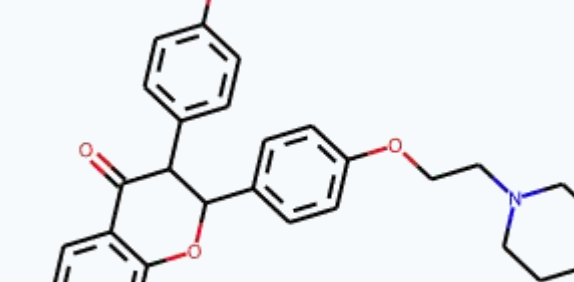
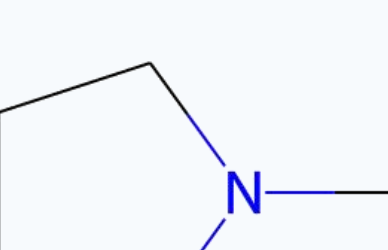
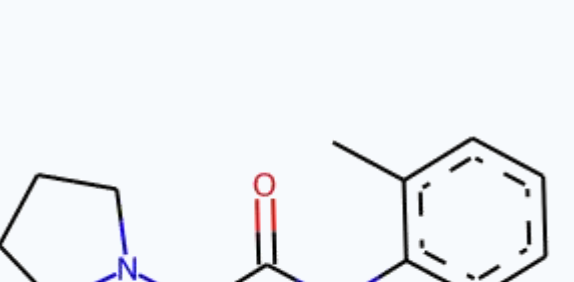
AR

SA	Representative Structure	SMILES	p-value
		<chem>O=C1CCC2(C(=C1)CCC1C2CCC2(C1C2C2(=O)COC(=O)C(C)C)C)C</chem>	1,6453E-39
		<chem>COC=CC12CCC(C1=CC1C2CCC2(C1CCC2O)C)O</chem>	8,8649E-17
		<chem>CC1(C)N(CCCCc2cn(CCCC(=O)NO)n2)C(=O)N(C1=O)c1ccc(C#N)c(c1)C(F)F)F</chem>	2,0294E-17

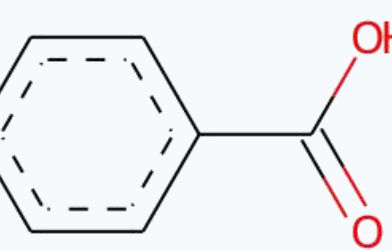
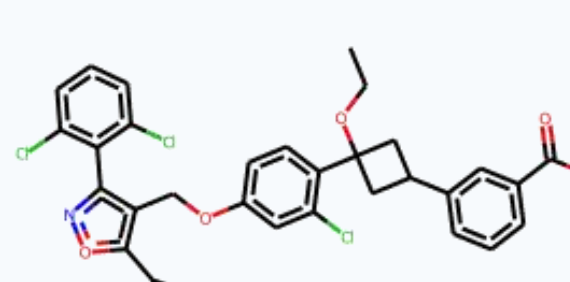
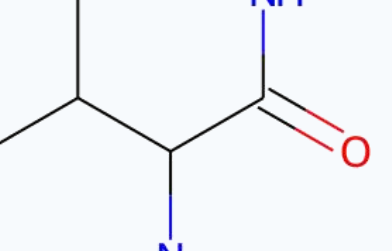
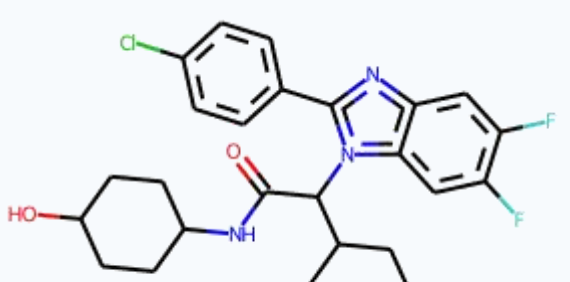
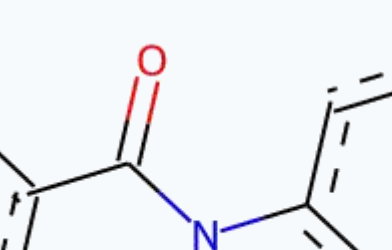
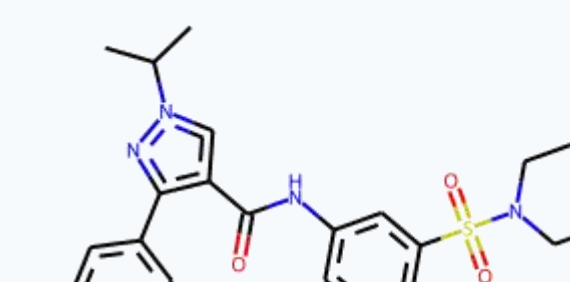
ERA

