Full wwPDB X-ray Structure Validation Report (i)

Oct 18, 2023 – 12:06 AM EDT

PDB ID	:	2YVV
Title	:	Crystal structure of hyluranidase complexed with lactose at 2.6 A resolution reveals three specific sugar recognition sites
Authors	:	Mishra, P.; Prem Kumar, R.; Singh, N.; Sharma, S.; Kaur, P.; Perbandt, M.; Betzel, C.; Bhakuni, V.; Singh, T.P.
Deposited on	:	2007-04-16
Resolution	:	2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) were used in the production of this report:

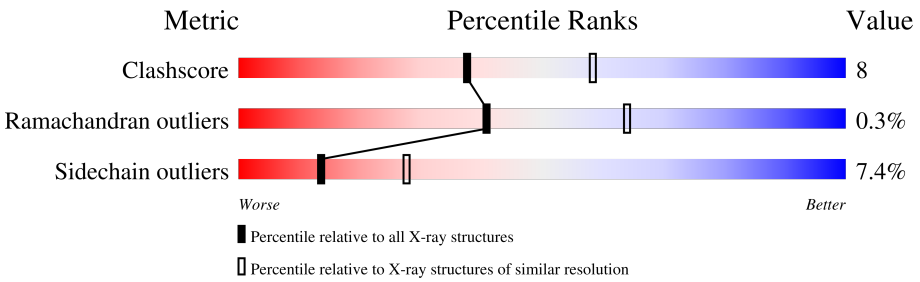
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	338	83% 11% • • •
2	B	2	100%
2	C	2	50% 50%
2	D	2	100%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2694 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Hyaluronidase, phage associated.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	332	2515	1565	439	506	5	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	338	HIS	-	expression tag	UNP Q9A0M7

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-alpha-D-glucopyranose.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	B	2	23	12	11	0	0	0
2	C	2	23	12	11	0	0	0
2	D	2	23	12	11	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	110	110	110	0	0

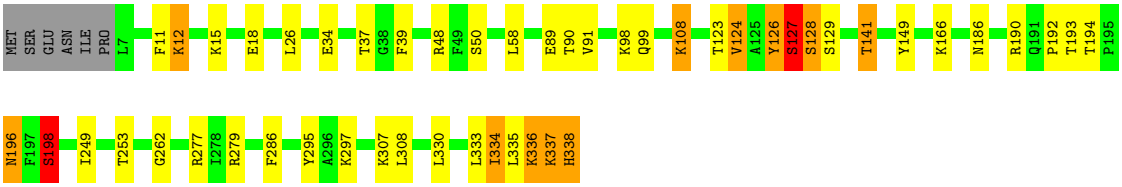
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

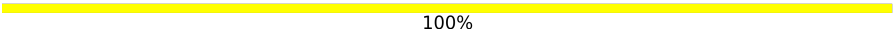
• Molecule 1: Hyaluronidase, phage associated

Chain A:



• Molecule 2: beta-D-galactopyranose-(1-4)-alpha-D-glucopyranose

Chain B:



• Molecule 2: beta-D-galactopyranose-(1-4)-alpha-D-glucopyranose

Chain C:

• Molecule 2: beta-D-galactopyranose-(1-4)-alpha-D-glucopyranose

Chain D:

