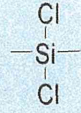
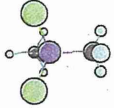
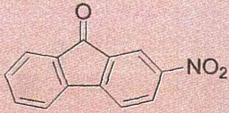
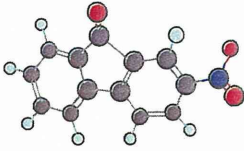
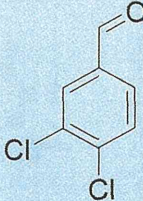
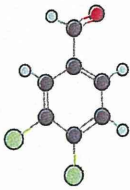

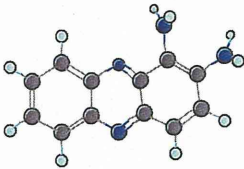
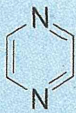
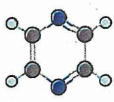
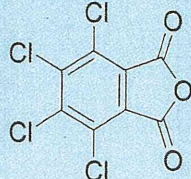
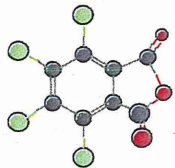
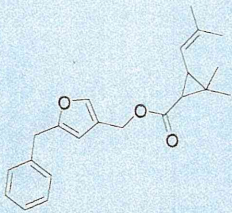
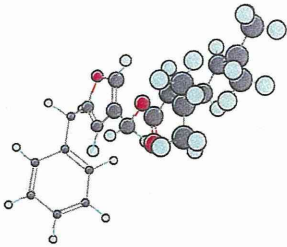

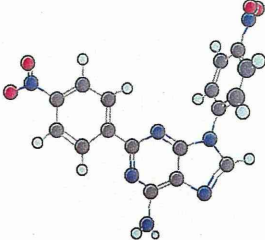
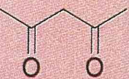
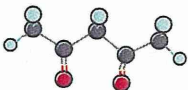

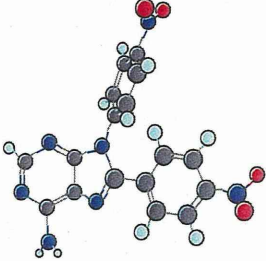
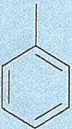
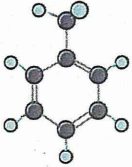
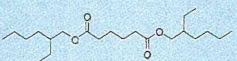
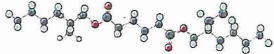
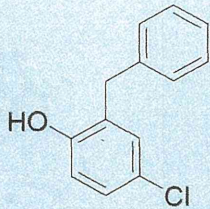
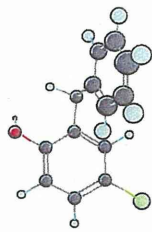
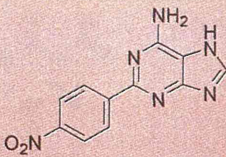
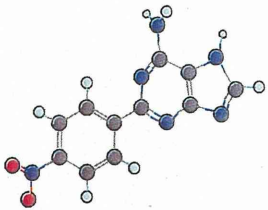
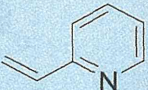
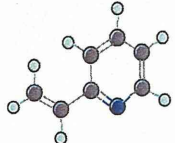
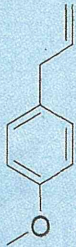
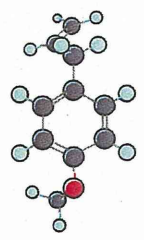
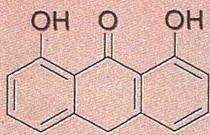
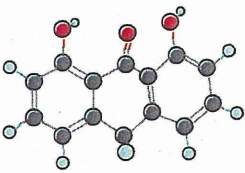
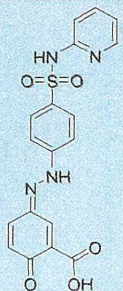


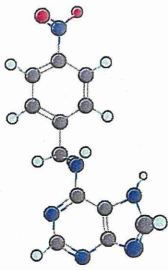
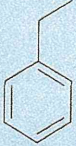
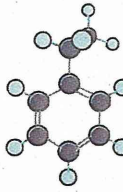