SA	Representative Structure	SMILES	p-value
		<chem>OCC(CN1C(C)Cc2c(C1c1ccc(cc1)C=CC(=O)O)[nH]c1c2cccc1)C</chem>	5,6428E-21
		<chem>CCC12CCC(=O)C(C(=C1c1c(C2=O)c2cn[nH]c2cc1)C</chem>	1,6971E-33
		<chem>Oc1ccc2c(c1)CCC(c1cccc1)N2Cc1ccc(OCCN2CCC(CC2)cc1</chem>	3,0141E-36

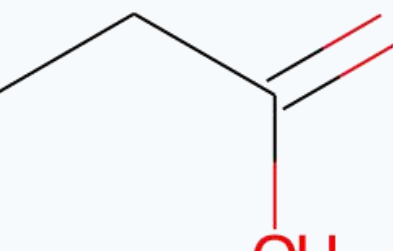
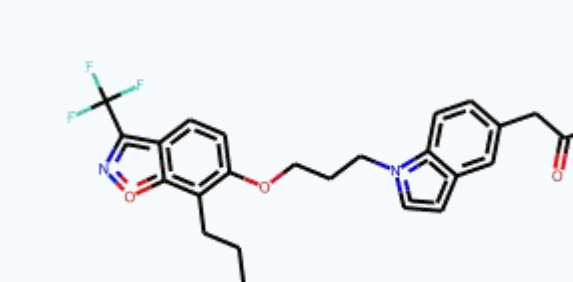
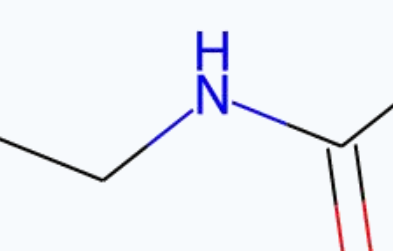
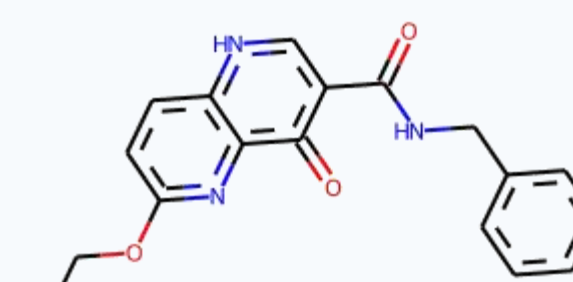
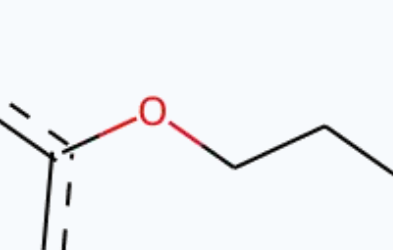
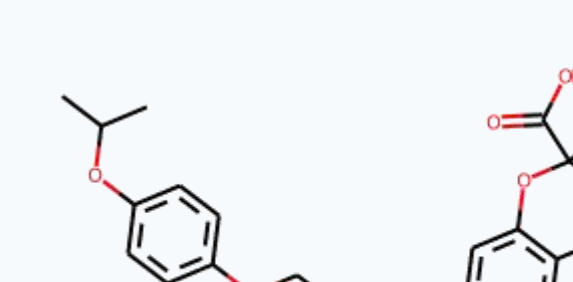
ERB