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α, β, γ	59.06Å 59.06Å 588.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.60	Depositor
% Data completeness (in resolution range)	98.9 (50.00-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.191 , 0.226	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2694	wwPDB-VP
Average B, all atoms (Å^2)	47.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: GLC, GAL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\|>5$	RMSZ	$\#\|Z\|>5$
1	A	0.74	0/2549	0.93	8/3434 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	336	LYS	N-CA-C	14.30	149.60	111.00
1	A	337	LYS	N-CA-C	13.96	148.70	111.00
1	A	198	SER	C-N-CA	8.62	143.26	121.70
1	A	127	SER	CB-CA-C	8.38	126.02	110.10
1	A	198	SER	N-CA-C	-8.29	88.60	111.00
1	A	336	LYS	CB-CA-C	-7.63	95.14	110.40
1	A	279	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	A	127	SER	N-CA-C	-5.19	97.00	111.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	337	LYS	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	198	SER	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2515	0	2531	38	0
2	B	23	0	21	0	0
2	C	23	0	21	1	0
2	D	23	0	21	4	0
3	A	110	0	0	2	0
All	All	2694	0	2594	40	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (40) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:LYS:HE3	3:A:3028:HOH:O	1.50	1.10
1:A:126:TYR:HD2	1:A:127:SER:N	1.80	0.79
2:D:1:GLC:H62	2:D:2:GAL:O2	1.82	0.78
1:A:126:TYR:CD2	1:A:127:SER:N	2.57	0.72
1:A:98:LYS:HB2	1:A:98:LYS:NZ	2.07	0.70
1:A:48:ARG:CG	1:A:48:ARG:HH11	2.06	0.68
1:A:338:HIS:N	1:A:338:HIS:CD2	2.62	0.67
1:A:48:ARG:HH11	1:A:48:ARG:HG2	1.59	0.67
1:A:196:ASN:OD1	1:A:198:SER:HB2	2.04	0.58
1:A:48:ARG:CG	1:A:48:ARG:NH1	2.67	0.57
1:A:330:LEU:O	1:A:334:ILE:HB	2.04	0.57
1:A:333:LEU:O	1:A:336:LYS:HA	2.05	0.57
1:A:338:HIS:CD2	1:A:338:HIS:H	2.25	0.55
1:A:141:THR:HG22	3:A:3061:HOH:O	2.07	0.55
1:A:333:LEU:O	1:A:333:LEU:HD23	2.07	0.55
1:A:34:GLU:HB2	1:A:37:THR:OG1	2.07	0.54
1:A:98:LYS:HB2	1:A:98:LYS:HZ2	1.71	0.54
1:A:196:ASN:OD1	1:A:198:SER:N	2.40	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:LYS:NZ	1:A:98:LYS:CB	2.73	0.52
1:A:249:ILE:HD11	1:A:262:GLY:HA2	1.91	0.51
1:A:15:LYS:HB2	1:A:18:GLU:HG3	1.93	0.50
1:A:338:HIS:N	1:A:338:HIS:HD2	2.08	0.48
1:A:108:LYS:HE3	1:A:108:LYS:HB2	1.48	0.47
1:A:128:SER:HB3	1:A:129:SER:H	1.55	0.47
1:A:333:LEU:HD23	1:A:333:LEU:C	2.36	0.46
1:A:277:ARG:NH1	2:D:1:GLC:O6	2.50	0.44
2:D:1:GLC:C6	2:D:2:GAL:O2	2.60	0.44
1:A:337:LYS:CG	1:A:337:LYS:O	2.64	0.44
1:A:277:ARG:NH2	2:D:1:GLC:O1	2.50	0.43
1:A:286:PHE:HA	1:A:295:TYR:O	2.19	0.43
1:A:26:LEU:HD23	1:A:26:LEU:HA	1.77	0.42
1:A:90:THR:O	1:A:91:VAL:HG23	2.20	0.42
1:A:190:ARG:O	1:A:192:PRO:HD3	2.20	0.42
1:A:149:TYR:CD1	1:A:149:TYR:N	2.88	0.41
1:A:123:THR:O	1:A:124:VAL:HG23	2.20	0.41
1:A:98:LYS:HB2	1:A:98:LYS:HZ3	1.83	0.41
1:A:98:LYS:CB	1:A:98:LYS:HZ3	2.34	0.41
1:A:186:ASN:HD21	2:C:2:GAL:H61	1.85	0.40
1:A:11:PHE:O	1:A:12:LYS:C	2.60	0.40
1:A:37:THR:HB	1:A:39:PHE:CD2	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	330/338 (98%)	310 (94%)	19 (6%)	1 (0%)	41 64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	VAL

