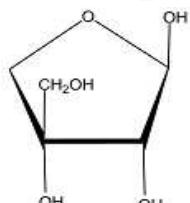
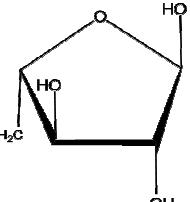
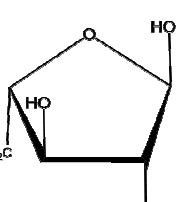
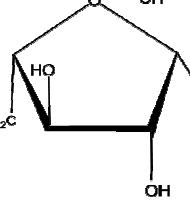
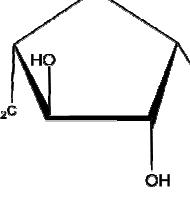
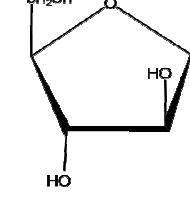
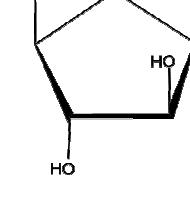
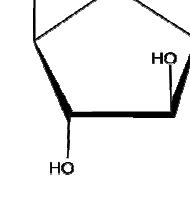
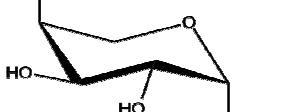
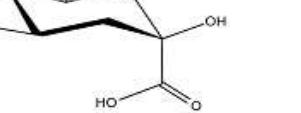
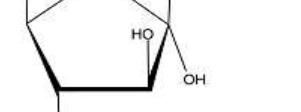
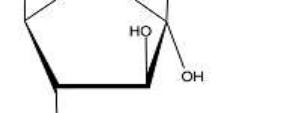
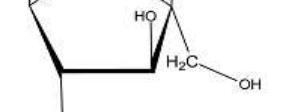
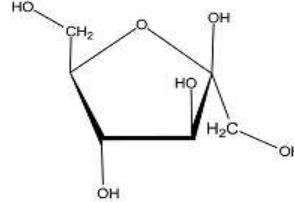
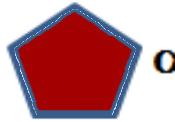
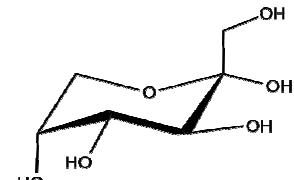
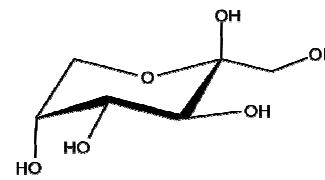
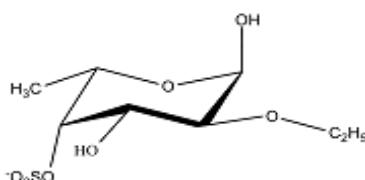
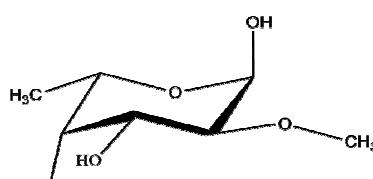
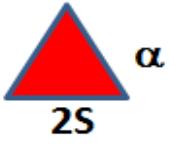
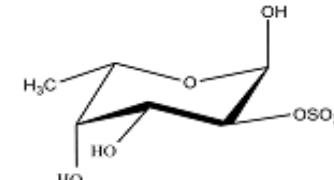
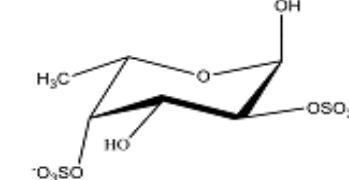
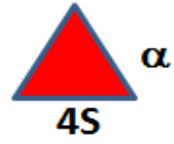
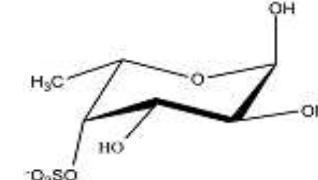
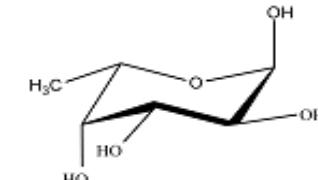


Nomenclature	Symbol	Structure 3D	File Name
αD_Abe			Abe_aD
$\beta L_Acef A$			AceA_bL(S)
$\beta L_Acef A$			AceA_bL(N)
$\alpha L_Acef A$			AceA_aL(N)
$\alpha L_Acef A$			AceA_aL(S)
$\alpha D_Alt p$			Alt_aD
$\beta D_Alt p$			Alt_bD
βD_Apif			Api_bD(N)