SA	Representative Structure	SMILES	p-value
		<chem>OCC(=O)C1(O)CCC2C1(C)CC(O)C1C2CC2=CC(=O)C)CCC12C</chem>	2,5157E-18
		<chem>Oc1cc2OC(c3ccc(cc3)OCCN3CCCC3)C(C(=O)c2c(c1)C)c1ccc(cc1)O</chem>	2,9580E-108
		<chem>O=C(Nc1c(C)cccc1C)N2CCC2</chem>	1,9932E-76

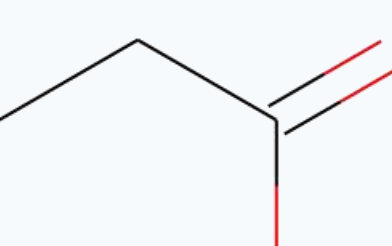
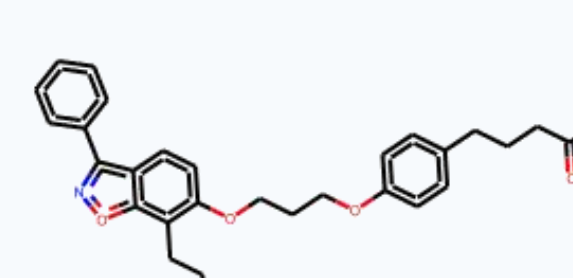
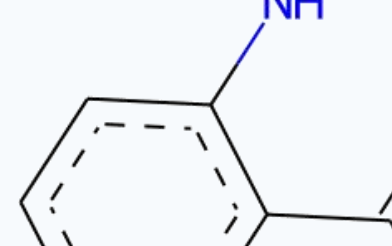
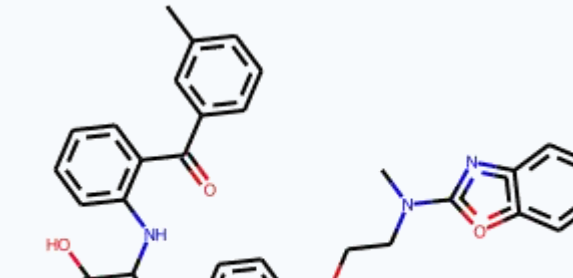
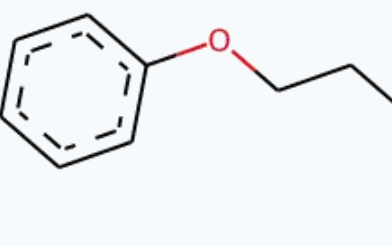
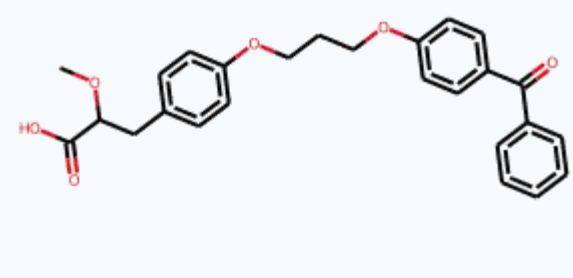
FXR

SA	Representative Structure	SMILES	p-value
		<chem>Cl1cc(ccc1C1(O)CC(C1)c1ccc(cc1)C(=O)O)OCc1c(onc1c(Cl)cccc1C)C1CC1</chem>	1,1107E-41
		<chem>OC1CCC(CC1)NC(=O)C(n1c(nc2c1cc(F)c2)F)c1ccc(cc1C1)C1CCCC1</chem>	2,8113E-31
		<chem>COc1cccc(c1)c1nn(cc1C(=O)Nc1ccc(c1)S(=O)(=O)N1CCOC(C1)C)C)C</chem>	1,3146E-30