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	270/276 (98%)	250 (93%)	20 (7%)	13 28

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LYS
1	A	50	SER
1	A	58	LEU
1	A	89	GLU
1	A	99	GLN
1	A	108	LYS
1	A	126	TYR
1	A	127	SER
1	A	128	SER
1	A	141	THR
1	A	166	LYS
1	A	193	THR
1	A	194	THR
1	A	196	ASN
1	A	253	THR
1	A	307	LYS
1	A	308	LEU
1	A	334	ILE
1	A	335	LEU
1	A	338	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	GLN

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Mol	Chain	Res	Type
1	A	186	ASN
1	A	202	ASN
1	A	338	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

6 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	GLC	B	1	2	12,12,12	1.74	3 (25%)	17,17,17	1.77	2 (11%)
2	GAL	B	2	2	11,11,12	2.44	7 (63%)	15,15,17	1.93	6 (40%)
2	GLC	C	1	2	12,12,12	0.60	0	17,17,17	1.14	2 (11%)
2	GAL	C	2	2	11,11,12	0.72	0	15,15,17	1.07	1 (6%)
2	GLC	D	1	2	12,12,12	0.58	0	17,17,17	1.06	1 (5%)
2	GAL	D	2	2	11,11,12	0.57	0	15,15,17	0.97	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	B	1	2	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GAL	B	2	2	-	2/2/19/22	0/1/1/1
2	GLC	C	1	2	-	2/2/22/22	0/1/1/1
2	GAL	C	2	2	-	2/2/19/22	0/1/1/1
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GAL	D	2	2	-	2/2/19/22	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	GAL	O3-C3	3.93	1.52	1.43
2	B	2	GAL	O5-C1	3.57	1.49	1.43
2	B	2	GAL	O5-C5	3.19	1.49	1.43
2	B	1	GLC	C3-C2	3.18	1.60	1.52
2	B	2	GAL	C4-C5	2.86	1.59	1.53
2	B	2	GAL	C1-C2	2.58	1.58	1.52
2	B	1	GLC	O4-C4	2.51	1.48	1.43
2	B	2	GAL	C4-C3	2.31	1.58	1.52
2	B	2	GAL	C6-C5	2.16	1.59	1.51
2	B	1	GLC	O1-C1	2.08	1.46	1.39

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1	GLC	O4-C4-C3	6.04	124.30	110.35
2	B	2	GAL	C2-C3-C4	-3.26	105.26	110.89
2	B	2	GAL	O5-C1-C2	-2.81	106.43	110.77
2	B	2	GAL	O5-C5-C4	-2.76	104.11	110.83
2	B	2	GAL	O5-C5-C6	2.53	111.17	107.20
2	B	2	GAL	C1-C2-C3	2.52	112.77	109.67
2	B	2	GAL	C6-C5-C4	-2.50	107.14	113.00
2	C	1	GLC	C3-C4-C5	-2.46	105.84	110.24
2	D	1	GLC	O4-C4-C3	-2.31	105.02	110.35
2	C	2	GAL	O5-C1-C2	-2.18	107.41	110.77
2	C	1	GLC	O5-C5-C6	2.13	111.73	106.44
2	D	2	GAL	C1-O5-C5	2.10	115.03	112.19
2	B	1	GLC	C4-C3-C2	-2.06	107.23	110.82

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	2	GAL	C4-C5-C6-O6
2	B	2	GAL	O5-C5-C6-O6
2	D	2	GAL	C4-C5-C6-O6
2	C	2	GAL	O5-C5-C6-O6
2	D	2	GAL	O5-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	C	2	GAL	C4-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6