βD_Apif			Api_bD(S)
αL_Araf			Ara_aL(S)
αL_Araf			Ara_aL(N)
βL_Araf			Ara_bL(N)
βL_Araf			Ara_bL(S)
αD_Araf			Ara_aD(S)
αD_Araf			Ara_aD(N)
βD_Araf			Ara_bD(S)

β D_Araf					Ara_bD(N)
α L_Ara					Ara_aL
β L_Ara					Ara_bL
β D_Arap					Ara_bD
α D_Arap				Ara_aD	
DHA				DHA_bD	
α D_Fruf				Fru_aD(N)	
α D_Fruf				Fru_aD(S)	
β D_Fruf				Fru_bD(N)	

$\beta\text{D}_\text{-Fruf}$			Fru_bD(S)
$\alpha\text{D}_\text{-Frup}$			
$\beta\text{D}_\text{-Frup}$			
$\alpha\text{L}_\text{-Fucp [2Et4S]}$			Fuc[2Et4S]_aL
$\alpha\text{L}_\text{-Fucp [2Me]}$			Fuc[2Me]_aL
$\alpha\text{L}_\text{-Fucp [2S]}$			Fuc[2S]_aL
$\alpha\text{L}_\text{-Fucp [2S4S]}$			Fuc[2S4S]_aL
$\alpha\text{L}_\text{-Fucp [4S]}$			Fuc[4S]_aL
$\alpha\text{L}_\text{-Fucp}$			Fuc_aL

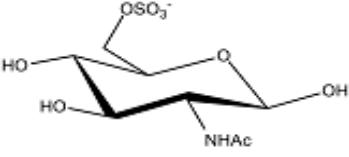
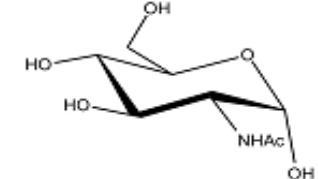
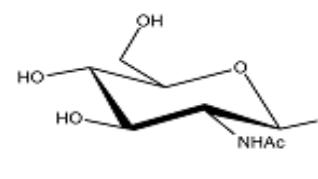
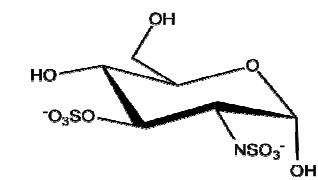
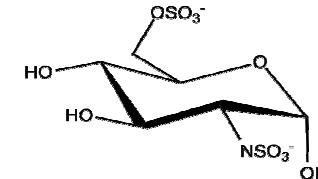
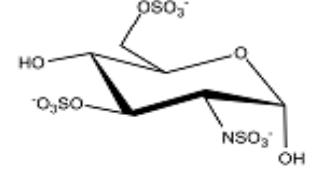
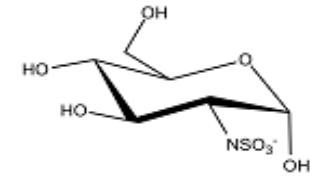
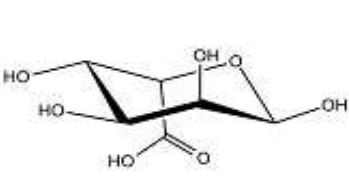
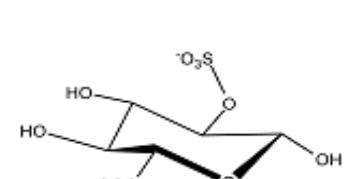
αD_Galp [2S]	α		Gal[2S]_aD
αD_Galp [2S3S]	α		Gal[2S3S]_aD
αD_Galp [2S6S]	α		Gal[2S6S]_aD
αD_Galp [36Anh]	α		Gal[36Anh]_aD
βD_Galp [3S]	β		Gal[3S]_bD
βD_Galp [3S4S]	β		Gal[3S4S]_bD
βD_Galp [3S6S]	β		Gal[3S6S]_bD
βD_Galp [4S]	β		Gal[4S]_bD
bD_Galp [4S6S]	β		Gal[4S6S]_bD
βD_Galp [6S]	β		Gal[6S]_bD

