



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:18 PM UTC

PDB ID : 6X3M / pdb_00006x3m
Title : Crystal structure of full-length Streptococcal bacteriophage hyaluronidase in complex with unsaturated hyaluronan octa-saccharides
Authors : Deivanayagam, C.; Schormann, N.
Deposited on : 2020-05-21
Resolution : 3.58 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

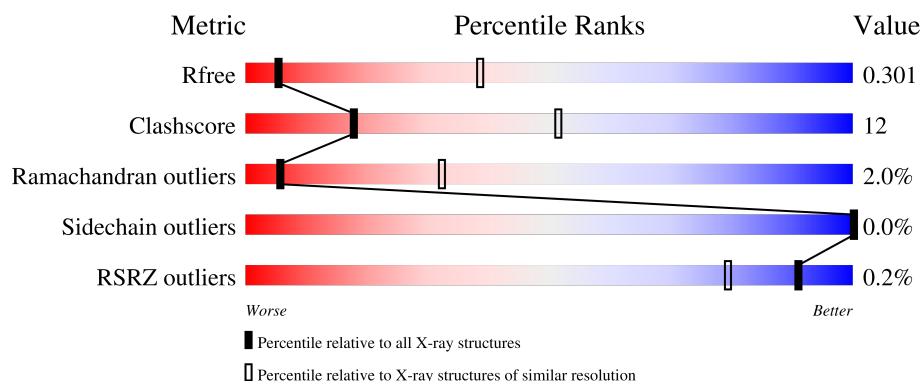
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.58 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.



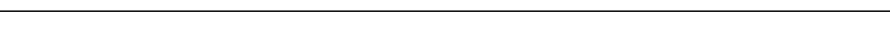
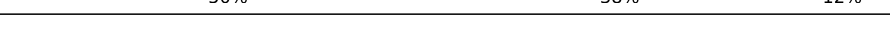
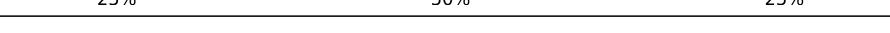
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1521 (3.66-3.50)
Clashscore	190562	1595 (3.66-3.50)
Ramachandran outliers	187476	1551 (3.66-3.50)
Sidechain outliers	187428	1551 (3.66-3.50)
RSRZ outliers	180081	1520 (3.66-3.50)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	379	 69% 25% • 5%
1	B	379	 72% 22% • 5%
1	C	379	 74% 21% 5%
1	D	379	 65% 26% • 8%
1	E	379	 68% 25% • 6%

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Mol	Chain	Length	Quality of chain
1	F	379	
1	G	379	
1	H	379	
1	I	379	
2	J	4	
2	N	4	
2	O	4	
2	Q	4	
2	T	4	
2	U	4	
2	V	4	
2	Z	4	
2	a	4	
3	K	8	
3	L	8	
3	M	8	
3	P	8	
3	R	8	
3	S	8	
3	W	8	
3	X	8	
3	Y	8	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	M	1	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 25574 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 Hyaluronoglucosaminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2693	1671	469	545	8			
1	B	361	Total	C	N	O	S	0	0	0
			2692	1670	469	545	8			
1	C	361	Total	C	N	O	S	0	0	0
			2692	1670	469	545	8			
1	D	349	Total	C	N	O	S	0	0	0
			2599	1610	454	527	8			
1	E	357	Total	C	N	O	S	0	0	0
			2662	1650	464	540	8			
1	F	366	Total	C	N	O	S	0	0	0
			2733	1697	475	553	8			
1	G	363	Total	C	N	O	S	0	0	0
			2708	1682	471	547	8			
1	H	361	Total	C	N	O	S	0	0	0
			2693	1673	468	544	8			
1	I	362	Total	C	N	O	S	0	0	0
			2700	1676	470	546	8			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	372	LEU	-	expression tag	UNP P15316
A	373	GLU	-	expression tag	UNP P15316
A	374	HIS	-	expression tag	UNP P15316
A	375	HIS	-	expression tag	UNP P15316
A	376	HIS	-	expression tag	UNP P15316
A	377	HIS	-	expression tag	UNP P15316
A	378	HIS	-	expression tag	UNP P15316
A	379	HIS	-	expression tag	UNP P15316
B	372	LEU	-	expression tag	UNP P15316
B	373	GLU	-	expression tag	UNP P15316
B	374	HIS	-	expression tag	UNP P15316

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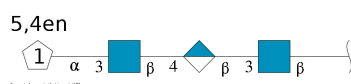
Chain	Residue	Modelled	Actual	Comment	Reference
B	375	HIS	-	expression tag	UNP P15316
B	376	HIS	-	expression tag	UNP P15316
B	377	HIS	-	expression tag	UNP P15316
B	378	HIS	-	expression tag	UNP P15316
B	379	HIS	-	expression tag	UNP P15316
C	372	LEU	-	expression tag	UNP P15316
C	373	GLU	-	expression tag	UNP P15316
C	374	HIS	-	expression tag	UNP P15316
C	375	HIS	-	expression tag	UNP P15316
C	376	HIS	-	expression tag	UNP P15316
C	377	HIS	-	expression tag	UNP P15316
C	378	HIS	-	expression tag	UNP P15316
C	379	HIS	-	expression tag	UNP P15316
D	372	LEU	-	expression tag	UNP P15316
D	373	GLU	-	expression tag	UNP P15316
D	374	HIS	-	expression tag	UNP P15316
D	375	HIS	-	expression tag	UNP P15316
D	376	HIS	-	expression tag	UNP P15316
D	377	HIS	-	expression tag	UNP P15316
D	378	HIS	-	expression tag	UNP P15316
D	379	HIS	-	expression tag	UNP P15316
E	372	LEU	-	expression tag	UNP P15316
E	373	GLU	-	expression tag	UNP P15316
E	374	HIS	-	expression tag	UNP P15316
E	375	HIS	-	expression tag	UNP P15316
E	376	HIS	-	expression tag	UNP P15316
E	377	HIS	-	expression tag	UNP P15316
E	378	HIS	-	expression tag	UNP P15316
E	379	HIS	-	expression tag	UNP P15316
F	372	LEU	-	expression tag	UNP P15316
F	373	GLU	-	expression tag	UNP P15316
F	374	HIS	-	expression tag	UNP P15316
F	375	HIS	-	expression tag	UNP P15316
F	376	HIS	-	expression tag	UNP P15316
F	377	HIS	-	expression tag	UNP P15316
F	378	HIS	-	expression tag	UNP P15316
F	379	HIS	-	expression tag	UNP P15316
G	372	LEU	-	expression tag	UNP P15316
G	373	GLU	-	expression tag	UNP P15316
G	374	HIS	-	expression tag	UNP P15316
G	375	HIS	-	expression tag	UNP P15316
G	376	HIS	-	expression tag	UNP P15316

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Chain	Residue	Modelled	Actual	Comment	Reference
G	377	HIS	-	expression tag	UNP P15316
G	378	HIS	-	expression tag	UNP P15316
G	379	HIS	-	expression tag	UNP P15316
H	372	LEU	-	expression tag	UNP P15316
H	373	GLU	-	expression tag	UNP P15316
H	374	HIS	-	expression tag	UNP P15316
H	375	HIS	-	expression tag	UNP P15316
H	376	HIS	-	expression tag	UNP P15316
H	377	HIS	-	expression tag	UNP P15316
H	378	HIS	-	expression tag	UNP P15316
H	379	HIS	-	expression tag	UNP P15316
I	372	LEU	-	expression tag	UNP P15316
I	373	GLU	-	expression tag	UNP P15316
I	374	HIS	-	expression tag	UNP P15316
I	375	HIS	-	expression tag	UNP P15316
I	376	HIS	-	expression tag	UNP P15316
I	377	HIS	-	expression tag	UNP P15316
I	378	HIS	-	expression tag	UNP P15316
I	379	HIS	-	expression tag	UNP P15316

- Molecule 2 is an oligosaccharide called 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose.



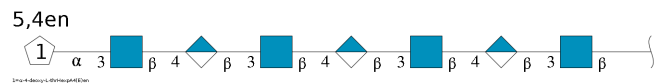
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	J	4	Total	C	N	O	0	0	0
			51	28	2	21			
2	N	4	Total	C	N	O	0	0	0
			52	28	2	22			
2	O	4	Total	C	N	O	0	0	0
			52	28	2	22			
2	Q	4	Total	C	N	O	0	0	0
			52	28	2	22			
2	T	4	Total	C	N	O	0	0	0
			52	28	2	22			
2	U	4	Total	C	N	O	0	0	0
			52	28	2	22			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	V	4	Total	C	N	O	0	0	0
			51	28	2	21			
2	Z	4	Total	C	N	O	0	0	0
			52	28	2	22			
2	a	4	Total	C	N	O	0	0	0
			52	28	2	22			

- Molecule 3 is an oligosaccharide called 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose.

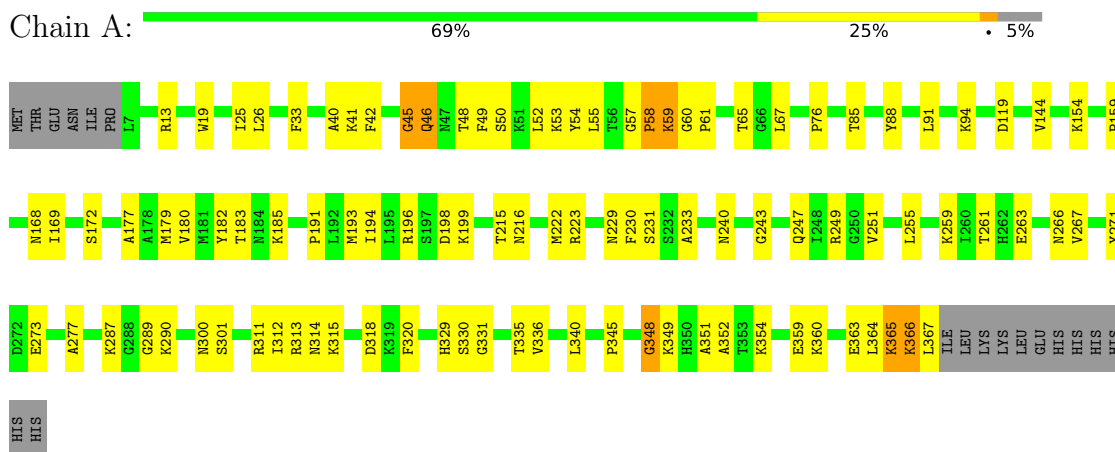


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	K	8	Total	C	N	O	0	0	0
			104	56	4	44			
3	L	8	Total	C	N	O	0	0	0
			104	56	4	44			
3	M	8	Total	C	N	O	0	0	0
			104	56	4	44			
3	P	8	Total	C	N	O	0	0	0
			104	56	4	44			
3	R	8	Total	C	N	O	0	0	0
			104	56	4	44			
3	S	8	Total	C	N	O	0	0	0
			104	56	4	44			
3	W	8	Total	C	N	O	0	0	0
			104	56	4	44			
3	X	8	Total	C	N	O	0	0	0
			104	56	4	44			
3	Y	8	Total	C	N	O	0	0	0
			104	56	4	44			

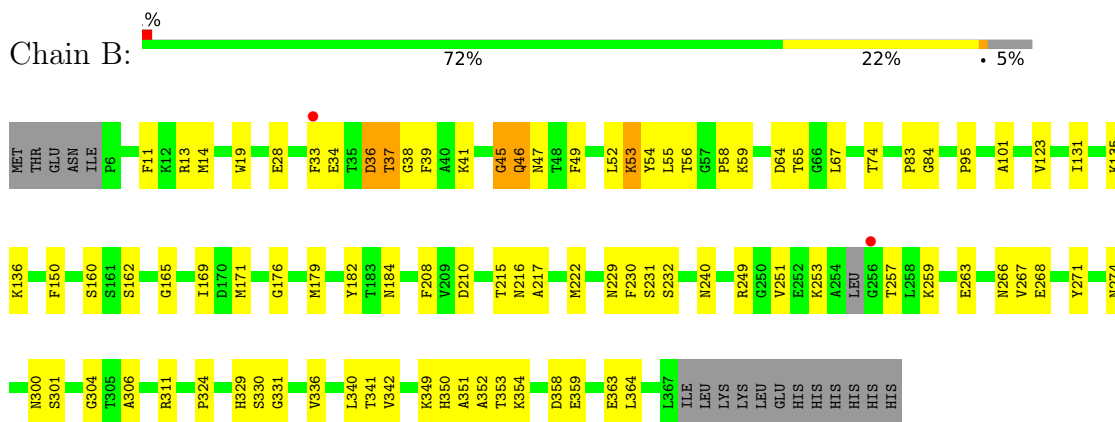
3 Residue-property plots [i](#)

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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