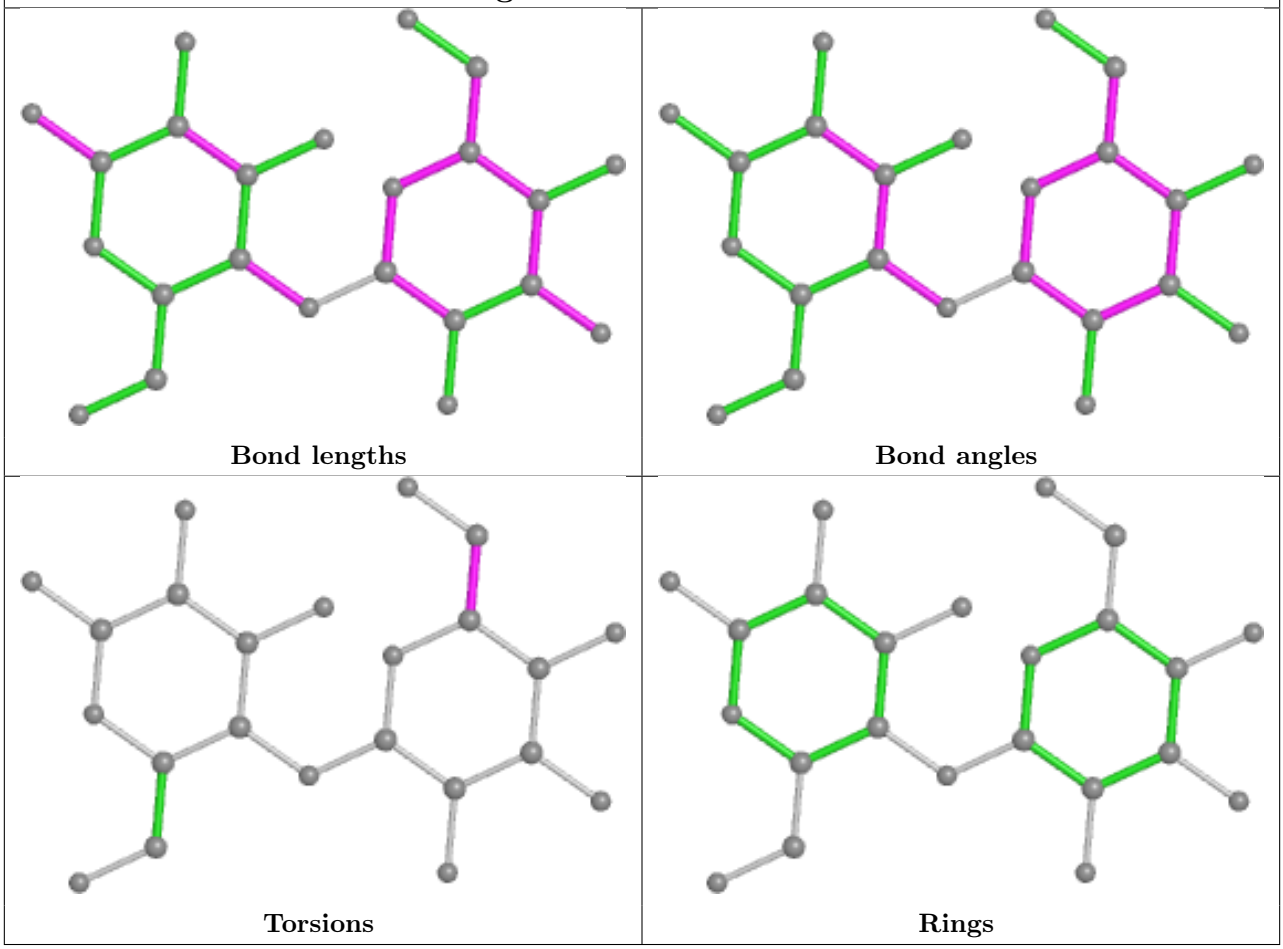
There are no ring outliers.