$\beta\text{D}_\text{Galp}$ [6Ac]			Gal[6Ac]_bD
$\alpha\text{D}_\text{Galp}$ [6OMe]			Gal[6OMe]_aD
$\alpha\text{D}_\text{Galp}$			Gal_aD
aL_Galp			Gal_aL
$\beta\text{D}_\text{Galp}$			Gal_bD
$\alpha\text{D}_\text{Galp A}$			GalA_aD
$\alpha\text{L}_\text{Galp A}$			GalA_aL
$\beta\text{D}_\text{Galp A}$			GalA_bD
$\alpha\text{D}_\text{Galp A}[2S3S6Me]$			GalA[2S3S6Me]_aD

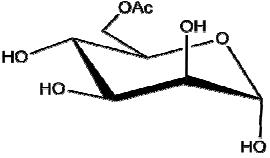
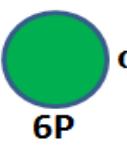
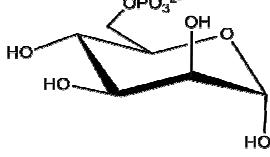
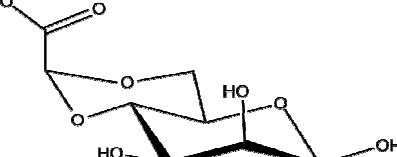
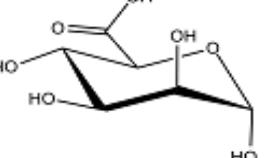
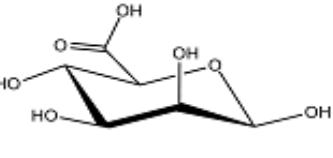
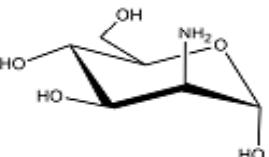
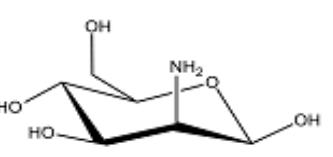
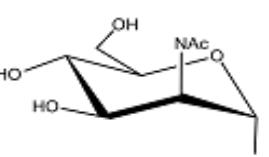
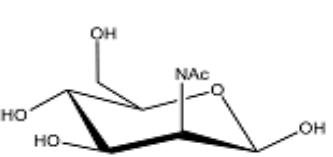
β D_Galp A[2S3S6Me]		β	 A chemical structure of a galactose residue in a chair conformation. It has a methyl group (CH ₃) at C1, a hydroxyl group (OH) at C2, two sulfate groups (-O ₃ SO ₃ ⁻) at C3 and C6, and a hydroxyl group (OH) at C4.	GalA[2S3S6Me]_bD
α D_Galp N		α	 A chemical structure of a galactose residue in a chair conformation. It has a hydroxyl group (OH) at C1, a hydroxyl group (OH) at C2, a hydroxyl group (OH) at C3, an amino group (NH ₂) at C4, and a hydroxyl group (OH) at C6.	GalN_aD
β D_Galp N		β	 A chemical structure of a galactose residue in a chair conformation. It has a hydroxyl group (OH) at C1, a hydroxyl group (OH) at C2, a hydroxyl group (OH) at C3, an amino group (NH ₂) at C4, and a hydroxyl group (OH) at C6.	GalN_bD
β D_Galp NAc[4S]		β	 A chemical structure of a galactose residue in a chair conformation. It has a hydroxyl group (OH) at C1, a hydroxyl group (OH) at C2, a sulfate group (-O ₃ SO ₃ ⁻) at C3, an acetyl group (NHAc) at C4, and a hydroxyl group (OH) at C6.	GalNAc[4S]_bD
β D_Galp NAc[6S]		β	 A chemical structure of a galactose residue in a chair conformation. It has a hydroxyl group (OH) at C1, a hydroxyl group (OH) at C2, a sulfate group (-OSO ₃ ⁻) at C3, an acetyl group (NHAc) at C4, and a hydroxyl group (OH) at C6.	GalNAc[6S]_bD
β D_Galp NAc[4S6S]		β	 A chemical structure of a galactose residue in a chair conformation. It has a hydroxyl group (OH) at C1, a hydroxyl group (OH) at C2, a sulfate group (-OSO ₃ ⁻) at C3, an acetyl group (NHAc) at C4, and a hydroxyl group (OH) at C6.	GalNAc[4S6S]_bD
α D_Galp NAc		α	 A chemical structure of a galactose residue in a chair conformation. It has a hydroxyl group (OH) at C1, a hydroxyl group (OH) at C2, a sulfate group (-OSO ₃ ⁻) at C3, an acetyl group (NHAc) at C4, and a hydroxyl group (OH) at C6.	GalNAc_aD
β D_Galp NAc		β	 A chemical structure of a galactose residue in a chair conformation. It has a hydroxyl group (OH) at C1, a hydroxyl group (OH) at C2, a sulfate group (-OSO ₃ ⁻) at C3, an acetyl group (NHAc) at C4, and a hydroxyl group (OH) at C6.	GalNAc_bD
α D_GlcP [2S3Me6S]		α	 A chemical structure of a glucose residue in a chair conformation. It has a methyl group (H ₃ C) at C1, a hydroxyl group (OH) at C2, a sulfate group (-OSO ₃ ⁻) at C3, and two sulfate groups (-OSO ₃ ⁻) at C6.	Glc[2S3Me6S]_aD