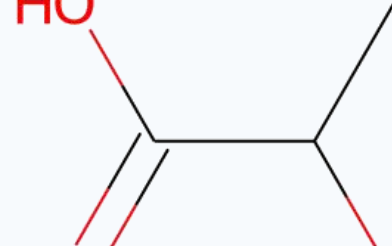
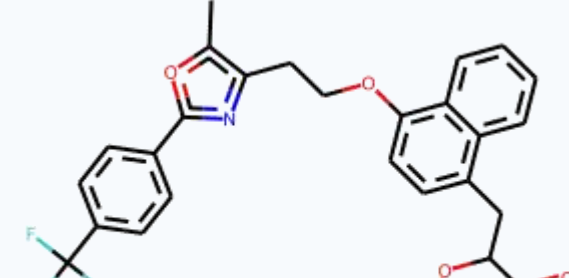
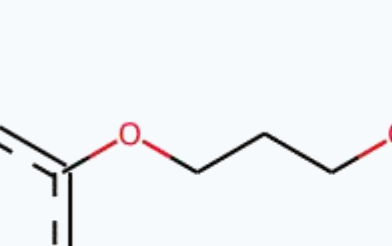
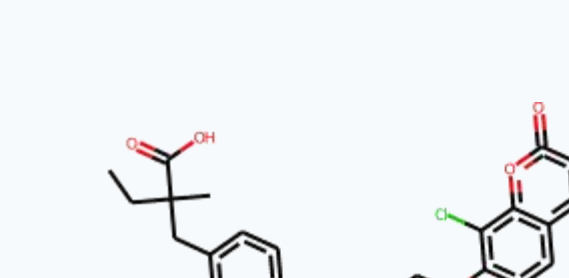
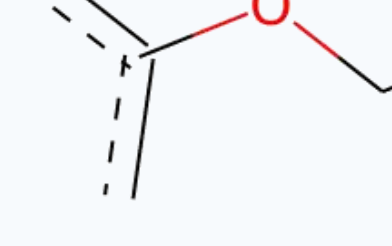
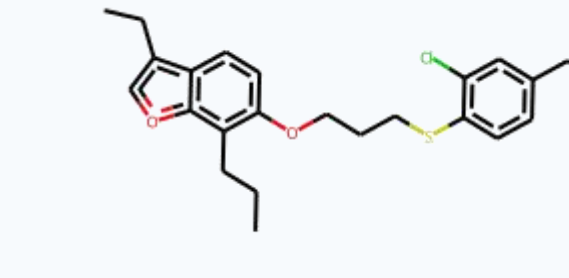
PPARD

SA	Representative Structure	SMILES	p-value
		<chem>CCCc1c(OCCCn2cnc3cc(CC(=O)O)ccc32)ccc2c(C(F)F)F)noc12</chem>	4,4501E-40
		<chem>CCOc1ccc2c(n1)c(=O)c(c[nH]2)C(=O)NCC1cccc1</chem>	3,1489E-30
		<chem>CC(Oc1cccc(c1)C)OCCCOc1ccc2c(c1)OC(CC2)C(C(=O)O)C</chem>	8,6807E-50

PPARG

SA	Representative Structure	SMILES	p-value
		<chem>CCCc1c(OCCCOc2cccc2CCC(=O)O)ccc2c1onc2c1cccc1</chem>	7,7458E-23
		<chem>Cc1ccc(c(c1)C(=O)c1cccc1)NC(C(=O)O)Cc1ccc(cc1)OCCN(c1nc2c(o1)cccc2)C</chem>	2,0328E-34
		<chem>COC(C(=O)O)Cc1ccc(cc1)OCCCOC1cccc(c1)C(=O)c1cccc1</chem>	2,3729E-30

PPARA

SA	Representative Structure	SMILES	p-value
		<chem>CCOC(C(=O)O)Cc1ccc(c2c1cccc2)OCCc1nc1oc1c1ccc(cc1)C(F)F)F</chem>	3,0577E-09
		<chem>CCC(C(=O)O)(Cc1ccc(cc1)OCCCOc1ccc2c(c1)oc(=O)cc2C(F)F)F)C</chem>	7,3175E-04
		<chem>CCCc1c(OCCCS2cccc(cc2C1)CC(=O)O)ccc2c1occ2CC</chem>	9,6182E-06