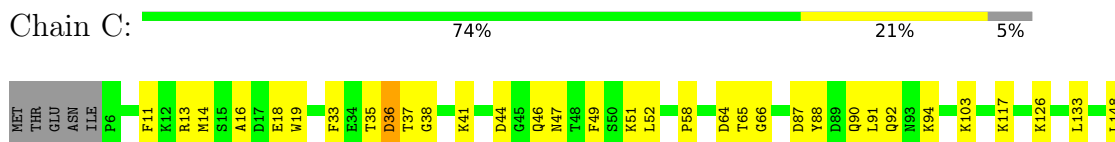
• Molecule 1: Hyaluronoglucosaminidase

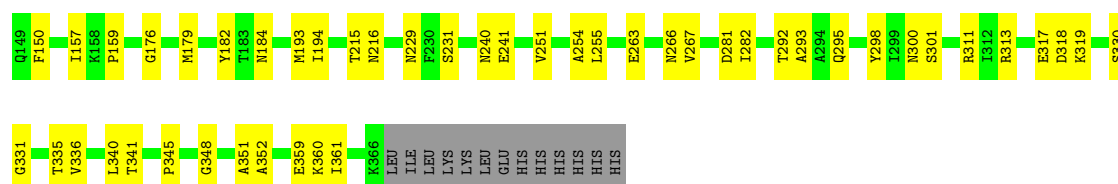


• Molecule 1: Hyaluronoglucosaminidase



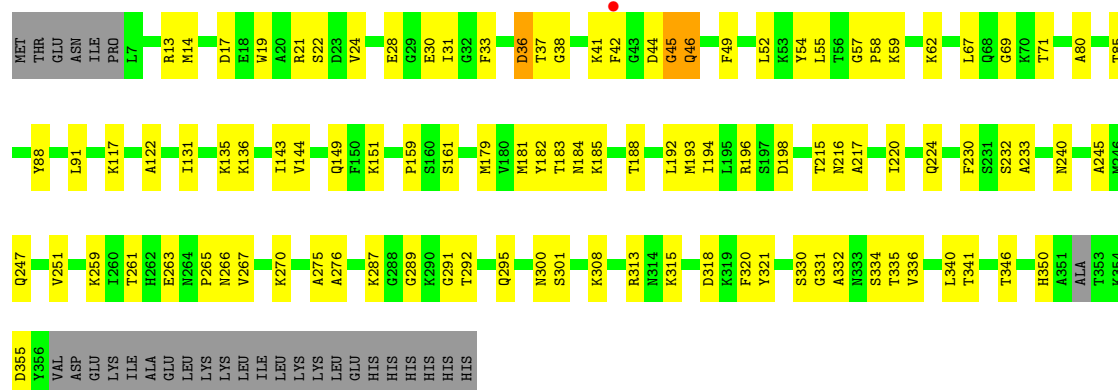
• Molecule 1: Hyaluronoglucosaminidase





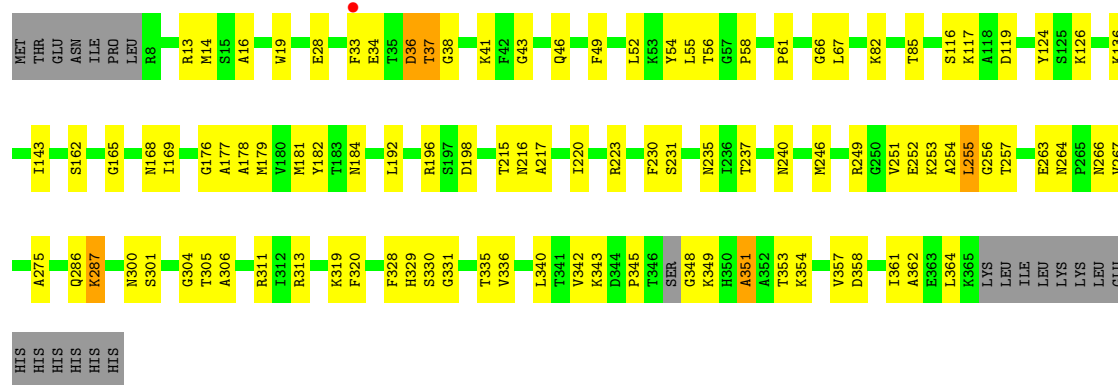
• Molecule 1: Hyaluronoglucosaminidase

Chain D: 65% 26% 8%



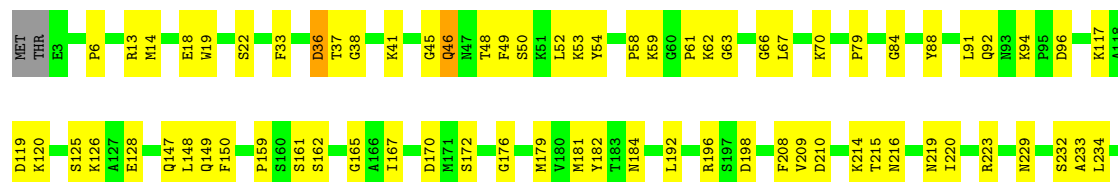
• Molecule 1: Hyaluronoglucosaminidase

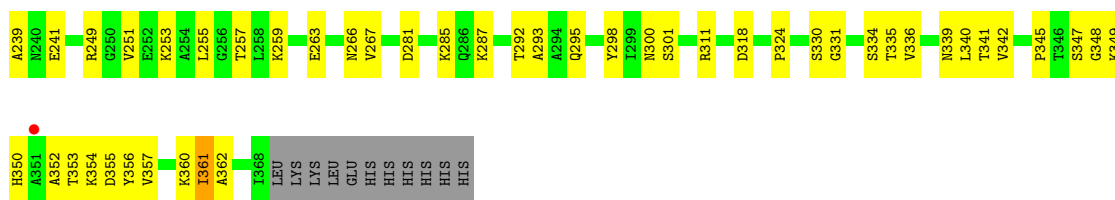
Chain E: 68% 25% 6%



• Molecule 1: Hyaluronoglucosaminidase

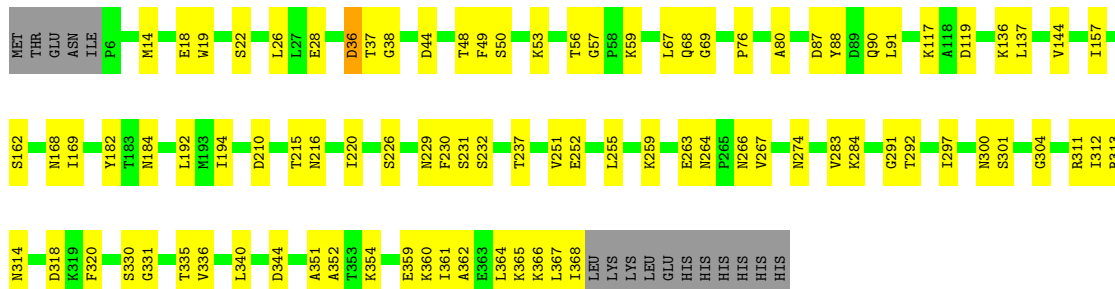
Chain F: 65% 30% 5%





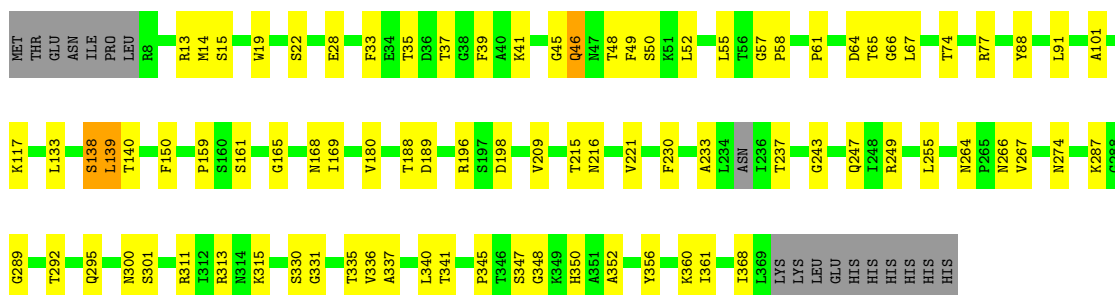
- Molecule 1: Hyaluronoglucosaminidase

Chain G: 72% 23%



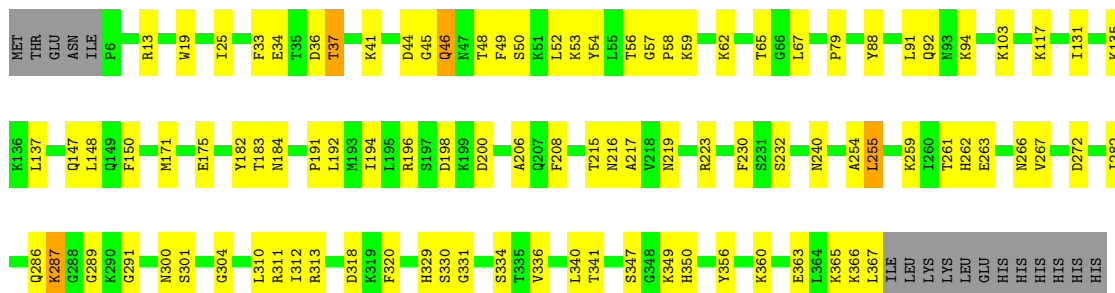
- Molecule 1: Hyaluronoglucosaminidase

Chain H: 73% 22% 5%



- Molecule 1: Hyaluronoglucosaminidase

Chain I: 70% 25%



- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  75% 25%



- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  25% 25% 50%



- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  25% 25% 50%



- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  75% 25%



- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:  100%



- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain U:  50% 50%



- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

anose

Chain V: 

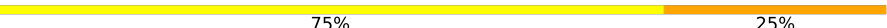


- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z: 



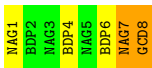
- Molecule 2: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain a: 



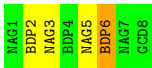
- Molecule 3: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K: 

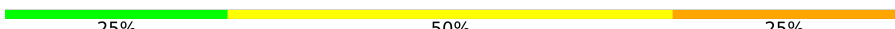


- Molecule 3: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L: 



- Molecule 3: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M: 



• Molecule 3: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose



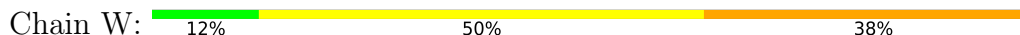
• Molecule 3: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose



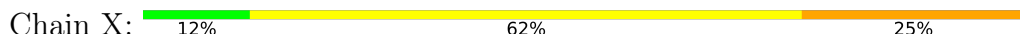
• Molecule 3: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 3: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 3: 4-deoxy-alpha-L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1	BDP2	MAG3	BDP4	MAG5	BDP6	MAG7	GCD8
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- Molecule 3: 4-deoxy- α -L-threo-hex-4-enopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Y: 

MAG1	BDP2	MAG3	BDP4	MAG5	BDP6	MAG7	GCD8
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	338.17Å 56.80Å 276.95Å 90.00° 126.68° 90.00°	Depositor
Resolution (Å)	29.87 – 3.58 29.87 – 3.58	Depositor EDS
% Data completeness (in resolution range)	95.8 (29.87-3.58) 95.6 (29.87-3.58)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 3.56Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.247 , 0.299 0.249 , 0.301	Depositor DCC
R_{free} test set	2519 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	91.9	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 51.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	25574	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GCD, BDP, NAG

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	0.14	0/2732	0.43	0/3670
1	B	0.14	0/2731	0.41	0/3667
1	C	0.16	0/2732	0.44	0/3670
1	D	0.15	0/2637	0.44	0/3542
1	E	0.14	0/2700	0.44	0/3626
1	F	0.14	0/2773	0.46	0/3727
1	G	0.15	0/2748	0.44	0/3692
1	H	0.13	0/2731	0.42	0/3667
1	I	0.18	0/2740	0.45	1/3681 (0.0%)
All	All	0.15	0/24524	0.44	1/32942 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	46	GLN	CA-CB-CG	5.01	124.13	114.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	2693	0	2702	95	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2692	0	2698	89	0
1	C	2692	0	2699	77	0
1	D	2599	0	2594	95	0
1	E	2662	0	2661	101	0
1	F	2733	0	2743	108	0
1	G	2708	0	2721	93	0
1	H	2693	0	2706	83	0
1	I	2700	0	2710	89	0
2	J	51	0	35	0	0
2	N	52	0	37	3	0
2	O	52	0	37	6	0
2	Q	52	0	37	0	0
2	T	52	0	37	4	0
2	U	52	0	37	6	0
2	V	51	0	35	4	0
2	Z	52	0	37	2	0
2	a	52	0	38	2	0
3	K	104	0	73	5	0
3	L	104	0	73	1	0
3	M	104	0	73	8	0
3	P	104	0	73	11	0
3	R	104	0	73	7	0
3	S	104	0	73	6	0
3	W	104	0	73	7	0
3	X	104	0	73	6	0
3	Y	104	0	73	10	0
All	All	25574	0	25221	627	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 12.