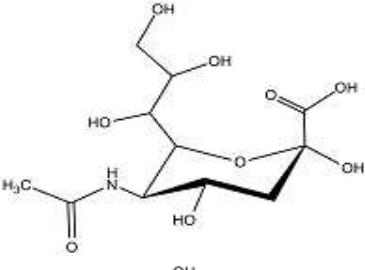
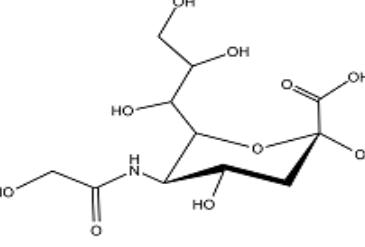
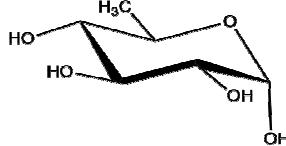
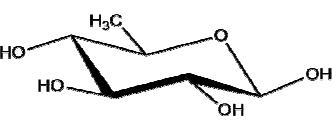
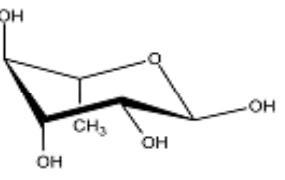
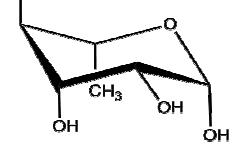
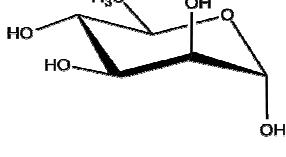
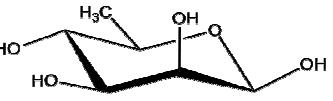
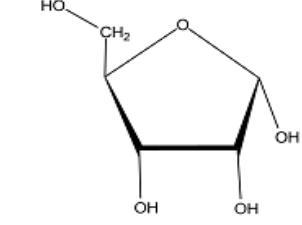
$\alpha_D\text{-Glc}p\ [2S3S6S]$	α 2S3S6S		Glc[2S3S6S]_aD
$\alpha_D\text{-Glc}p\ [2S6S]$	α 2S6S		Glc[2S6S]_aD
$\alpha_D\text{-Glc}p\ [3P]$	α 3P		Glc[3P]_aD
$\beta_D\text{-Glc}p\ [6S]$	β 6S		Glc[6S]_bD
$\alpha_D\text{-Glc}p$	α		Glc_aD
$\beta_D\text{-Glc}p$	β		Glc_bD
$\beta_D\text{-Glc}p\ [6Ac]$	β 6Ac		Glc[6Ac]_bD
$\alpha_D\text{-Glc}p\ [6P]$	α 6P		Glc[6P]_aD
$\beta_D\text{-Glc}p\ A[4S]$	β 4S		GlcA[4S]_bD
$\beta_D\text{-Glc}p\ A[6S]$	β 6S		GlcA[6S]_bD