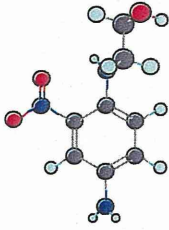

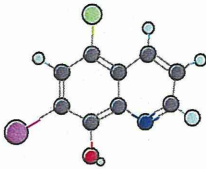
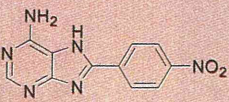
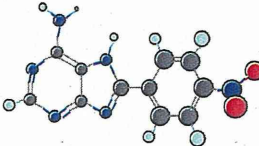
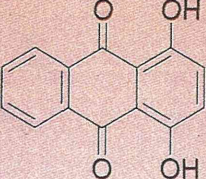
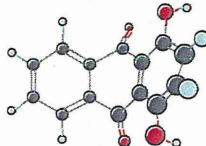
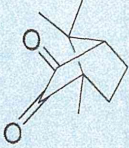
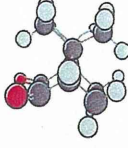
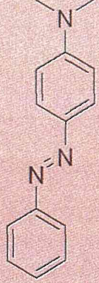
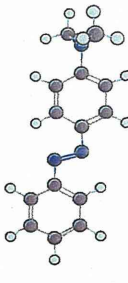
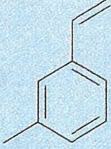
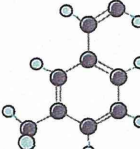
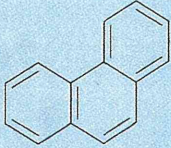
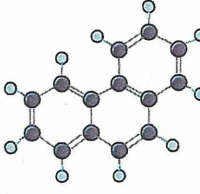
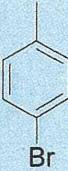
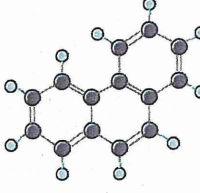
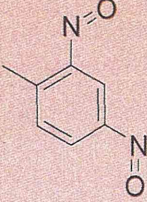
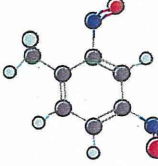


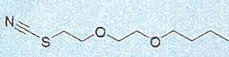

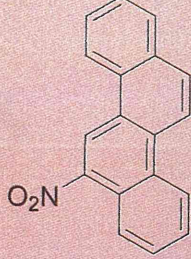
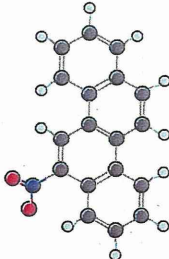
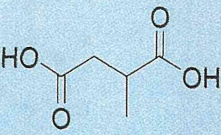

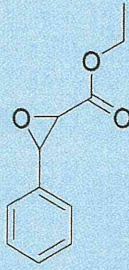
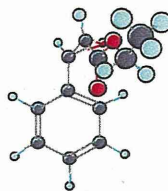
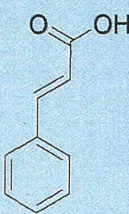
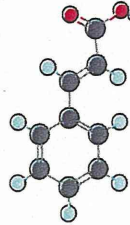
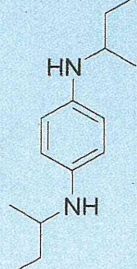
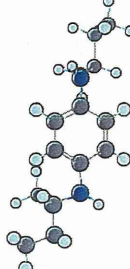
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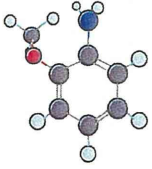
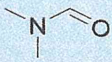
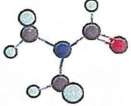