All (627) 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:55:LEU:HD12	1:A:55:LEU:O	1.30	1.30
1:G:364:LEU:HD22	1:H:361:ILE:HG23	1.36	1.01
1:A:55:LEU:O	1:A:55:LEU:CD1	2.13	0.96
1:A:364:LEU:HD12	1:A:364:LEU:O	1.72	0.87
1:H:237:THR:HG21	3:Y:3:NAG:H62	1.56	0.85
1:H:274:ASN:O	2:Z:3:NAG:H83	1.81	0.80
1:E:49:PHE:HA	1:E:52:LEU:HD12	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:340:LEU:HD11	1:F:336:VAL:HG23	1.66	0.77
1:G:292:THR:HG22	1:H:264:ASN:HD22	1.48	0.77
1:A:336:VAL:HG23	1:C:340:LEU:HD11	1.68	0.76
1:C:14:MET:HE2	1:C:18:GLU:HB3	1.68	0.76
1:E:19:TRP:NE1	1:E:33:PHE:O	2.18	0.76
1:C:49:PHE:HA	1:C:52:LEU:HD12	1.68	0.75
1:I:88:TYR:HA	1:I:91:LEU:HD12	1.68	0.75
1:A:311:ARG:NH2	1:A:318:ASP:OD2	2.20	0.75
1:I:19:TRP:NE1	1:I:33:PHE:O	2.20	0.74
1:A:41:LYS:HD2	1:A:52:LEU:HD12	1.70	0.73
1:B:19:TRP:NE1	1:B:33:PHE:O	2.21	0.73
1:B:49:PHE:HA	1:B:52:LEU:HD12	1.70	0.73
1:G:292:THR:HG22	1:H:264:ASN:HB2	1.70	0.73
1:D:42:PHE:CZ	1:D:55:LEU:HD13	2.23	0.73
1:G:335:THR:HG23	1:I:341:THR:HB	1.70	0.73
1:F:41:LYS:HD2	1:F:52:LEU:HD12	1.69	0.72
1:H:58:PRO:HG2	1:I:59:LYS:HG3	1.70	0.72
1:B:259:LYS:HG2	3:M:1:NAG:C8	2.20	0.72
1:D:263:GLU:OE1	1:E:249:ARG:NH1	2.24	0.71
1:H:19:TRP:NE1	1:H:33:PHE:O	2.23	0.71
1:D:136:LYS:NZ	1:E:136:LYS:O	2.22	0.71
1:B:259:LYS:CG	3:M:1:NAG:H81	2.21	0.71
1:D:196:ARG:NH2	1:D:198:ASP:OD2	2.23	0.71
1:B:274:ASN:HB3	2:O:3:NAG:H62	1.72	0.70
1:F:311:ARG:NH2	1:F:318:ASP:OD2	2.23	0.70
1:H:49:PHE:HA	1:H:52:LEU:HD12	1.74	0.70
1:I:259:LYS:HG3	3:Y:1:NAG:H83	1.74	0.70
1:D:247:GLN:HB3	3:P:1:NAG:H83	1.73	0.70
1:A:229:ASN:O	1:C:240:ASN:ND2	2.21	0.69
1:G:80:ALA:O	1:H:77:ARG:NH1	2.25	0.69
1:B:11:PHE:O	1:B:13:ARG:NH1	2.25	0.69
1:A:85:THR:HG23	1:C:92:GLN:HG3	1.74	0.69
1:C:88:TYR:HA	1:C:91:LEU:HD12	1.75	0.69
1:A:196:ARG:NH2	1:A:198:ASP:OD2	2.24	0.69
2:a:1:NAG:O7	2:a:1:NAG:O1	2.10	0.68
1:E:182:TYR:HE1	1:E:184:ASN:HB2	1.59	0.68
1:G:36:ASP:O	1:G:38:GLY:N	2.26	0.68
1:G:263:GLU:OE1	1:H:249:ARG:NH1	2.27	0.67
1:D:240:ASN:ND2	1:E:231:SER:OG	2.28	0.67
1:I:196:ARG:NH2	1:I:198:ASP:OD2	2.27	0.67
1:I:261:THR:HB	3:Y:1:NAG:H81	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:GLY:HA2	1:C:159:PRO:HB3	1.76	0.66
1:D:88:TYR:HA	1:D:91:LEU:HD12	1.76	0.66
1:F:36:ASP:O	1:F:38:GLY:N	2.27	0.66
1:G:76:PRO:HD2	1:H:74:THR:HA	1.77	0.66
1:H:340:LEU:HD11	1:I:336:VAL:HG23	1.78	0.66
1:I:44:ASP:O	1:I:46:GLN:N	2.29	0.66
1:D:36:ASP:O	1:D:38:GLY:N	2.30	0.65
1:A:40:ALA:HB3	1:A:55:LEU:HD21	1.79	0.65
1:D:57:GLY:HA3	1:E:55:LEU:O	1.97	0.65
1:B:182:TYR:HE1	1:B:184:ASN:HB2	1.62	0.65
1:C:36:ASP:O	1:C:38:GLY:N	2.29	0.65
1:I:286:GLN:O	1:I:287:LYS:HG3	1.96	0.65
1:G:231:SER:OG	1:I:240:ASN:ND2	2.29	0.65
1:G:284:LYS:HA	1:G:292:THR:CG2	2.26	0.65
1:D:259:LYS:NZ	1:D:261:THR:OG1	2.30	0.65
1:C:44:ASP:HB2	1:C:52:LEU:HD21	1.79	0.64
1:C:19:TRP:NE1	1:C:33:PHE:O	2.30	0.64
1:A:183:THR:HG22	1:A:185:LYS:H	1.63	0.64
1:F:196:ARG:NH2	1:F:198:ASP:OD2	2.26	0.64
1:D:259:LYS:HG2	3:S:1:NAG:H82	1.80	0.64
1:G:336:VAL:HG23	1:I:340:LEU:HD11	1.79	0.64
1:A:340:LEU:HD11	1:B:336:VAL:HG23	1.80	0.64
1:A:364:LEU:O	1:A:366:LYS:N	2.30	0.63
1:E:169:ILE:HA	1:F:150:PHE:HB2	1.81	0.63
1:A:365:LYS:O	1:A:367:LEU:N	2.30	0.63
1:D:159:PRO:HB3	1:F:176:GLY:HA2	1.80	0.63
1:E:313:ARG:NE	2:U:1:NAG:O1	2.31	0.63
1:G:229:ASN:O	1:I:240:ASN:ND2	2.30	0.63
1:A:41:LYS:HE2	1:A:54:TYR:CZ	2.33	0.63
1:A:58:PRO:HD2	1:B:56:THR:HA	1.79	0.63
1:A:19:TRP:NE1	1:A:33:PHE:O	2.31	0.63
1:E:286:GLN:O	1:E:287:LYS:HG3	1.98	0.63
1:A:345:PRO:HG3	1:A:352:ALA:HA	1.80	0.63
1:B:259:LYS:HG2	3:M:1:NAG:H81	1.80	0.63
1:H:14:MET:SD	1:H:22:SER:OG	2.52	0.63
1:F:334:SER:OG	1:F:335:THR:N	2.24	0.62
1:D:71:THR:HG22	1:F:70:LYS:HB2	1.79	0.62
1:G:14:MET:HB2	1:G:19:TRP:CD1	2.34	0.62
1:H:295:GLN:NE2	2:a:2:BDP:O6B	2.31	0.62
1:D:41:LYS:HD2	1:D:52:LEU:HD12	1.80	0.62
1:F:219:ASN:ND2	3:P:3:NAG:O7	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:300:ASN:OD1	1:B:301:SER:N	2.32	0.62
1:B:46:GLN:O	1:B:47:ASN:ND2	2.32	0.62
1:F:259:LYS:NZ	2:T:4:GCD:O6A	2.32	0.61
1:H:311:ARG:HH12	2:V:1:NAG:H4	1.65	0.61
1:G:59:LYS:HG3	1:I:58:PRO:HG2	1.82	0.61
2:T:3:NAG:H4	2:T:4:GCD:H2	1.82	0.61
1:B:259:LYS:HG3	3:M:1:NAG:H81	1.82	0.61
1:A:169:ILE:HA	1:B:150:PHE:HB2	1.82	0.61
1:I:363:GLU:OE2	1:I:363:GLU:N	2.31	0.61
1:H:300:ASN:OD1	1:H:301:SER:N	2.32	0.61
1:G:169:ILE:HA	1:H:150:PHE:HB2	1.81	0.61
1:E:354:LYS:HA	1:E:357:VAL:HG22	1.83	0.61
1:G:136:LYS:NZ	1:H:133:LEU:O	2.34	0.61
1:D:185:LYS:HD3	1:D:188:THR:HB	1.83	0.60
1:I:266:ASN:OD1	1:I:267:VAL:N	2.34	0.60
1:G:88:TYR:HA	1:G:91:LEU:HD12	1.83	0.60
1:A:263:GLU:OE1	1:B:249:ARG:NH1	2.35	0.60
1:G:364:LEU:CD2	1:H:361:ILE:HG23	2.22	0.60
1:G:182:TYR:HE1	1:G:184:ASN:HB2	1.67	0.60
1:A:300:ASN:OD1	1:A:301:SER:N	2.34	0.60
1:B:34:GLU:OE1	1:B:37:THR:OG1	2.20	0.60
1:F:48:THR:HG23	1:F:50:SER:H	1.67	0.60
1:A:271:TYR:CZ	1:B:253:LYS:HE3	2.37	0.59
1:B:266:ASN:OD1	1:B:267:VAL:N	2.35	0.59
1:F:14:MET:HE1	1:F:22:SER:HB2	1.83	0.59
3:S:4:BDP:O3	3:S:5:NAG:C7	2.50	0.59
1:B:311:ARG:NH1	2:N:2:BDP:O6A	2.35	0.59
1:D:41:LYS:HE2	1:D:54:TYR:CZ	2.37	0.59
1:F:14:MET:HE2	1:F:18:GLU:HB3	1.83	0.59
1:D:117:LYS:NZ	1:E:119:ASP:OD1	2.36	0.59
1:E:263:GLU:HB3	1:F:251:VAL:HG12	1.84	0.59
3:R:1:NAG:O3	3:R:2:BDP:O6A	2.20	0.59
1:C:254:ALA:HB1	2:O:3:NAG:H61	1.84	0.59
1:E:176:GLY:HA2	1:F:159:PRO:HB3	1.84	0.59
1:F:352:ALA:HB3	1:F:356:TYR:CD2	2.37	0.59
1:G:220:ILE:HA	1:H:209:VAL:HB	1.85	0.58
1:H:345:PRO:HG3	1:H:352:ALA:HA	1.84	0.58
1:H:289:GLY:HA2	1:I:267:VAL:HA	1.85	0.58
1:C:300:ASN:OD1	1:C:301:SER:N	2.36	0.58
1:D:149:GLN:HB2	1:E:143:ILE:HD13	1.85	0.58
1:E:58:PRO:HG2	1:F:59:LYS:HG3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:144:VAL:HG23	1:F:148:LEU:HD11	1.84	0.58
1:I:259:LYS:CG	3:Y:1:NAG:H83	2.33	0.58
1:H:48:THR:HG23	1:H:50:SER:H	1.68	0.58
1:H:61:PRO:HG2	1:I:62:LYS:HG3	1.86	0.58
1:B:36:ASP:O	1:B:38:GLY:N	2.36	0.58
1:E:66:GLY:HA3	1:F:67:LEU:O	2.04	0.58
1:F:295:GLN:OE1	2:U:1:NAG:H82	2.04	0.58
1:A:41:LYS:HE3	1:A:49:PHE:CE2	2.39	0.58
1:D:341:THR:HB	1:E:335:THR:HG23	1.86	0.58
1:E:300:ASN:OD1	1:E:301:SER:N	2.37	0.58
1:G:26:LEU:HD11	1:G:49:PHE:CE1	2.39	0.58
1:A:48:THR:HG23	1:A:50:SER:H	1.69	0.57
1:A:290:LYS:HD2	1:B:268:GLU:HA	1.86	0.57
1:E:357:VAL:HG23	1:E:358:ASP:H	1.69	0.57
1:G:292:THR:HG22	1:H:264:ASN:ND2	2.17	0.57
1:E:336:VAL:HA	1:F:330:SER:HB3	1.86	0.57
1:D:42:PHE:CE1	1:D:55:LEU:HD13	2.40	0.57
1:B:259:LYS:HG2	3:M:1:NAG:H82	1.85	0.57
1:E:345:PRO:HB3	1:E:351:ALA:O	2.05	0.57
1:B:340:LEU:HD11	1:C:336:VAL:HG23	1.85	0.57
1:E:61:PRO:HG2	1:F:62:LYS:HG2	1.85	0.57
1:A:249:ARG:NH2	3:K:1:NAG:O3	2.38	0.57
1:E:348:GLY:O	1:F:355:ASP:N	2.36	0.57
1:I:13:ARG:HB3	1:I:33:PHE:HB3	1.87	0.57
1:I:300:ASN:OD1	1:I:301:SER:N	2.38	0.57
1:F:266:ASN:OD1	1:F:267:VAL:N	2.38	0.57
1:E:263:GLU:OE1	1:F:249:ARG:NH1	2.38	0.57
1:A:88:TYR:HA	1:A:91:LEU:HD12	1.86	0.56
1:D:30:GLU:OE1	1:D:31:ILE:N	2.38	0.56
1:G:14:MET:HB2	1:G:19:TRP:HD1	1.69	0.56
1:G:364:LEU:HD22	1:H:361:ILE:CG2	2.24	0.56
1:A:119:ASP:OD1	1:C:117:LYS:NZ	2.33	0.56
1:D:69:GLY:HA3	1:E:67:LEU:O	2.05	0.56
1:E:266:ASN:OD1	1:E:267:VAL:N	2.38	0.56
1:A:61:PRO:HG2	1:B:59:LYS:HG3	1.88	0.56
1:B:64:ASP:HB2	1:C:65:THR:HG22	1.87	0.56
1:C:266:ASN:OD1	1:C:267:VAL:N	2.38	0.56
1:D:336:VAL:HG13	1:E:330:SER:HB3	1.88	0.56
1:E:182:TYR:CD2	1:F:162:SER:HA	2.40	0.56
1:H:165:GLY:O	1:I:147:GLN:NE2	2.38	0.56
1:H:216:ASN:ND2	3:X:3:NAG:O7	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:MET:HB2	1:B:19:TRP:CD1	2.41	0.56
1:H:315:LYS:HE3	1:I:304:GLY:HA3	1.88	0.56
1:D:216:ASN:HD22	1:F:232:SER:H	1.52	0.56
1:E:14:MET:HB2	1:E:19:TRP:CD1	2.40	0.56
1:H:313:ARG:NH2	2:V:2:BDP:O3	2.39	0.56
1:I:25:ILE:HG13	1:I:25:ILE:O	2.04	0.56
1:A:59:LYS:HG2	1:C:58:PRO:HG2	1.86	0.56
1:E:41:LYS:HE3	1:E:49:PHE:CE2	2.40	0.56
1:H:57:GLY:HA2	1:I:56:THR:HB	1.87	0.56
1:H:45:GLY:O	1:H:46:GLN:HB2	2.05	0.56
1:C:255:LEU:HA	2:O:3:NAG:O4	2.06	0.55
1:D:183:THR:HG22	1:D:185:LYS:H	1.71	0.55
1:D:335:THR:OG1	1:E:329:HIS:ND1	2.31	0.55
1:I:41:LYS:HE3	1:I:49:PHE:CE2	2.41	0.55
1:I:34:GLU:OE1	1:I:37:THR:OG1	2.16	0.55
1:I:41:LYS:HD2	1:I:52:LEU:HB3	1.87	0.55
1:A:60:GLY:HA3	1:B:58:PRO:O	2.07	0.55
1:D:17:ASP:OD1	1:D:21:ARG:NH1	2.39	0.55
1:E:196:ARG:NH2	1:E:198:ASP:OD2	2.38	0.55
1:G:56:THR:HB	1:I:57:GLY:HA2	1.87	0.55
1:G:266:ASN:OD1	1:G:267:VAL:N	2.40	0.55
1:H:159:PRO:O	3:W:5:NAG:H81	2.06	0.55
1:D:340:LEU:HD11	1:E:336:VAL:HG23	1.88	0.55
1:E:237:THR:OG1	3:P:2:BDP:O3	2.20	0.55
1:G:48:THR:HG23	1:G:50:SER:H	1.70	0.55
1:G:57:GLY:HA3	1:H:55:LEU:O	2.06	0.55
1:A:348:GLY:HA2	1:B:354:LYS:HD2	1.89	0.55
1:C:311:ARG:NH2	1:C:318:ASP:OD2	2.40	0.55
1:E:342:VAL:HG23	1:E:353:THR:HG22	1.88	0.55
1:H:266:ASN:OD1	1:H:267:VAL:N	2.40	0.55
3:Y:7:NAG:O4	3:Y:8:GCD:O5	2.22	0.55
1:D:161:SER:HB3	3:P:5:NAG:H61	1.88	0.54
1:E:36:ASP:O	1:E:38:GLY:N	2.40	0.54
1:E:237:THR:HG21	3:P:3:NAG:H62	1.89	0.54
1:F:170:ASP:OD1	1:F:172:SER:OG	2.21	0.54
1:F:300:ASN:OD1	1:F:301:SER:N	2.39	0.54
1:G:300:ASN:OD1	1:G:301:SER:N	2.36	0.54
1:A:42:PHE:HB3	1:B:33:PHE:CE2	2.43	0.54
1:I:183:THR:HG21	1:I:192:LEU:HD11	1.88	0.54
1:F:347:SER:OG	1:F:348:GLY:N	2.40	0.54