3 monomers are involved in 5 short contacts:

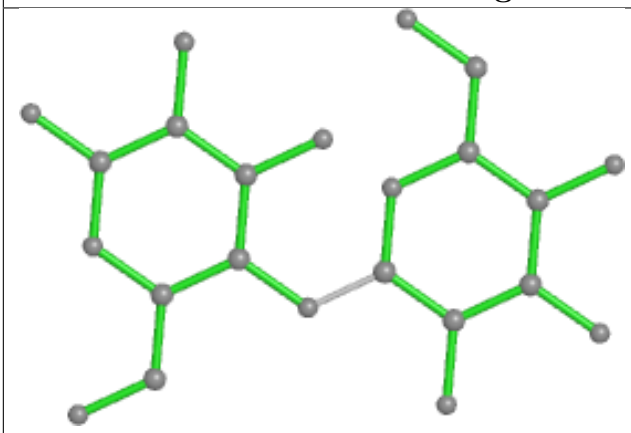
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	GAL	2	0
2	C	2	GAL	1	0
2	D	1	GLC	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

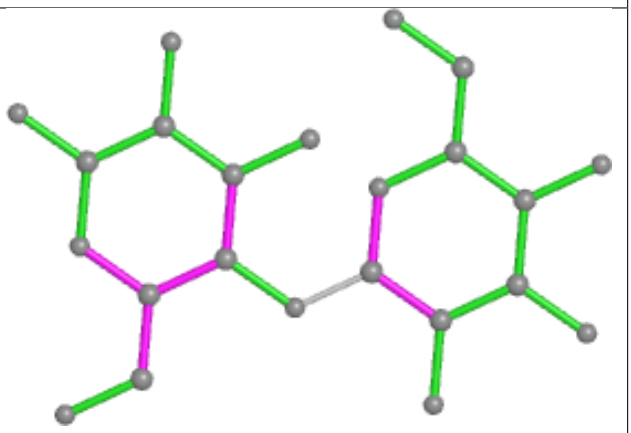