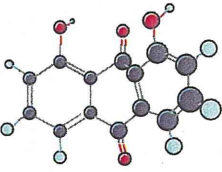

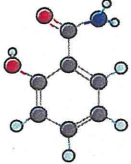
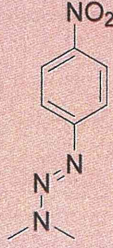
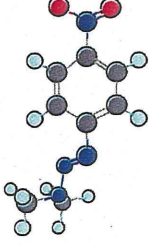
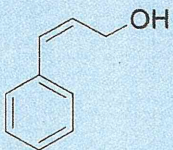
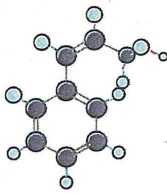
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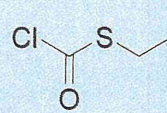
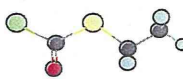
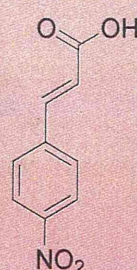
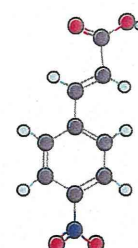
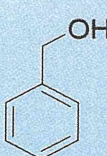
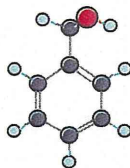
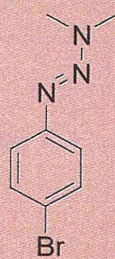
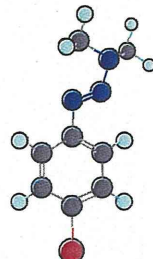
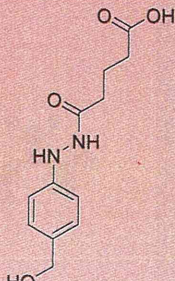
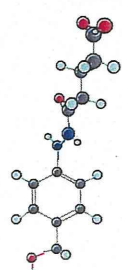
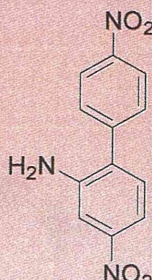
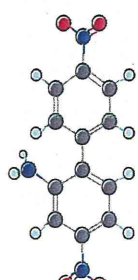
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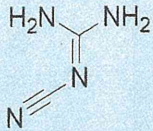
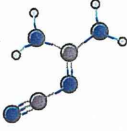
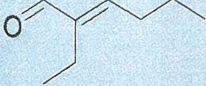
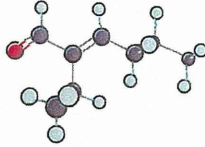
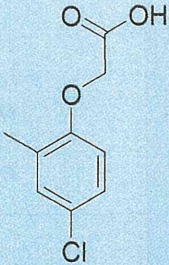
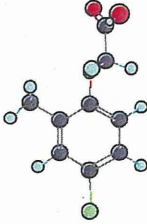
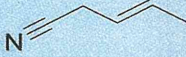
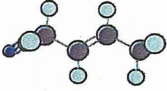
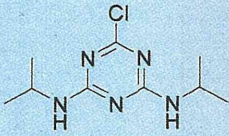
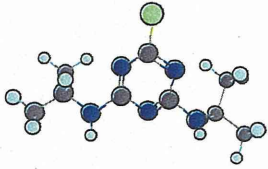
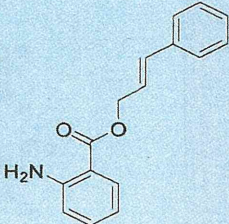
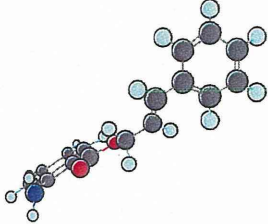
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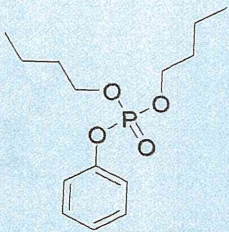
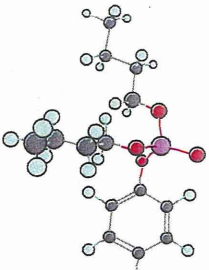
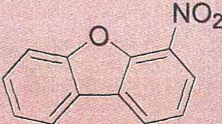
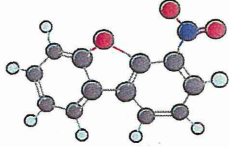
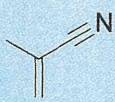
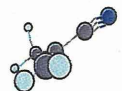
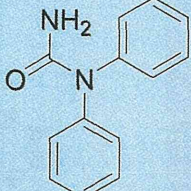
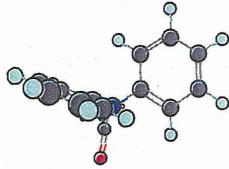

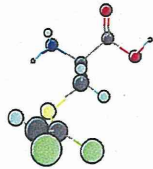
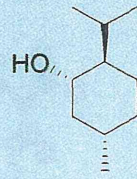
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114	 <chem>NC(=N)N#N</chem>	
115	 <chem>CCCC=C(C)C=O</chem>	
116	 <chem>CC1=CC=C(C=C1)OCC(=O)OCl</chem>	
117	 <chem>CCCC#N</chem>	
118	 <chem>CC(C)N1C=NC(Cl)=NC1N(C)C</chem>	
119	 <chem>NC1=CC=C(C=C1)C(=O)OCC=Cc2ccccc2</chem>	

120	 <chem>CCOP(=O)(OCC)Cc1ccccc1</chem>	
121	 <chem>O=C1C=CC=C2C(=O)N(C2)C1[N+](=O)[O-]</chem>	
122	 <chem>C=CC#N</chem>	
123	 <chem>NC(=O)N(Cc1ccccc1)c2ccccc2</chem>	
124	 <chem>NC(C=C(Cl)Cl)CO</chem>	
125	 <chem>CC1(C)CCCC(O)C1</chem>	