1:F:360:LYS:C	1:F:362:ALA:H	2.15	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:219:ASN:ND2	3:Y:3:NAG:O7	2.40	0.54
1:F:96:ASP:OD1	1:F:96:ASP:N	2.35	0.54
1:G:56:THR:O	1:H:39:PHE:HA	2.08	0.54
1:F:13:ARG:HB3	1:F:33:PHE:HB3	1.88	0.54
1:H:169:ILE:HA	1:I:150:PHE:HB2	1.89	0.54
1:F:19:TRP:NE1	1:F:33:PHE:O	2.38	0.54
1:B:354:LYS:HE2	1:B:358:ASP:OD2	2.08	0.53
1:D:230:PHE:HE2	3:R:1:NAG:H5	1.72	0.53
1:I:48:THR:HG23	1:I:50:SER:H	1.73	0.53
1:B:240:ASN:ND2	1:C:229:ASN:O	2.31	0.53
1:B:101:ALA:HB2	1:C:103:LYS:HA	1.90	0.53
1:G:292:THR:CG2	1:H:264:ASN:HB2	2.38	0.53
1:D:19:TRP:NE1	1:D:33:PHE:O	2.42	0.53
1:H:196:ARG:NH2	1:H:198:ASP:OD2	2.41	0.53
1:D:196:ARG:HH12	3:R:6:BDP:C6	2.22	0.52
1:C:182:TYR:HE1	1:C:184:ASN:HB2	1.73	0.52
1:H:292:THR:OG1	1:I:272:ASP:OD2	2.21	0.52
1:H:15:SER:HA	1:H:35:THR:HG22	1.91	0.52
1:D:131:ILE:O	1:D:135:LYS:HG3	2.09	0.52
1:E:34:GLU:OE1	1:E:37:THR:OG1	2.25	0.52
1:A:335:THR:HG23	1:C:341:THR:HB	1.92	0.52
2:O:2:BDP:O4	2:O:3:NAG:O6	2.19	0.52
1:D:300:ASN:OD1	1:D:301:SER:N	2.42	0.52
1:E:177:ALA:O	1:F:165:GLY:HA2	2.10	0.52
1:H:243:GLY:HA2	1:I:230:PHE:HD1	1.75	0.52
1:F:223:ARG:NH2	3:P:4:BDP:O2	2.40	0.51
1:D:295:GLN:NE2	2:T:2:BDP:O6A	2.43	0.51
1:H:301:SER:HB2	1:I:282:ILE:HB	1.92	0.51
1:A:13:ARG:HB3	1:A:33:PHE:HB3	1.91	0.51
1:C:313:ARG:NE	1:C:318:ASP:OD1	2.43	0.51
1:G:162:SER:HA	1:I:182:TYR:HD2	1.76	0.51
1:D:313:ARG:NE	1:D:318:ASP:OD1	2.44	0.51
1:F:41:LYS:HE2	1:F:54:TYR:CZ	2.45	0.51
1:F:208:PHE:CE2	3:R:4:BDP:H4	2.45	0.51
2:U:1:NAG:H1	2:U:1:NAG:O7	2.09	0.51
1:A:199:LYS:HG2	1:B:184:ASN:OD1	2.11	0.51
1:D:117:LYS:O	1:F:117:LYS:NZ	2.39	0.51
1:A:349:LYS:HB3	1:B:341:THR:HG23	1.92	0.51
1:C:41:LYS:HE3	1:C:49:PHE:CE2	2.46	0.51
1:D:320:PHE:HD2	1:D:330:SER:HA	1.75	0.51
1:H:41:LYS:HE3	1:H:49:PHE:CE2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:GLN:HB3	3:K:1:NAG:H83	1.93	0.51
1:C:11:PHE:O	1:C:13:ARG:NH1	2.41	0.50
1:G:168:ASN:O	1:H:150:PHE:N	2.39	0.50
1:G:230:PHE:CZ	3:X:1:NAG:H1	2.45	0.50
1:A:240:ASN:ND2	1:B:229:ASN:O	2.31	0.50
1:A:311:ARG:HE	1:A:313:ARG:HE	1.59	0.50
1:E:165:GLY:O	1:F:147:GLN:NE2	2.44	0.50
1:G:119:ASP:HA	1:I:117:LYS:HD3	1.93	0.50
1:B:240:ASN:ND2	1:C:231:SER:OG	2.45	0.50
1:D:42:PHE:HZ	1:D:55:LEU:HD13	1.72	0.50
1:G:157:ILE:HB	1:I:175:GLU:O	2.12	0.50
1:G:230:PHE:HZ	3:X:1:NAG:H1	1.77	0.50
3:P:7:NAG:O4	3:P:8:GCD:O5	2.26	0.50
1:A:277:ALA:O	1:B:257:THR:HB	2.12	0.50
1:E:319:LYS:HD3	1:F:324:PRO:O	2.12	0.50
1:E:357:VAL:HG23	1:E:358:ASP:OD1	2.12	0.50
1:C:44:ASP:OD2	1:C:47:ASN:ND2	2.45	0.50
1:D:67:LEU:O	1:F:66:GLY:HA3	2.12	0.50
1:D:295:GLN:N	1:E:275:ALA:O	2.33	0.50
1:A:45:GLY:O	1:A:46:GLN:HB2	2.11	0.50
1:D:216:ASN:ND2	1:F:232:SER:H	2.10	0.50
1:G:67:LEU:HB2	1:H:65:THR:HG22	1.94	0.50
1:G:194:ILE:HB	1:H:180:VAL:HG22	1.93	0.50
1:H:14:MET:HB2	1:H:19:TRP:CD1	2.47	0.50
1:H:335:THR:HB	1:I:329:HIS:ND1	2.27	0.50
1:A:311:ARG:CZ	1:A:313:ARG:HH21	2.24	0.49
1:A:67:LEU:O	1:C:66:GLY:HA3	2.11	0.49
1:C:298:TYR:CD1	2:N:3:NAG:H83	2.47	0.49
1:E:252:GLU:OE1	1:E:257:THR:OG1	2.23	0.49
1:G:36:ASP:C	1:G:38:GLY:H	2.20	0.49
1:G:340:LEU:HD11	1:H:336:VAL:HG23	1.93	0.49
1:G:361:ILE:O	1:G:365:LYS:N	2.45	0.49
1:A:289:GLY:HA2	1:B:267:VAL:HA	1.94	0.49
1:H:340:LEU:HD13	1:I:334:SER:HB2	1.94	0.49
1:D:41:LYS:HE3	1:D:49:PHE:CE2	2.48	0.49
1:D:44:ASP:O	1:D:46:GLN:HG2	2.12	0.49
1:E:240:ASN:ND2	1:F:229:ASN:O	2.40	0.49
1:G:284:LYS:HA	1:G:292:THR:HG21	1.93	0.49
1:H:64:ASP:HB2	1:I:65:THR:HG22	1.94	0.49
1:C:313:ARG:HG2	1:C:318:ASP:HA	1.93	0.49
1:D:41:LYS:HE2	1:D:54:TYR:CE1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:36:ASP:C	1:C:38:GLY:H	2.19	0.49
1:D:36:ASP:C	1:D:38:GLY:H	2.20	0.49
1:E:246:MET:HA	1:F:234:LEU:O	2.13	0.49
1:G:117:LYS:HD2	1:H:117:LYS:HB2	1.95	0.48
2:O:3:NAG:O7	2:O:3:NAG:H3	2.12	0.48
1:D:276:ALA:O	1:E:256:GLY:HA2	2.13	0.48
1:D:13:ARG:HB3	1:D:33:PHE:HB3	1.95	0.48
1:G:259:LYS:CG	3:W:1:NAG:H82	2.43	0.48
1:D:220:ILE:HG22	1:D:233:ALA:HB1	1.95	0.48
1:E:124:TYR:CE2	1:F:126:LYS:HG3	2.49	0.48
1:G:119:ASP:OD1	1:G:119:ASP:N	2.45	0.48
1:B:230:PHE:HE2	3:K:1:NAG:H1	1.79	0.48
1:D:330:SER:OG	1:D:331:GLY:N	2.43	0.48
1:F:356:TYR:OH	1:F:361:ILE:HG21	2.14	0.48
1:C:215:THR:OG1	1:C:216:ASN:N	2.47	0.48
1:E:82:LYS:O	1:F:84:GLY:N	2.42	0.48
1:E:275:ALA:HB2	1:F:253:LYS:O	2.13	0.48
1:I:330:SER:OG	1:I:331:GLY:N	2.47	0.48
1:A:266:ASN:OD1	1:A:267:VAL:N	2.46	0.48
1:G:362:ALA:HA	1:G:365:LYS:HB3	1.96	0.48
3:P:4:BDP:O3	3:P:5:NAG:O7	2.28	0.48
1:A:57:GLY:HA3	1:B:55:LEU:O	2.13	0.48
1:F:36:ASP:C	1:F:38:GLY:H	2.22	0.48
1:I:365:LYS:O	1:I:367:LEU:N	2.47	0.48
1:B:208:PHE:HB3	1:C:194:ILE:HG12	1.96	0.48
1:E:117:LYS:HD2	1:F:117:LYS:HB2	1.95	0.48
1:H:88:TYR:HA	1:H:91:LEU:HD12	1.96	0.48
1:I:360:LYS:HD3	1:I:360:LYS:HA	1.66	0.48
1:B:222:MET:HE1	1:B:231:SER:O	2.14	0.47
1:B:359:GLU:O	1:B:363:GLU:HB2	2.15	0.47
1:H:168:ASN:O	1:I:150:PHE:N	2.38	0.47
1:H:215:THR:OG1	1:H:216:ASN:N	2.47	0.47
1:A:42:PHE:CE2	1:A:53:LYS:HB2	2.50	0.47
1:A:65:THR:HG22	1:C:64:ASP:HB2	1.96	0.47
1:D:275:ALA:HB2	1:E:253:LYS:O	2.14	0.47
1:D:315:LYS:NZ	1:E:304:GLY:HA3	2.29	0.47
1:B:330:SER:OG	1:B:331:GLY:N	2.47	0.47
1:F:354:LYS:HA	1:F:357:VAL:HG22	1.96	0.47
2:V:1:NAG:O7	2:V:1:NAG:H3	2.14	0.47
1:A:177:ALA:O	1:B:165:GLY:HA2	2.14	0.47
1:B:215:THR:OG1	1:B:216:ASN:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:59:LYS:HG3	1:F:58:PRO:HG2	1.96	0.47
1:E:43:GLY:HA2	1:E:52:LEU:HD22	1.96	0.47
1:G:80:ALA:HB2	1:I:79:PRO:HG2	1.97	0.47
1:G:237:THR:HG23	1:H:221:VAL:HG22	1.96	0.47
1:C:87:ASP:OD2	1:C:90:GLN:NE2	2.46	0.47
1:H:347:SER:HB3	1:H:350:HIS:CD2	2.50	0.47
1:H:347:SER:OG	1:H:348:GLY:N	2.47	0.47
1:F:330:SER:OG	1:F:331:GLY:N	2.48	0.47
1:G:26:LEU:HD11	1:G:49:PHE:HE1	1.78	0.47
1:H:101:ALA:HB2	1:I:103:LYS:HA	1.97	0.47
1:H:243:GLY:HA2	1:I:230:PHE:CD1	2.50	0.47
1:H:356:TYR:O	1:H:360:LYS:HG2	2.14	0.47
1:A:196:ARG:HH22	3:M:7:NAG:H83	1.79	0.47
1:B:169:ILE:HA	1:C:150:PHE:HB2	1.96	0.47
1:B:259:LYS:CG	3:M:1:NAG:C8	2.86	0.47
1:D:143:ILE:HG12	1:F:149:GLN:HB2	1.95	0.47
1:D:251:VAL:HA	1:F:263:GLU:O	2.13	0.47
1:E:41:LYS:HE2	1:E:54:TYR:CE2	2.49	0.47
1:H:138:SER:O	1:H:140:THR:N	2.48	0.47
1:I:208:PHE:HE2	3:X:4:BDP:H2	1.80	0.47
1:D:261:THR:HA	1:F:281:ASP:O	2.15	0.47
1:D:334:SER:HB2	1:E:328:PHE:O	2.15	0.47
1:E:28:GLU:HA	1:E:43:GLY:O	2.15	0.47
1:H:330:SER:OG	1:H:331:GLY:N	2.48	0.47
1:A:329:HIS:CD2	1:C:335:THR:HB	2.49	0.47
1:B:41:LYS:HE2	1:B:54:TYR:CE1	2.49	0.47
1:G:267:VAL:HA	1:I:289:GLY:HA2	1.95	0.47
1:G:283:VAL:O	1:G:292:THR:HG21	2.14	0.47
1:I:215:THR:OG1	1:I:216:ASN:N	2.48	0.47
1:B:268:GLU:HG3	1:B:271:TYR:HB2	1.97	0.47
1:G:352:ALA:N	1:H:352:ALA:O	2.48	0.47
1:I:347:SER:HB3	1:I:350:HIS:CD2	2.49	0.47
2:U:2:BDP:O4	2:U:3:NAG:O7	2.32	0.47
1:D:122:ALA:O	1:E:126:LYS:N	2.48	0.46
1:D:266:ASN:OD1	1:D:267:VAL:N	2.48	0.46
1:E:254:ALA:HB1	2:U:2:BDP:H3	1.98	0.46
1:G:274:ASN:HA	2:V:2:BDP:O2	2.15	0.46
1:D:45:GLY:O	1:D:46:GLN:HB3	2.14	0.46
1:F:41:LYS:HE2	1:F:54:TYR:CE1	2.50	0.46
3:Y:5:NAG:H2	3:Y:5:NAG:H61	1.97	0.46
1:A:182:TYR:CD2	1:B:162:SER:HA	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:GLU:O	1:C:251:VAL:HA	2.15	0.46
1:F:41:LYS:HD2	1:F:52:LEU:CD1	2.40	0.46
1:F:88:TYR:HA	1:F:91:LEU:HD12	1.97	0.46
1:H:247:GLN:HB3	3:W:1:NAG:H83	1.98	0.46
1:A:251:VAL:HA	1:C:263:GLU:O	2.14	0.46
1:C:295:GLN:HE22	2:N:3:NAG:C7	2.28	0.46
1:E:169:ILE:HG21	1:E:178:ALA:HB1	1.98	0.46
1:G:162:SER:HA	1:I:182:TYR:CD2	2.51	0.46
1:G:330:SER:OG	1:G:331:GLY:N	2.48	0.46
1:I:311:ARG:NH2	1:I:318:ASP:OD2	2.46	0.46
3:S:4:BDP:O4	3:S:5:NAG:O7	2.34	0.46
1:C:13:ARG:HB3	1:C:33:PHE:HB3	1.96	0.46
1:D:62:LYS:HG2	1:F:63:GLY:HA2	1.98	0.46
1:F:70:LYS:HB3	1:F:70:LYS:HE2	1.72	0.46
1:G:264:ASN:ND2	1:I:291:GLY:O	2.47	0.46
1:F:53:LYS:HB3	1:F:53:LYS:HE3	1.70	0.46
1:H:368:ILE:HD11	1:I:367:LEU:HD11	1.98	0.46
1:A:154:LYS:HD3	1:D:270:LYS:NZ	2.30	0.46
1:C:182:TYR:CE1	1:C:184:ASN:HB2	2.51	0.46
1:I:208:PHE:CE2	3:X:4:BDP:H2	2.51	0.46
3:R:2:BDP:O4	3:R:2:BDP:O2	2.30	0.46
1:A:359:GLU:O	1:A:363:GLU:HB2	2.15	0.46
1:C:16:ALA:HB2	1:C:36:ASP:HB2	1.97	0.46
1:D:224:GLN:HB2	1:F:239:ALA:HB1	1.98	0.46
1:F:45:GLY:O	1:F:46:GLN:HB2	2.16	0.46
1:G:216:ASN:HA	1:I:232:SER:O	2.15	0.46
1:G:251:VAL:HA	1:I:263:GLU:O	2.16	0.46
1:A:196:ARG:HE	1:A:198:ASP:CG	2.24	0.46
1:D:215:THR:OG1	1:D:216:ASN:N	2.49	0.46
1:F:214:LYS:HE2	1:F:214:LYS:HB3	1.69	0.46
1:E:311:ARG:HH11	1:E:313:ARG:HH21	1.64	0.45
3:P:6:BDP:C6	3:P:7:NAG:H2	2.46	0.45
2:U:2:BDP:H5	2:U:3:NAG:O5	2.16	0.45
1:A:215:THR:OG1	1:A:216:ASN:N	2.48	0.45
1:C:359:GLU:C	1:C:361:ILE:H	2.24	0.45
1:E:343:LYS:N	1:F:339:ASN:OD1	2.45	0.45
1:A:58:PRO:O	1:A:59:LYS:HG3	2.17	0.45
1:G:87:ASP:HA	1:I:92:GLN:HB2	1.99	0.45
1:A:144:VAL:HG23	1:C:148:LEU:HD11	1.98	0.45
1:D:336:VAL:HB	1:F:341:THR:O	2.16	0.45
1:E:305:THR:HA	1:F:292:THR:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:215:THR:OG1	1:F:216:ASN:N	2.50	0.45
1:F:350:HIS:H	1:F:350:HIS:CD2	2.33	0.45
1:I:41:LYS:HE2	1:I:54:TYR:CE1	2.52	0.45
1:C:330:SER:OG	1:C:331:GLY:N	2.50	0.45
1:G:366:LYS:O	1:G:368:ILE:N	2.48	0.45
1:A:354:LYS:HD3	1:C:348:GLY:HA2	1.99	0.45
1:C:345:PRO:HG3	1:C:352:ALA:HA	1.97	0.45
1:D:181:MET:HE1	1:D:193:MET:SD	2.56	0.45
1:B:171:MET:HE1	1:C:157:ILE:HG21	1.97	0.45