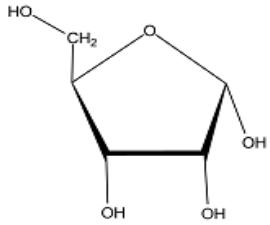
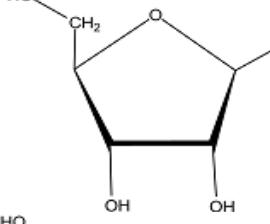
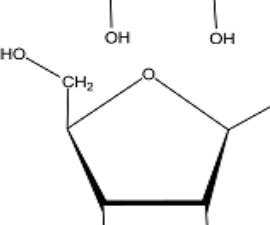
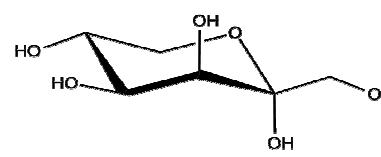
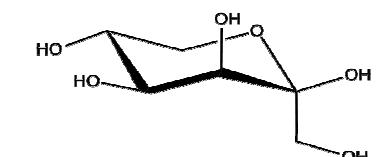
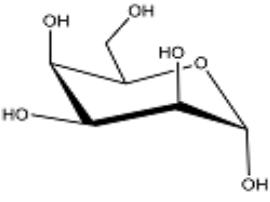
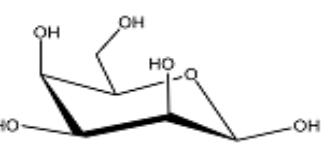
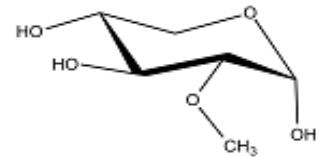
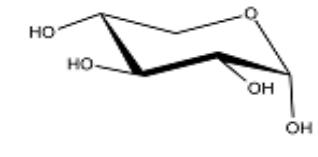
α D_GlcP A		α		GlcA_aD
β D_GlcP A		β		GlcA_bD
α D_GlcP N[6S]		α 6S		GlcN[6S]_aD
α D_GlcP N		α		GlcN_aD
β D_GlcP N		β		GlcN_bD
α D_GlcP NAc[3S]		α 3S		GlcNAc[3S]_aD
β D_GlcP NAc[3S]		β 3S		GlcNAc[3S]_bD
β D_GlcP NAc[3S6S]		β 3S6S		GlcNAc[3S6S]_bD
β D_GlcP NAc[4S]		β 4S		GlcNAc[4S]_bD
α D_GlcP NAc[6S]		α 6S		GlcNAc[6S]_aD

βD_Glcp NAc[6S]	 β 6S		GlcNAc[6S]_bD
αD_Glcp NAc	 α		GlcNAc_aD
βD_Glcp NAc	 β		GlcNAc_bD
aD_GlcpNS[3S]	 α 3S		GlcNS[3S]_aD
aD_GlcpNS[6S]	 α 6S		GlcNS[6S]_aD
αD_Glcp NS[3S6S]	 α 3S6S		GlcNS[3S6S]_aD
αD_Glcp NS	 α		GlcNS_aD
αL_Gulp A	 α		GulA_aL
αL_Idop A[2S]	 α 2S 1C4		IdoA[2S]_aL(1C4)

$\alpha L_IdoP\ A[2S]$			$IdoA[2S]_aL\ (2S0)$
$\alpha L_IdoP\ A$			$IdoA_aL(1C4)$
$\alpha L_IdoP\ A$			$IdoA_aL(2S0)$
α_KDN			KDN_a
α_KDO			KDO_a
αD_Lyxp			Lyx_aD
βD_Lyxp			Lyx_bD
αD_Manp			Man_aD
βD_Manp			Man_bD

α D_Manp [6Ac]			Man[6Ac]_aD
α D_Manp [6P]			Man[6P]_aD
β D_Manp[Pyr]			Man[Pyr]_bD
α D_ManpA			ManA_aD
β D_Manp A			ManA_bD
α D_Manp N			ManN_aD
β D_Manp N			ManN_bD
α D_Manp NAc			ManNAc_aD
β D_Manp NAc			ManNAc_bD

α D_Neu5Ac	 α		Neu5Ac_aD
α D_Neu5Gc	 α		Neu5Gc_aD
α D_Qui	 α		Qui_aD
β D_Qui	 β		Qui_bD
α L_Rhap	 α		Rha_aL
β L_Rhap	 β		Rha_bL
α D_Rhap	 α		aD_Rhap
β D_Rhap	 β		bD_Rhap
α D_Ribf	 N α		Rib_aD(N)

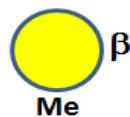
α D_Ribf			Rib_aD(S)
β D_Ribf			Rib_bD(N)
β D_Ribf			Rib_bD(S)
α D-Tagp			Tag_aD
β D-Tagp			Tag_bD
α D_Talp			Tal_aD
β D_Talp			Tal_bD
α D_Xylp [2Me]			Xyl[2Me]_aD
α D_Xylp			Xyl_aD

β D_Xylp



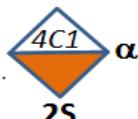
Xyl_bD

β D_Galp [Me]



Gal[Me]_bD

α L_Idop A[2S]



IdoA[2S]_aL (4C1)

α L_Idop A



IdoA_aL (4C1)