Oligosaccharide Chain B



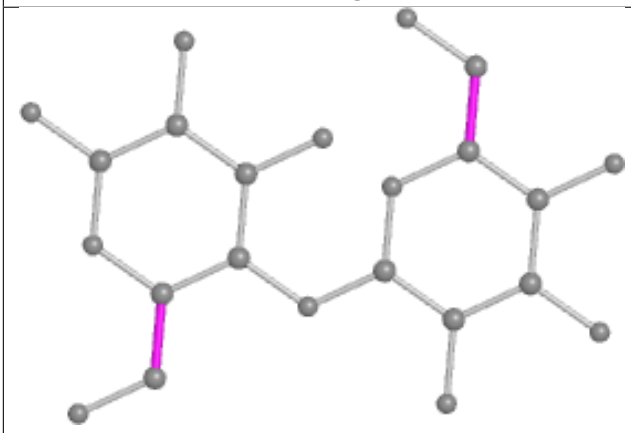
Oligosaccharide Chain C



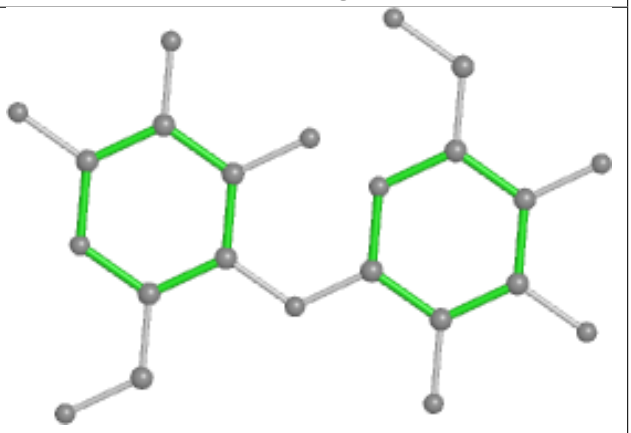
Bond lengths



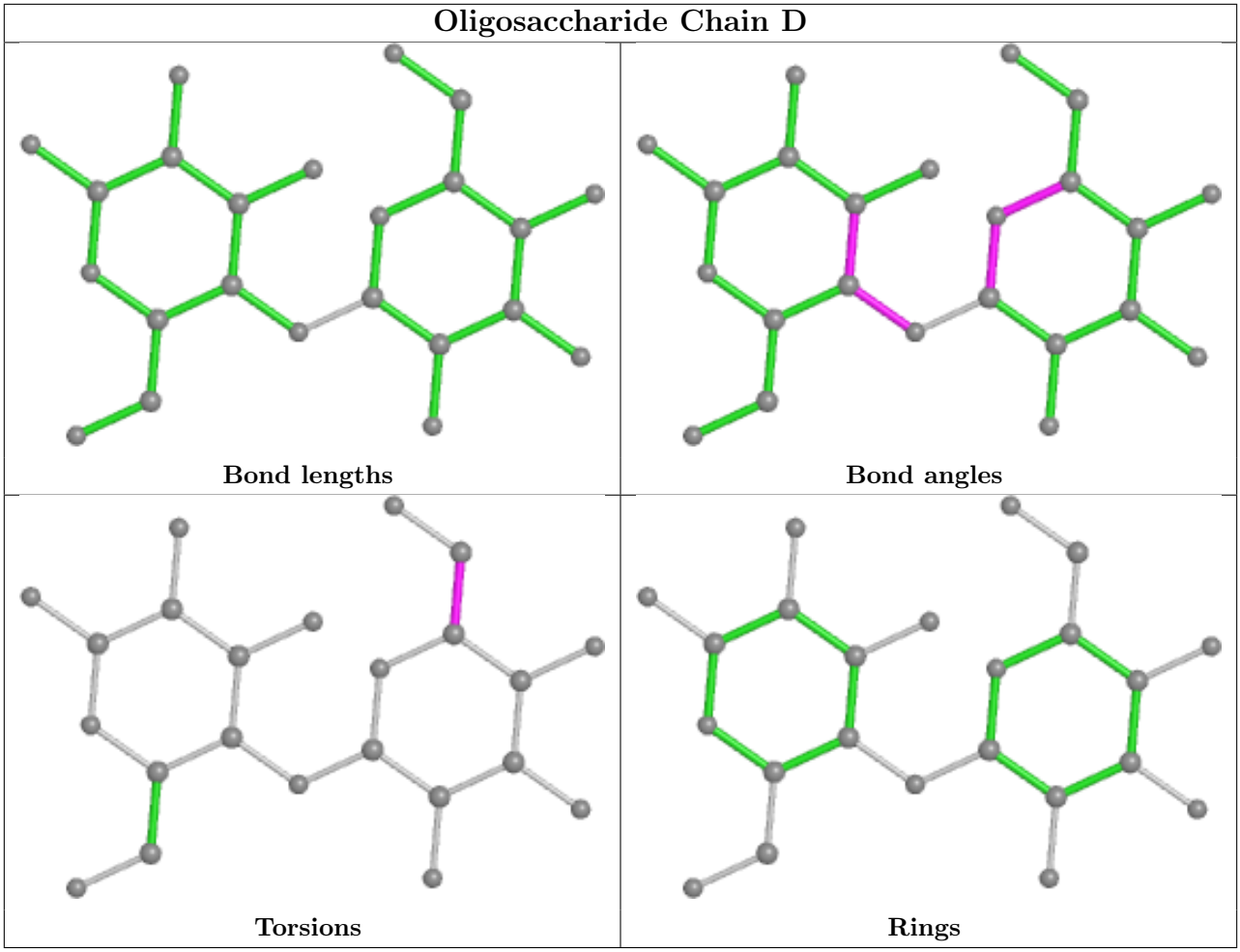
Bond angles



Torsions



Rings



5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.