1:C:44:ASP:HB3	1:C:47:ASN:HB2	1.99	0.45
1:D:117:LYS:C	1:F:117:LYS:HZ2	2.23	0.45
1:G:291:GLY:C	1:G:292:THR:HG23	2.42	0.45
1:H:66:GLY:HA3	1:I:67:LEU:O	2.17	0.45
1:A:168:ASN:O	1:B:150:PHE:N	2.38	0.45
1:A:313:ARG:NE	1:A:318:ASP:OD1	2.50	0.45
1:G:313:ARG:NE	1:G:318:ASP:OD1	2.49	0.45
1:F:220:ILE:HG22	1:F:233:ALA:HB1	1.99	0.45
1:G:53:LYS:NZ	1:H:37:THR:O	2.50	0.45
1:E:330:SER:OG	1:E:331:GLY:N	2.49	0.44
1:G:18:GLU:O	1:G:22:SER:OG	2.29	0.44
1:I:254:ALA:O	1:I:255:LEU:HG	2.18	0.44
1:B:179:MET:HE3	1:B:179:MET:HB2	1.82	0.44
1:E:320:PHE:HA	1:E:329:HIS:O	2.17	0.44
1:A:26:LEU:O	1:A:45:GLY:HA2	2.17	0.44
1:E:182:TYR:HD2	1:F:162:SER:HA	1.81	0.44
1:E:251:VAL:HG23	1:F:241:GLU:HA	1.98	0.44
1:G:300:ASN:HA	1:I:313:ARG:O	2.17	0.44
1:A:67:LEU:HB2	1:B:65:THR:HG23	1.99	0.44
1:A:159:PRO:HB3	1:C:176:GLY:HA2	1.98	0.44
1:A:312:ILE:HB	1:A:320:PHE:HB3	2.00	0.44
1:G:297:ILE:HB	1:I:310:LEU:HD13	1.99	0.44
1:A:222:MET:HE1	1:A:231:SER:O	2.16	0.44
1:B:300:ASN:O	1:C:281:ASP:HA	2.17	0.44
1:C:94:LYS:HB3	1:C:94:LYS:HE2	1.72	0.44
1:E:342:VAL:HB	1:F:339:ASN:OD1	2.18	0.44
1:G:344:ASP:OD1	1:I:349:LYS:NZ	2.50	0.44
1:H:139:LEU:HA	1:H:139:LEU:HD12	1.74	0.44
1:B:41:LYS:HE2	1:B:54:TYR:CD1	2.52	0.44
1:E:263:GLU:HB2	1:F:249:ARG:HH11	1.83	0.44
1:A:360:LYS:HD3	1:A:360:LYS:HA	1.77	0.44
1:B:306:ALA:O	1:B:324:PRO:HG3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:261:THR:HB	3:Y:1:NAG:C8	2.46	0.44
1:A:193:MET:HE1	1:C:193:MET:HE1	1.99	0.44
1:A:259:LYS:HE2	1:A:261:THR:OG1	2.18	0.44
1:C:19:TRP:HB3	1:C:49:PHE:CD2	2.53	0.44
1:F:181:MET:HB3	1:F:192:LEU:HD23	1.99	0.44
1:H:360:LYS:HA	1:H:360:LYS:HD3	1.75	0.44
1:I:191:PRO:HD2	1:I:194:ILE:HD11	2.00	0.44
1:A:199:LYS:HE2	1:B:184:ASN:OD1	2.18	0.43
1:D:230:PHE:CE2	3:R:1:NAG:H5	2.51	0.43
1:G:313:ARG:HG2	1:G:318:ASP:HA	1.99	0.43
3:R:7:NAG:O4	3:R:8:GCD:O5	2.30	0.43
1:B:41:LYS:HG3	1:B:49:PHE:CZ	2.54	0.43
1:B:131:ILE:O	1:B:135:LYS:HG3	2.17	0.43
1:C:13:ARG:HB2	1:C:35:THR:HG23	2.00	0.43
1:A:25:ILE:HD13	1:A:46:GLN:HA	2.00	0.43
1:B:45:GLY:O	1:B:46:GLN:HB2	2.17	0.43
1:C:311:ARG:HE	1:C:313:ARG:HE	1.66	0.43
1:E:19:TRP:HB3	1:E:49:PHE:CD2	2.53	0.43
1:E:162:SER:OG	3:S:6:BDP:O6A	2.26	0.43
1:C:14:MET:HB2	1:C:19:TRP:CD1	2.54	0.43
1:A:315:LYS:NZ	1:B:304:GLY:HA3	2.34	0.43
1:B:136:LYS:NZ	1:C:133:LEU:O	2.52	0.43
1:E:13:ARG:HA	1:E:33:PHE:HB2	2.00	0.43
1:I:53:LYS:HB3	1:I:53:LYS:HE3	1.63	0.43
3:Y:3:NAG:H3	3:Y:4:BDP:H2	1.50	0.43
1:A:58:PRO:HA	1:B:39:PHE:CZ	2.53	0.43
1:B:160:SER:OG	1:B:162:SER:OG	2.31	0.43
1:D:335:THR:O	1:E:329:HIS:HA	2.19	0.43
1:D:346:THR:O	1:D:346:THR:OG1	2.37	0.43
1:E:116:SER:O	1:F:120:LYS:N	2.37	0.43
1:E:349:LYS:HD2	1:E:349:LYS:N	2.34	0.43
1:H:161:SER:H	3:W:6:BDP:HB	1.66	0.43
1:E:124:TYR:CZ	1:F:126:LYS:HG3	2.54	0.43
1:E:181:MET:HB3	1:E:192:LEU:HD23	2.01	0.43
1:F:311:ARG:HH21	2:T:1:NAG:H82	1.83	0.43
3:K:7:NAG:O4	3:K:8:GCD:O6A	2.37	0.43
1:A:94:LYS:HB3	1:A:94:LYS:HE2	1.71	0.43
1:D:22:SER:OG	1:D:24:VAL:HG12	2.19	0.43
1:E:41:LYS:HG3	1:E:49:PHE:CZ	2.53	0.43
1:E:215:THR:OG1	1:E:216:ASN:N	2.48	0.43
1:G:366:LYS:HE2	1:G:366:LYS:HB3	1.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:249:ARG:NE	3:W:2:BDP:O2	2.48	0.43
3:S:1:NAG:O7	3:S:1:NAG:O3	2.37	0.43
1:A:233:ALA:O	1:B:217:ALA:HB3	2.19	0.43
1:D:62:LYS:HG3	1:F:61:PRO:HG2	2.00	0.43
1:G:28:GLU:HA	1:G:44:ASP:HA	2.00	0.43
1:G:354:LYS:HE3	1:I:356:TYR:OH	2.18	0.43
1:F:345:PRO:HD3	1:F:352:ALA:HB2	2.01	0.42
1:G:144:VAL:HG23	1:I:148:LEU:HD11	2.00	0.42
1:G:215:THR:OG1	1:G:216:ASN:N	2.52	0.42
1:I:36:ASP:OD1	1:I:37:THR:N	2.45	0.42
1:I:312:ILE:HD12	1:I:320:PHE:HB3	2.00	0.42
1:C:317:GLU:HG3	1:C:319:LYS:HE2	2.00	0.42
1:D:181:MET:HB3	1:D:192:LEU:HD13	2.00	0.42
1:E:179:MET:SD	1:F:179:MET:HE1	2.59	0.42
1:G:69:GLY:HA3	1:H:67:LEU:O	2.20	0.42
1:B:123:VAL:N	1:C:126:LYS:HG3	2.34	0.42
1:D:217:ALA:HB3	1:F:233:ALA:O	2.19	0.42
1:H:341:THR:O	1:I:336:VAL:N	2.51	0.42
1:B:352:ALA:N	1:C:352:ALA:O	2.52	0.42
1:C:36:ASP:C	1:C:38:GLY:N	2.77	0.42
1:G:311:ARG:NE	1:G:313:ARG:HH21	2.18	0.42
1:G:312:ILE:N	1:G:320:PHE:O	2.47	0.42
1:G:314:ASN:OD1	1:H:301:SER:HB3	2.19	0.42
1:A:76:PRO:HD2	1:B:74:THR:HA	2.01	0.42
1:D:80:ALA:HB2	1:F:79:PRO:HG2	2.02	0.42
1:H:188:THR:OG1	1:H:189:ASP:N	2.53	0.42
1:B:41:LYS:HD3	1:B:53:LYS:C	2.44	0.42
1:E:306:ALA:H	1:F:293:ALA:HA	1.85	0.42
1:F:298:TYR:CD1	1:F:298:TYR:C	2.97	0.42
1:G:28:GLU:OE2	1:H:13:ARG:CD	2.67	0.42
1:G:68:GLN:HG3	1:I:67:LEU:HD12	2.02	0.42
1:I:131:ILE:O	1:I:135:LYS:HG3	2.19	0.42
1:D:85:THR:HB	1:F:92:GLN:OE1	2.19	0.42
1:A:349:LYS:HD3	1:A:349:LYS:HA	1.63	0.42
1:B:342:VAL:HG23	1:B:353:THR:HG22	2.01	0.42
1:C:313:ARG:NH1	2:O:2:BDP:O6A	2.47	0.42
1:G:291:GLY:O	1:G:292:THR:CG2	2.68	0.42
1:G:301:SER:O	1:G:304:GLY:N	2.52	0.42
3:K:7:NAG:O7	3:K:7:NAG:O3	2.37	0.42
1:A:330:SER:OG	1:A:331:GLY:N	2.49	0.42
1:G:259:LYS:HG2	3:W:1:NAG:H82	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:171:MET:HE3	1:I:171:MET:HB3	1.95	0.42
3:W:1:NAG:O4	3:W:2:BDP:O5	2.31	0.42
1:A:172:SER:HB3	1:C:184:ASN:HB3	2.01	0.42
1:G:192:LEU:HB2	1:I:206:ALA:O	2.20	0.42
1:H:337:ALA:HB3	1:I:331:GLY:O	2.20	0.42
3:X:5:NAG:H3	3:X:6:BDP:H2	1.45	0.42
1:C:282:ILE:HG22	1:C:292:THR:HG21	2.02	0.41
1:D:151:LYS:HG3	1:E:143:ILE:HG23	2.02	0.41
1:D:232:SER:O	1:E:216:ASN:HA	2.20	0.41
1:E:230:PHE:CZ	3:P:1:NAG:H1	2.54	0.41
1:E:254:ALA:O	1:E:255:LEU:HG	2.20	0.41
1:F:340:LEU:O	1:F:342:VAL:HG13	2.20	0.41
1:G:359:GLU:HG2	1:G:360:LYS:HE3	2.02	0.41
1:I:313:ARG:NH2	2:Z:2:BDP:C1	2.83	0.41
1:B:36:ASP:C	1:B:38:GLY:H	2.28	0.41
1:D:58:PRO:HG2	1:E:56:THR:HG22	2.01	0.41
1:D:182:TYR:CE2	1:D:184:ASN:HB2	2.55	0.41
1:E:235:ASN:ND2	1:F:219:ASN:OD1	2.52	0.41
1:A:179:MET:HE1	1:C:179:MET:SD	2.60	0.41
1:A:180:VAL:HG22	1:C:194:ILE:HB	2.00	0.41
1:B:232:SER:O	1:C:216:ASN:HA	2.20	0.41
1:D:14:MET:H	1:D:19:TRP:HE1	1.67	0.41
1:D:308:LYS:HD3	1:D:321:TYR:CD1	2.55	0.41
1:E:16:ALA:HB2	1:E:36:ASP:HB2	2.02	0.41
3:S:5:NAG:H61	3:S:5:NAG:H2	2.01	0.41
1:A:243:GLY:HA2	1:B:230:PHE:HD1	1.85	0.41
1:A:364:LEU:HD22	1:B:364:LEU:HD11	2.01	0.41
1:C:361:ILE:HD13	1:C:361:ILE:HA	1.94	0.41
1:D:28:GLU:O	1:D:28:GLU:HG3	2.21	0.41
1:D:291:GLY:O	1:E:264:ASN:ND2	2.49	0.41
1:E:230:PHE:CE2	3:P:1:NAG:H1	2.55	0.41
1:F:349:LYS:HA	1:F:349:LYS:HD3	1.68	0.41
1:B:349:LYS:HD3	1:B:349:LYS:HA	1.76	0.41
1:D:55:LEU:O	1:D:55:LEU:HG	2.21	0.41
1:D:194:ILE:HG21	1:F:161:SER:HB2	2.03	0.41
1:E:168:ASN:O	1:F:150:PHE:N	2.44	0.41
1:F:125:SER:N	1:F:128:GLU:HB2	2.35	0.41
1:G:226:SER:OG	1:I:200:ASP:OD2	2.29	0.41
1:D:183:THR:OG1	1:D:192:LEU:HD21	2.21	0.41
1:D:265:PRO:O	1:F:285:LYS:NZ	2.46	0.41
1:A:41:LYS:HD2	1:A:52:LEU:CD1	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:PRO:HG2	1:A:194:ILE:HD11	2.01	0.41
1:E:223:ARG:HH21	1:F:210:ASP:CG	2.29	0.41
1:G:210:ASP:CG	1:I:223:ARG:HH11	2.29	0.41
1:A:42:PHE:HB3	1:B:33:PHE:HE2	1.84	0.41
1:B:41:LYS:HD3	1:B:53:LYS:O	2.21	0.41
1:B:364:LEU:HD12	1:B:364:LEU:O	2.21	0.41
1:D:122:ALA:C	1:E:126:LYS:HB2	2.46	0.41
1:D:192:LEU:HD23	1:D:192:LEU:HA	1.82	0.41
1:F:361:ILE:HD12	1:F:361:ILE:HA	1.86	0.41
1:G:252:GLU:O	1:I:262:HIS:NE2	2.43	0.41
1:A:223:ARG:HE	1:B:210:ASP:CG	2.29	0.41
1:C:51:LYS:HB3	1:C:51:LYS:HE2	1.79	0.41
1:D:179:MET:HE1	1:F:179:MET:SD	2.60	0.41
1:F:88:TYR:O	1:F:94:LYS:HD2	2.19	0.41
1:F:345:PRO:HD3	1:F:352:ALA:CB	2.51	0.41
1:G:28:GLU:OE2	1:H:13:ARG:HD3	2.21	0.41
1:I:182:TYR:HE1	1:I:184:ASN:HB2	1.85	0.41
1:A:230:PHE:CE2	3:M:1:NAG:H5	2.56	0.41
1:B:251:VAL:O	1:C:241:GLU:HA	2.21	0.40
1:D:289:GLY:HA2	1:E:267:VAL:HA	2.02	0.40
1:E:357:VAL:O	1:E:361:ILE:HB	2.21	0.40
1:E:362:ALA:C	1:E:364:LEU:H	2.29	0.40
1:G:137:LEU:HD23	1:I:137:LEU:HD23	2.03	0.40
1:H:233:ALA:O	1:I:217:ALA:HB3	2.21	0.40
1:A:273:GLU:HA	1:C:293:ALA:O	2.21	0.40
1:D:233:ALA:O	1:E:217:ALA:HB3	2.21	0.40
1:E:179:MET:HA	1:F:167:ILE:O	2.21	0.40
1:G:87:ASP:OD1	1:G:90:GLN:HB2	2.21	0.40
1:H:230:PHE:CE2	3:Y:1:NAG:H1	2.57	0.40
1:I:94:LYS:HB3	1:I:94:LYS:HE2	1.84	0.40
1:I:182:TYR:CE1	1:I:184:ASN:HB2	2.55	0.40
1:A:19:TRP:HB3	1:A:49:PHE:CD2	2.57	0.40
1:A:314:ASN:OD1	1:B:301:SER:HB3	2.21	0.40
1:A:335:THR:HB	1:B:329:HIS:CD2	2.57	0.40
1:C:254:ALA:O	1:C:255:LEU:HB2	2.20	0.40
1:D:245:ALA:N	1:F:257:THR:O	2.54	0.40
1:E:220:ILE:HA	1:F:209:VAL:HB	2.01	0.40
1:A:85:THR:O	1:B:83:PRO:HB3	2.22	0.40
1:B:95:PRO:HD3	1:C:88:TYR:CD2	2.56	0.40
1:B:162:SER:OG	3:L:6:BDP:O6A	2.37	0.40
1:C:311:ARG:NH2	1:C:313:ARG:HH21	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:36:ASP:C	1:F:38:GLY:N	2.79	0.40
1:F:182:TYR:CE1	1:F:184:ASN:HB2	2.57	0.40
1:F:19:TRP:HB3	1:F:49:PHE:CD2	2.57	0.40
1:G:232:SER:O	1:H:216:ASN:HA	2.20	0.40
1:G:291:GLY:O	1:G:292:THR:HG23	2.21	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	359/379 (95%)	323 (90%)	26 (7%)	10 (3%)	4	26
1	B	357/379 (94%)	321 (90%)	26 (7%)	10 (3%)	4	26
1	C	359/379 (95%)	322 (90%)	32 (9%)	5 (1%)	9	38
1	D	345/379 (91%)	311 (90%)	26 (8%)	8 (2%)	5	29
1	E	353/379 (93%)	317 (90%)	29 (8%)	7 (2%)	6	32
1	F	364/379 (96%)	325 (89%)	30 (8%)	9 (2%)	4	28
1	G	361/379 (95%)	331 (92%)	25 (7%)	5 (1%)	9	38
1	H	357/379 (94%)	322 (90%)	29 (8%)	6 (2%)	7	35
1	I	360/379 (95%)	324 (90%)	31 (9%)	5 (1%)	9	38
All	All	3215/3411 (94%)	2896 (90%)	254 (8%)	65 (2%)	6	32

All (65) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	45	GLY
1	A	46	GLN
1	A	287	LYS

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Mol	Chain	Res	Type
1	A	366	LYS
1	B	37	THR
1	B	46	GLN
1	B	351	ALA
1	C	37	THR
1	D	37	THR
1	D	46	GLN
1	D	287	LYS
1	E	37	THR
1	E	287	LYS
1	E	351	ALA
1	F	37	THR
1	F	46	GLN
1	F	255	LEU
1	F	287	LYS
1	F	361	ILE
1	G	37	THR
1	G	351	ALA
1	H	46	GLN
1	H	139	LEU
1	H	287	LYS
1	I	45	GLY
1	I	287	LYS
1	A	59	LYS
1	A	255	LEU
1	B	45	GLY
1	B	53	LYS
1	B	67	LEU
1	B	84	GLY
1	C	351	ALA
1	D	45	GLY
1	F	6	PRO
1	G	255	LEU
1	H	255	LEU
1	I	37	THR
1	I	366	LYS
1	A	365	LYS
1	B	350	HIS
1	D	332	ALA
1	D	350	HIS
1	E	46	GLN
1	E	85	THR

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Mol	Chain	Res	Type
1	E	255	LEU
1	F	353	THR
1	H	138	SER
1	A	58	PRO
1	B	36	ASP
1	C	36	ASP
1	C	360	LYS
1	D	36	ASP
1	D	355	ASP
1	E	36	ASP
1	F	119	ASP
1	G	36	ASP
1	G	367	LEU
1	I	255	LEU
1	C	46	GLN
1	F	36	ASP
1	A	351	ALA
1	B	28	GLU
1	H	28	GLU
1	A	348	GLY

5.3.2 Protein sidechains ⓘ

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	288/306 (94%)	288 (100%)	0	100	100
1	B	288/306 (94%)	288 (100%)	0	100	100
1	C	288/306 (94%)	288 (100%)	0	100	100
1	D	278/306 (91%)	277 (100%)	1 (0%)	84	81
1	E	284/306 (93%)	284 (100%)	0	100	100
1	F	293/306 (96%)	293 (100%)	0	100	100
1	G	290/306 (95%)	290 (100%)	0	100	100
1	H	288/306 (94%)	288 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	I	289/306 (94%)	289 (100%)	0	100	100
All	All	2586/2754 (94%)	2585 (100%)	1 (0%)	100	100

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	292	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	229	ASN
1	A	339	ASN
1	B	147	GLN
1	B	149	GLN
1	B	339	ASN
1	C	295	GLN
1	C	339	ASN
1	D	240	ASN
1	E	10	GLN
1	E	247	GLN
1	E	274	ASN
1	G	90	GLN
1	H	68	GLN
1	I	247	GLN
1	I	339	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

108 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	NAG	J	1	2	14,14,15	0.26	0	17,19,21	0.44	0
2	BDP	J	2	2	12,12,13	1.22	2 (16%)	14,17,19	0.98	1 (7%)
2	NAG	J	3	2	14,14,15	0.25	0	17,19,21	0.77	0
2	GCD	J	4	2	10,11,12	0.34	0	12,15,17	0.45	0
3	NAG	K	1	3	15,15,15	0.12	0	21,21,21	0.51	0
3	BDP	K	2	3	12,12,13	0.15	0	14,17,19	0.41	0
3	NAG	K	3	3	14,14,15	0.34	0	17,19,21	0.58	0
3	BDP	K	4	3	12,12,13	1.23	2 (16%)	14,17,19	0.92	1 (7%)
3	NAG	K	5	3	14,14,15	0.25	0	17,19,21	0.76	0
3	BDP	K	6	3	12,12,13	1.21	2 (16%)	14,17,19	1.32	3 (21%)
3	NAG	K	7	3	14,14,15	0.31	0	17,19,21	0.97	1 (5%)
3	GCD	K	8	3	10,11,12	0.61	0	12,15,17	1.23	3 (25%)
3	NAG	L	1	3	15,15,15	0.11	0	21,21,21	0.47	0
3	BDP	L	2	3	12,12,13	0.41	0	14,17,19	1.60	3 (21%)
3	NAG	L	3	3	14,14,15	0.33	0	17,19,21	1.00	1 (5%)
3	BDP	L	4	3	12,12,13	0.55	0	14,17,19	0.60	0
3	NAG	L	5	3	14,14,15	0.26	0	17,19,21	0.84	1 (5%)
3	BDP	L	6	3	12,12,13	1.24	2 (16%)	14,17,19	0.89	0
3	NAG	L	7	3	14,14,15	0.26	0	17,19,21	0.64	0
3	GCD	L	8	3	10,11,12	0.42	0	12,15,17	0.46	0
3	NAG	M	1	3	15,15,15	0.21	0	21,21,21	0.99	2 (9%)
3	BDP	M	2	3	12,12,13	1.27	2 (16%)	14,17,19	1.08	0
3	NAG	M	3	3	14,14,15	0.28	0	17,19,21	0.74	0
3	BDP	M	4	3	12,12,13	1.23	2 (16%)	14,17,19	1.25	2 (14%)
3	NAG	M	5	3	14,14,15	0.30	0	17,19,21	1.72	4 (23%)
3	BDP	M	6	3	12,12,13	0.23	0	14,17,19	0.38	0
3	NAG	M	7	3	14,14,15	0.28	0	17,19,21	1.51	3 (17%)
3	GCD	M	8	3	10,11,12	1.57	3 (30%)	12,15,17	1.58	2 (16%)
2	NAG	N	1	2	15,15,15	0.17	0	21,21,21	0.63	0
2	BDP	N	2	2	12,12,13	1.20	2 (16%)	14,17,19	1.56	2 (14%)
2	NAG	N	3	2	14,14,15	0.33	0	17,19,21	2.08	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GCD	N	4	2	10,11,12	0.50	0	12,15,17	1.25	2 (16%)
2	NAG	O	1	2	15,15,15	0.14	0	21,21,21	0.43	0
2	BDP	O	2	2	12,12,13	1.21	2 (16%)	14,17,19	1.10	2 (14%)
2	NAG	O	3	2	14,14,15	0.66	1 (7%)	17,19,21	2.20	5 (29%)
2	GCD	O	4	2	10,11,12	1.43	2 (20%)	12,15,17	1.83	3 (25%)
3	NAG	P	1	3	15,15,15	0.14	0	21,21,21	0.21	0
3	BDP	P	2	3	12,12,13	0.20	0	14,17,19	0.53	0
3	NAG	P	3	3	14,14,15	0.43	0	17,19,21	1.41	3 (17%)
3	BDP	P	4	3	12,12,13	1.27	2 (16%)	14,17,19	1.02	1 (7%)
3	NAG	P	5	3	14,14,15	0.34	0	17,19,21	1.75	3 (17%)
3	BDP	P	6	3	12,12,13	0.34	0	14,17,19	1.07	2 (14%)
3	NAG	P	7	3	14,14,15	0.52	0	17,19,21	1.21	2 (11%)
3	GCD	P	8	3	10,11,12	0.52	0	12,15,17	0.82	0
2	NAG	Q	1	2	15,15,15	0.16	0	21,21,21	0.56	0
2	BDP	Q	2	2	12,12,13	1.37	2 (16%)	14,17,19	1.90	3 (21%)
2	NAG	Q	3	2	14,14,15	0.28	0	17,19,21	0.79	0
2	GCD	Q	4	2	10,11,12	0.46	0	12,15,17	0.38	0
3	NAG	R	1	3	15,15,15	0.28	0	21,21,21	0.85	0
3	BDP	R	2	3	12,12,13	1.35	2 (16%)	14,17,19	1.39	2 (14%)
3	NAG	R	3	3	14,14,15	0.44	0	17,19,21	0.75	0
3	BDP	R	4	3	12,12,13	1.29	2 (16%)	14,17,19	1.50	2 (14%)
3	NAG	R	5	3	14,14,15	0.31	0	17,19,21	0.92	1 (5%)
3	BDP	R	6	3	12,12,13	0.25	0	14,17,19	0.58	0
3	NAG	R	7	3	14,14,15	0.23	0	17,19,21	0.43	0
3	GCD	R	8	3	10,11,12	0.60	0	12,15,17	1.54	3 (25%)
3	NAG	S	1	3	15,15,15	0.20	0	21,21,21	0.84	2 (9%)
3	BDP	S	2	3	12,12,13	0.40	0	14,17,19	1.11	2 (14%)
3	NAG	S	3	3	14,14,15	0.49	0	17,19,21	1.80	2 (11%)
3	BDP	S	4	3	12,12,13	1.28	2 (16%)	14,17,19	1.69	3 (21%)
3	NAG	S	5	3	14,14,15	1.06	1 (7%)	17,19,21	2.21	5 (29%)
3	BDP	S	6	3	12,12,13	0.43	0	14,17,19	0.69	0
3	NAG	S	7	3	14,14,15	0.27	0	17,19,21	0.77	0
3	GCD	S	8	3	10,11,12	1.45	2 (20%)	12,15,17	1.08	1 (8%)
2	NAG	T	1	2	15,15,15	0.23	0	21,21,21	1.35	1 (4%)
2	BDP	T	2	2	12,12,13	0.48	0	14,17,19	1.33	3 (21%)
2	NAG	T	3	2	14,14,15	0.61	0	17,19,21	1.73	3 (17%)
2	GCD	T	4	2	10,11,12	1.58	2 (20%)	12,15,17	1.20	1 (8%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	U	1	2	15,15,15	0.22	0	21,21,21	0.70	1 (4%)
2	BDP	U	2	2	12,12,13	0.19	0	14,17,19	1.01	1 (7%)
2	NAG	U	3	2	14,14,15	0.28	0	17,19,21	0.87	0
2	GCD	U	4	2	10,11,12	1.56	2 (20%)	12,15,17	1.88	3 (25%)
2	NAG	V	1	2	14,14,15	0.31	0	17,19,21	0.92	1 (5%)
2	BDP	V	2	2	12,12,13	1.23	2 (16%)	14,17,19	1.41	3 (21%)
2	NAG	V	3	2	14,14,15	0.59	0	17,19,21	1.71	3 (17%)
2	GCD	V	4	2	10,11,12	0.49	0	12,15,17	1.13	1 (8%)
3	NAG	W	1	3	15,15,15	0.29	0	21,21,21	0.63	0
3	BDP	W	2	3	12,12,13	0.53	0	14,17,19	2.07	2 (14%)
3	NAG	W	3	3	14,14,15	0.44	0	17,19,21	0.94	1 (5%)
3	BDP	W	4	3	12,12,13	0.33	0	14,17,19	0.62	0
3	NAG	W	5	3	14,14,15	0.40	0	17,19,21	1.27	2 (11%)
3	BDP	W	6	3	12,12,13	1.46	2 (16%)	14,17,19	1.28	2 (14%)
3	NAG	W	7	3	14,14,15	0.28	0	17,19,21	0.95	1 (5%)
3	GCD	W	8	3	10,11,12	0.61	0	12,15,17	1.36	2 (16%)
3	NAG	X	1	3	15,15,15	0.13	0	21,21,21	0.49	0
3	BDP	X	2	3	12,12,13	0.29	0	14,17,19	0.83	1 (7%)
3	NAG	X	3	3	14,14,15	0.35	0	17,19,21	0.71	0
3	BDP	X	4	3	12,12,13	0.24	0	14,17,19	0.65	1 (7%)
3	NAG	X	5	3	14,14,15	0.36	0	17,19,21	0.91	1 (5%)
3	BDP	X	6	3	12,12,13	0.30	0	14,17,19	0.91	0
3	NAG	X	7	3	14,14,15	0.46	0	17,19,21	1.06	2 (11%)
3	GCD	X	8	3	10,11,12	0.55	0	12,15,17	0.64	0
3	NAG	Y	1	3	15,15,15	0.21	0	21,21,21	0.60	0
3	BDP	Y	2	3	12,12,13	1.22	2 (16%)	14,17,19	1.00	1 (7%)
3	NAG	Y	3	3	14,14,15	0.40	0	17,19,21	0.83	0
3	BDP	Y	4	3	12,12,13	1.23	2 (16%)	14,17,19	1.01	2 (14%)
3	NAG	Y	5	3	14,14,15	0.39	0	17,19,21	2.12	3 (17%)
3	BDP	Y	6	3	12,12,13	0.68	0	14,17,19	1.25	2 (14%)
3	NAG	Y	7	3	14,14,15	0.27	0	17,19,21	0.87	0
3	GCD	Y	8	3	10,11,12	1.95	3 (30%)	12,15,17	0.95	0
2	NAG	Z	1	2	15,15,15	0.13	0	21,21,21	0.43	0
2	BDP	Z	2	2	12,12,13	0.53	0	14,17,19	1.20	2 (14%)
2	NAG	Z	3	2	14,14,15	0.62	0	17,19,21	2.25	4 (23%)
2	GCD	Z	4	2	10,11,12	0.46	0	12,15,17	1.23	2 (16%)
2	NAG	a	1	2	15,15,15	0.22	0	21,21,21	0.84	1 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BDP	a	2	2	12,12,13	0.27	0	14,17,19	0.78	0
2	NAG	a	3	2	14,14,15	0.40	0	17,19,21	0.98	1 (5%)
2	GCD	a	4	2	10,11,12	1.09	1 (10%)	12,15,17	1.41	4 (33%)

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	NAG	J	1	2	-	4/6/23/26	0/1/1/1
2	BDP	J	2	2	-	0/4/21/24	0/1/1/1
2	NAG	J	3	2	-	0/6/23/26	0/1/1/1
2	GCD	J	4	2	-	0/4/17/20	0/1/1/1
3	NAG	K	1	3	-	3/6/26/26	0/1/1/1
3	BDP	K	2	3	-	2/4/21/24	0/1/1/1
3	NAG	K	3	3	-	2/6/23/26	0/1/1/1
3	BDP	K	4	3	-	0/4/21/24	0/1/1/1
3	NAG	K	5	3	-	4/6/23/26	0/1/1/1
3	BDP	K	6	3	-	0/4/21/24	0/1/1/1
3	NAG	K	7	3	-	5/6/23/26	0/1/1/1
3	GCD	K	8	3	-	0/4/17/20	0/1/1/1
3	NAG	L	1	3	-	0/6/26/26	0/1/1/1
3	BDP	L	2	3	-	2/4/21/24	0/1/1/1
3	NAG	L	3	3	-	4/6/23/26	0/1/1/1
3	BDP	L	4	3	-	0/4/21/24	1/1/1/1
3	NAG	L	5	3	-	0/6/23/26	0/1/1/1
3	BDP	L	6	3	-	0/4/21/24	0/1/1/1
3	NAG	L	7	3	-	4/6/23/26	0/1/1/1
3	GCD	L	8	3	-	0/4/17/20	0/1/1/1
3	NAG	M	1	3	-	4/6/26/26	0/1/1/1
3	BDP	M	2	3	-	0/4/21/24	0/1/1/1
3	NAG	M	3	3	-	1/6/23/26	0/1/1/1
3	BDP	M	4	3	-	3/4/21/24	0/1/1/1
3	NAG	M	5	3	-	3/6/23/26	0/1/1/1
3	BDP	M	6	3	-	0/4/21/24	1/1/1/1
3	NAG	M	7	3	-	3/6/23/26	0/1/1/1
3	GCD	M	8	3	-	0/4/17/20	0/1/1/1
2	NAG	N	1	2	-	6/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BDP	N	2	2	-	1/4/21/24	1/1/1/1
2	NAG	N	3	2	-	4/6/23/26	0/1/1/1
2	GCD	N	4	2	-	0/4/17/20	0/1/1/1
2	NAG	O	1	2	-	3/6/26/26	0/1/1/1
2	BDP	O	2	2	-	0/4/21/24	0/1/1/1
2	NAG	O	3	2	-	3/6/23/26	0/1/1/1
2	GCD	O	4	2	-	0/4/17/20	0/1/1/1
3	NAG	P	1	3	-	2/6/26/26	0/1/1/1
3	BDP	P	2	3	-	2/4/21/24	0/1/1/1
3	NAG	P	3	3	-	1/6/23/26	0/1/1/1
3	BDP	P	4	3	-	0/4/21/24	0/1/1/1
3	NAG	P	5	3	-	2/6/23/26	0/1/1/1
3	BDP	P	6	3	-	0/4/21/24	0/1/1/1
3	NAG	P	7	3	-	5/6/23/26	0/1/1/1
3	GCD	P	8	3	-	0/4/17/20	0/1/1/1
2	NAG	Q	1	2	-	4/6/26/26	0/1/1/1
2	BDP	Q	2	2	-	0/4/21/24	0/1/1/1
2	NAG	Q	3	2	-	2/6/23/26	0/1/1/1
2	GCD	Q	4	2	-	0/4/17/20	0/1/1/1
3	NAG	R	1	3	-	2/6/26/26	0/1/1/1
3	BDP	R	2	3	-	0/4/21/24	0/1/1/1
3	NAG	R	3	3	-	2/6/23/26	0/1/1/1
3	BDP	R	4	3	-	0/4/21/24	0/1/1/1
3	NAG	R	5	3	-	4/6/23/26	0/1/1/1
3	BDP	R	6	3	-	0/4/21/24	1/1/1/1
3	NAG	R	7	3	-	1/6/23/26	0/1/1/1
3	GCD	R	8	3	-	0/4/17/20	0/1/1/1
3	NAG	S	1	3	-	5/6/26/26	0/1/1/1
3	BDP	S	2	3	-	2/4/21/24	0/1/1/1
3	NAG	S	3	3	-	2/6/23/26	0/1/1/1
3	BDP	S	4	3	-	2/4/21/24	0/1/1/1
3	NAG	S	5	3	-	2/6/23/26	0/1/1/1
3	BDP	S	6	3	-	0/4/21/24	0/1/1/1
3	NAG	S	7	3	-	4/6/23/26	0/1/1/1
3	GCD	S	8	3	-	0/4/17/20	0/1/1/1
2	NAG	T	1	2	-	3/6/26/26	0/1/1/1
2	BDP	T	2	2	-	0/4/21/24	0/1/1/1
2	NAG	T	3	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GCD	T	4	2	-	0/4/17/20	0/1/1/1
2	NAG	U	1	2	-	3/6/26/26	0/1/1/1
2	BDP	U	2	2	-	0/4/21/24	1/1/1/1
2	NAG	U	3	2	-	4/6/23/26	0/1/1/1
2	GCD	U	4	2	-	0/4/17/20	0/1/1/1
2	NAG	V	1	2	-	5/6/23/26	0/1/1/1
2	BDP	V	2	2	-	2/4/21/24	1/1/1/1
2	NAG	V	3	2	-	2/6/23/26	0/1/1/1
2	GCD	V	4	2	-	0/4/17/20	0/1/1/1
3	NAG	W	1	3	-	6/6/26/26	0/1/1/1
3	BDP	W	2	3	-	0/4/21/24	0/1/1/1
3	NAG	W	3	3	-	4/6/23/26	0/1/1/1
3	BDP	W	4	3	-	2/4/21/24	0/1/1/1
3	NAG	W	5	3	-	5/6/23/26	0/1/1/1
3	BDP	W	6	3	-	2/4/21/24	0/1/1/1
3	NAG	W	7	3	-	4/6/23/26	0/1/1/1
3	GCD	W	8	3	-	0/4/17/20	0/1/1/1
3	NAG	X	1	3	-	1/6/26/26	0/1/1/1
3	BDP	X	2	3	-	2/4/21/24	0/1/1/1
3	NAG	X	3	3	-	4/6/23/26	0/1/1/1
3	BDP	X	4	3	-	0/4/21/24	0/1/1/1
3	NAG	X	5	3	-	2/6/23/26	0/1/1/1
3	BDP	X	6	3	-	2/4/21/24	1/1/1/1
3	NAG	X	7	3	-	4/6/23/26	0/1/1/1
3	GCD	X	8	3	-	0/4/17/20	0/1/1/1
3	NAG	Y	1	3	-	5/6/26/26	0/1/1/1
3	BDP	Y	2	3	-	0/4/21/24	0/1/1/1
3	NAG	Y	3	3	-	1/6/23/26	0/1/1/1
3	BDP	Y	4	3	-	2/4/21/24	1/1/1/1
3	NAG	Y	5	3	-	4/6/23/26	0/1/1/1
3	BDP	Y	6	3	-	0/4/21/24	1/1/1/1
3	NAG	Y	7	3	-	0/6/23/26	0/1/1/1
3	GCD	Y	8	3	-	0/4/17/20	0/1/1/1
2	NAG	Z	1	2	-	3/6/26/26	0/1/1/1
2	BDP	Z	2	2	-	0/4/21/24	0/1/1/1
2	NAG	Z	3	2	-	5/6/23/26	0/1/1/1
2	GCD	Z	4	2	-	0/4/17/20	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	a	1	2	-	6/6/26/26	0/1/1/1
2	BDP	a	2	2	-	2/4/21/24	0/1/1/1
2	NAG	a	3	2	-	4/6/23/26	0/1/1/1
2	GCD	a	4	2	-	3/4/17/20	0/1/1/1

All (51) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	Y	8	GCD	O6A-C6	4.08	1.32	1.22
2	T	4	GCD	O6A-C6	4.02	1.32	1.22
2	U	4	GCD	O6A-C6	3.65	1.31	1.22
3	S	8	GCD	O6A-C6	3.58	1.31	1.22
2	O	4	GCD	O6A-C6	3.56	1.31	1.22
3	Y	8	GCD	O6B-C6	-3.34	1.21	1.30
3	S	4	BDP	O6A-C6	3.25	1.31	1.22
3	S	5	NAG	C8-C7	3.22	1.57	1.50
3	P	4	BDP	O6A-C6	3.18	1.31	1.22
3	R	4	BDP	O6A-C6	3.16	1.31	1.22
3	M	8	GCD	C5-C6	-3.15	1.41	1.48
3	Y	8	GCD	C5-C6	-3.12	1.41	1.48
2	Q	2	BDP	O6A-C6	3.09	1.31	1.22
2	J	2	BDP	O6A-C6	3.08	1.31	1.22
3	K	4	BDP	O6A-C6	3.08	1.31	1.22
3	Y	4	BDP	O6A-C6	3.07	1.31	1.22
3	R	2	BDP	O6A-C6	3.07	1.31	1.22
3	W	6	BDP	O6A-C6	3.04	1.31	1.22
3	K	6	BDP	O6A-C6	3.03	1.31	1.22
2	O	2	BDP	O6A-C6	3.03	1.31	1.22
3	L	6	BDP	O6A-C6	3.03	1.31	1.22
3	M	4	BDP	O6A-C6	3.03	1.31	1.22
2	V	2	BDP	O6A-C6	3.01	1.31	1.22
3	M	2	BDP	O6A-C6	2.97	1.30	1.22
3	Y	2	BDP	O6A-C6	2.96	1.30	1.22
3	L	6	BDP	O6B-C6	-2.95	1.21	1.30
2	a	4	GCD	O6B-C6	2.94	1.38	1.30
2	N	2	BDP	O6A-C6	2.94	1.30	1.22
3	S	4	BDP	O6B-C6	-2.91	1.21	1.30
2	U	4	GCD	O6B-C6	-2.89	1.22	1.30
3	R	4	BDP	O6B-C6	-2.87	1.21	1.30
3	P	4	BDP	O6B-C6	-2.86	1.21	1.30
3	Y	4	BDP	O6B-C6	-2.85	1.21	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	W	6	BDP	O6B-C6	-2.83	1.21	1.30
3	Y	2	BDP	O6B-C6	-2.83	1.21	1.30
2	Q	2	BDP	O6B-C6	-2.83	1.21	1.30
2	N	2	BDP	O6B-C6	-2.82	1.21	1.30
3	S	8	GCD	O6B-C6	-2.82	1.22	1.30
2	J	2	BDP	O6B-C6	-2.78	1.21	1.30
3	K	6	BDP	O6B-C6	-2.78	1.21	1.30
3	M	4	BDP	O6B-C6	-2.78	1.21	1.30
3	R	2	BDP	O6B-C6	-2.76	1.21	1.30
2	O	2	BDP	O6B-C6	-2.76	1.21	1.30
3	K	4	BDP	O6B-C6	-2.75	1.21	1.30
2	V	2	BDP	O6B-C6	-2.74	1.21	1.30
3	M	8	GCD	O6B-C6	-2.67	1.23	1.30
3	M	2	BDP	O6B-C6	-2.66	1.22	1.30
3	M	8	GCD	O6A-C6	2.65	1.29	1.22
2	T	4	GCD	O6B-C6	-2.61	1.23	1.30
2	O	4	GCD	O6B-C6	-2.37	1.24	1.30
2	O	3	NAG	C1-C2	2.00	1.55	1.52

All (140) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Y	5	NAG	C1-O5-C5	7.37	122.07	112.19
3	S	5	NAG	C1-O5-C5	6.90	121.44	112.19
3	S	3	NAG	C1-O5-C5	6.62	121.06	112.19
3	W	2	BDP	O4-C4-C5	6.37	124.30	109.76
2	Z	3	NAG	C1-O5-C5	6.32	120.66	112.19
2	O	3	NAG	C1-O5-C5	6.11	120.37	112.19
2	T	3	NAG	C4-C3-C2	5.78	119.49	111.02
2	Z	3	NAG	C2-N2-C7	5.30	130.01	122.90
2	Q	2	BDP	C1-C2-C3	5.04	116.98	109.64
2	T	1	NAG	C4-C3-C2	4.91	117.55	110.40
2	V	3	NAG	C1-O5-C5	-4.80	105.75	112.19
2	N	2	BDP	O5-C5-C6	-4.75	89.61	106.59
3	P	5	NAG	C1-O5-C5	4.73	118.52	112.19
2	N	3	NAG	C4-C3-C2	-4.67	104.17	111.02
2	N	3	NAG	C1-O5-C5	4.56	118.29	112.19
3	L	2	BDP	C1-C2-C3	4.25	115.84	109.64
2	U	4	GCD	C1-C2-C3	3.99	115.45	109.64
3	S	4	BDP	C1-C2-C3	3.71	115.05	109.64
3	W	5	NAG	O3-C3-C2	3.66	117.01	109.40
3	M	5	NAG	C1-O5-C5	3.61	117.02	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	2	BDP	C2-C3-C4	3.51	117.04	110.86
3	M	8	GCD	O6A-C6-C5	-3.50	111.99	120.57
3	R	8	GCD	C1-C2-C3	3.37	114.56	109.64
3	M	8	GCD	O6B-C6-C5	3.34	121.62	114.06
2	O	4	GCD	C1-C2-C3	3.32	114.47	109.64
2	O	3	NAG	C4-C3-C2	3.30	115.86	111.02
3	M	7	NAG	C1-O5-C5	3.28	116.58	112.19
2	U	2	BDP	O4-C4-C5	3.26	117.21	109.76
3	P	3	NAG	C3-C4-C5	3.26	116.14	110.23
2	a	3	NAG	O3-C3-C2	3.24	116.12	109.40
2	N	3	NAG	O3-C3-C2	3.20	116.05	109.40
2	O	3	NAG	C2-N2-C7	3.19	127.18	122.90
3	M	5	NAG	C3-C4-C5	3.17	115.97	110.23
2	V	2	BDP	O4-C4-C5	3.16	116.97	109.76
3	W	8	GCD	C1-C2-C3	3.16	114.24	109.64
2	O	4	GCD	O6B-C6-C5	3.15	121.20	114.06
3	P	5	NAG	C3-C4-C5	3.14	115.93	110.23
3	P	7	NAG	C2-N2-C7	3.09	127.05	122.90
3	S	2	BDP	O4-C4-C3	3.08	117.64	110.38
2	N	3	NAG	C2-N2-C7	3.06	127.01	122.90
2	Z	2	BDP	O4-C4-C5	-3.05	102.79	109.76
3	M	1	NAG	O3-C3-C4	3.03	117.51	110.38
2	V	3	NAG	O4-C4-C5	-2.99	101.97	109.32
3	M	5	NAG	C1-C2-N2	2.98	115.13	110.43
3	S	5	NAG	C2-N2-C7	-2.94	118.95	122.90
2	U	4	GCD	C3-C4-C5	-2.93	116.49	121.45
3	M	7	NAG	C3-C4-C5	2.87	115.44	110.23
3	P	3	NAG	C4-C3-C2	2.85	115.20	111.02
3	W	2	BDP	C1-C2-C3	2.84	113.78	109.64
3	L	3	NAG	C1-O5-C5	2.82	115.97	112.19
3	P	7	NAG	C4-C3-C2	-2.82	106.89	111.02
3	K	8	GCD	C1-O5-C5	2.80	121.36	115.42
3	S	8	GCD	O6B-C6-C5	2.80	120.40	114.06
3	R	4	BDP	C1-C2-C3	2.78	113.69	109.64
2	V	1	NAG	C2-N2-C7	2.78	126.63	122.90
3	Y	6	BDP	C1-C2-C3	2.78	113.69	109.64
2	N	4	GCD	O3-C3-C2	2.77	114.38	109.44
3	R	4	BDP	O4-C4-C3	2.76	116.88	110.38
2	U	4	GCD	O6B-C6-C5	2.73	120.26	114.06
3	S	4	BDP	O4-C4-C5	2.72	115.97	109.76
3	K	6	BDP	O4-C4-C3	-2.72	103.97	110.38
2	T	3	NAG	C1-O5-C5	-2.70	108.56	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	2	BDP	O5-C5-C6	-2.66	97.09	106.59
3	R	2	BDP	O4-C4-C3	-2.62	104.19	110.38
3	M	5	NAG	C2-N2-C7	2.62	126.41	122.90
3	R	8	GCD	C3-C4-C5	-2.62	117.03	121.45
2	V	3	NAG	C3-C4-C5	2.61	114.96	110.23
2	T	2	BDP	C2-C3-C4	-2.57	106.34	110.86
3	Y	5	NAG	C4-C3-C2	2.56	114.78	111.02
2	O	4	GCD	O6A-C6-C5	-2.56	114.31	120.57
2	a	4	GCD	C1-C2-C3	-2.55	105.93	109.64
3	M	4	BDP	O6B-C6-C5	2.54	122.80	113.64
2	Z	3	NAG	C1-C2-N2	-2.53	106.45	110.43
3	W	7	NAG	C1-O5-C5	2.51	115.55	112.19
2	O	2	BDP	O4-C4-C5	-2.51	104.03	109.76
2	V	4	GCD	C1-C2-C3	-2.51	106.00	109.64
3	P	6	BDP	O5-C5-C6	-2.50	97.65	106.59
3	M	1	NAG	O3-C3-C2	-2.49	104.63	109.58
2	V	2	BDP	O6B-C6-C5	2.48	122.58	113.64
3	L	2	BDP	C3-C4-C5	-2.47	105.05	109.30
2	Z	4	GCD	C3-C4-C5	-2.43	117.35	121.45
3	X	5	NAG	C3-C4-C5	2.43	114.63	110.23
3	Y	4	BDP	O6B-C6-C5	2.41	122.36	113.64
3	P	5	NAG	O5-C5-C4	2.40	116.68	110.83
2	Z	3	NAG	O3-C3-C4	-2.38	104.77	110.38
2	T	2	BDP	O5-C1-C2	2.37	116.44	110.79
3	K	6	BDP	O6B-C6-C5	2.37	122.20	113.64
2	O	3	NAG	C1-C2-N2	2.37	114.16	110.43
2	a	1	NAG	C1-C2-N2	2.36	113.46	110.73
3	S	1	NAG	C4-C3-C2	-2.35	106.98	110.40
2	O	3	NAG	O3-C3-C2	2.31	114.19	109.40
3	R	8	GCD	O5-C5-C4	-2.30	122.83	124.94
2	Z	4	GCD	C1-O5-C5	2.29	120.27	115.42
3	P	6	BDP	O4-C4-C3	-2.28	105.00	110.38
2	N	4	GCD	C1-C2-C3	-2.26	106.35	109.64
3	P	3	NAG	C1-O5-C5	-2.26	109.16	112.19
3	X	2	BDP	O4-C4-C5	-2.24	104.65	109.76
3	W	6	BDP	O6B-C6-C5	2.23	121.69	113.64
3	K	8	GCD	C3-C4-C5	-2.23	117.68	121.45
2	J	2	BDP	O6B-C6-C5	2.22	121.66	113.64
2	a	4	GCD	O2-C2-C1	2.20	114.27	109.22
3	M	4	BDP	O6A-C6-C5	-2.20	112.87	120.81
3	R	5	NAG	C1-C2-N2	-2.20	106.97	110.43
3	L	2	BDP	O4-C4-C5	2.19	114.77	109.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	4	BDP	O6B-C6-C5	2.19	121.55	113.64
3	S	5	NAG	C4-C3-C2	2.19	114.22	111.02
3	Y	2	BDP	O6B-C6-C5	2.18	121.50	113.64
3	S	2	BDP	C1-C2-C3	2.17	112.81	109.64
2	T	3	NAG	O5-C5-C4	-2.16	105.57	110.83
2	U	1	NAG	C1-C2-N2	2.16	113.23	110.73
2	Q	2	BDP	O6B-C6-C5	2.16	121.42	113.64
3	S	1	NAG	C1-C2-N2	2.16	113.23	110.73
3	W	6	BDP	O4-C4-C5	2.15	114.67	109.76
3	L	5	NAG	C1-O5-C5	2.14	115.06	112.19
3	X	4	BDP	O5-C5-C6	-2.14	98.93	106.59
3	X	7	NAG	C1-O5-C5	2.14	115.06	112.19
3	S	4	BDP	C3-C4-C5	-2.13	105.64	109.30
3	Y	4	BDP	O6A-C6-C5	-2.13	113.13	120.81
3	M	7	NAG	C2-N2-C7	2.12	125.74	122.90
2	V	2	BDP	O6A-C6-C5	-2.12	113.16	120.81
3	K	7	NAG	C4-C3-C2	-2.11	107.92	111.02
2	a	4	GCD	C1-O5-C5	-2.11	110.95	115.42
2	N	2	BDP	O6B-C6-C5	2.11	121.25	113.64
3	R	2	BDP	C1-C2-C3	2.11	112.71	109.64
2	O	2	BDP	O6B-C6-C5	2.11	121.24	113.64
2	T	4	GCD	C1-O5-C5	2.10	119.88	115.42
3	S	5	NAG	C6-C5-C4	2.10	118.18	113.02
3	K	4	BDP	O6B-C6-C5	2.09	121.17	113.64
3	W	8	GCD	C1-O5-C5	-2.08	111.02	115.42
2	a	4	GCD	O6A-C6-C5	-2.08	115.48	120.57
3	Y	5	NAG	O3-C3-C4	-2.07	105.49	110.38
3	S	5	NAG	O3-C3-C4	-2.06	105.52	110.38
2	Z	2	BDP	C1-C2-C3	2.05	112.63	109.64
3	S	3	NAG	C4-C3-C2	-2.04	108.02	111.02
3	W	5	NAG	C4-C3-C2	2.04	114.01	111.02
3	K	8	GCD	C1-C2-C3	2.03	112.61	109.64
3	K	6	BDP	O6A-C6-C5	-2.03	113.47	120.81
3	W	3	NAG	O3-C3-C2	2.03	113.62	109.40
3	Y	6	BDP	O4-C4-C3	-2.01	105.63	110.38
3	X	7	NAG	C1-C2-N2	2.01	113.60	110.43

There are no chirality outliers.

All (199) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	N	1	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	N	1	NAG	O7-C7-N2-C2
2	O	3	NAG	C3-C2-N2-C7
2	Q	1	NAG	C1-C2-N2-C7
2	Q	1	NAG	C8-C7-N2-C2
2	Q	1	NAG	O7-C7-N2-C2
2	T	1	NAG	C1-C2-N2-C7
2	T	1	NAG	C8-C7-N2-C2
2	T	1	NAG	O7-C7-N2-C2
2	T	3	NAG	C8-C7-N2-C2
2	T	3	NAG	O7-C7-N2-C2
2	U	1	NAG	C1-C2-N2-C7
2	U	1	NAG	O7-C7-N2-C2
2	V	1	NAG	C3-C2-N2-C7
2	V	2	BDP	C4-C5-C6-O6A
2	V	2	BDP	C4-C5-C6-O6B
2	Z	1	NAG	C1-C2-N2-C7
2	Z	3	NAG	C3-C2-N2-C7
2	a	1	NAG	C1-C2-N2-C7
2	a	2	BDP	C4-C5-C6-O6A
2	a	2	BDP	C4-C5-C6-O6B
2	a	4	GCD	O5-C5-C6-O6A
2	a	4	GCD	O5-C5-C6-O6B
3	K	2	BDP	C4-C5-C6-O6A
3	K	2	BDP	C4-C5-C6-O6B
3	K	7	NAG	C8-C7-N2-C2
3	K	7	NAG	O7-C7-N2-C2
3	L	2	BDP	C4-C5-C6-O6A
3	L	2	BDP	C4-C5-C6-O6B
3	L	7	NAG	C8-C7-N2-C2
3	L	7	NAG	O7-C7-N2-C2
3	M	1	NAG	C1-C2-N2-C7
3	M	1	NAG	C8-C7-N2-C2
3	M	1	NAG	O7-C7-N2-C2
3	M	4	BDP	C4-C5-C6-O6A
3	M	4	BDP	C4-C5-C6-O6B
3	M	5	NAG	C1-C2-N2-C7
3	M	5	NAG	C8-C7-N2-C2
3	M	5	NAG	O7-C7-N2-C2
3	M	7	NAG	C3-C2-N2-C7
3	P	2	BDP	C4-C5-C6-O6A
3	P	2	BDP	C4-C5-C6-O6B
3	P	7	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	P	7	NAG	O7-C7-N2-C2
3	R	5	NAG	C3-C2-N2-C7
3	R	5	NAG	C8-C7-N2-C2
3	R	5	NAG	O7-C7-N2-C2
3	S	2	BDP	C4-C5-C6-O6A
3	S	2	BDP	C4-C5-C6-O6B
3	S	4	BDP	C4-C5-C6-O6A
3	S	4	BDP	C4-C5-C6-O6B
3	W	1	NAG	C1-C2-N2-C7
3	W	3	NAG	C8-C7-N2-C2
3	W	3	NAG	O7-C7-N2-C2
3	W	4	BDP	C4-C5-C6-O6A
3	W	5	NAG	C3-C2-N2-C7
3	W	6	BDP	C4-C5-C6-O6A
3	W	6	BDP	C4-C5-C6-O6B
3	W	7	NAG	C8-C7-N2-C2
3	W	7	NAG	O7-C7-N2-C2
3	X	2	BDP	C4-C5-C6-O6A
3	X	2	BDP	C4-C5-C6-O6B
3	Y	1	NAG	C1-C2-N2-C7
3	Y	1	NAG	C8-C7-N2-C2
3	Y	1	NAG	O7-C7-N2-C2
2	J	1	NAG	C4-C5-C6-O6
2	Z	3	NAG	O5-C5-C6-O6
2	O	1	NAG	C8-C7-N2-C2
2	O	1	NAG	O7-C7-N2-C2
2	U	1	NAG	C8-C7-N2-C2
2	U	3	NAG	C8-C7-N2-C2
2	U	3	NAG	O7-C7-N2-C2
2	V	1	NAG	O7-C7-N2-C2
2	V	3	NAG	O5-C5-C6-O6
3	L	3	NAG	O5-C5-C6-O6
3	P	7	NAG	O5-C5-C6-O6
2	J	1	NAG	O5-C5-C6-O6
2	V	1	NAG	C8-C7-N2-C2
2	Z	3	NAG	O7-C7-N2-C2
3	K	5	NAG	C4-C5-C6-O6
3	P	7	NAG	C4-C5-C6-O6
3	X	3	NAG	C4-C5-C6-O6
3	K	1	NAG	O5-C5-C6-O6
2	Z	3	NAG	C4-C5-C6-O6
3	X	7	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	X	7	NAG	C4-C5-C6-O6
3	L	3	NAG	C4-C5-C6-O6
2	V	1	NAG	O5-C5-C6-O6
3	L	7	NAG	O5-C5-C6-O6
2	V	3	NAG	C4-C5-C6-O6
3	S	3	NAG	C4-C5-C6-O6
2	N	1	NAG	O5-C5-C6-O6
3	R	3	NAG	C8-C7-N2-C2
3	X	7	NAG	C8-C7-N2-C2
3	L	7	NAG	C4-C5-C6-O6
3	W	1	NAG	C4-C5-C6-O6
3	K	3	NAG	O5-C5-C6-O6
3	X	3	NAG	O5-C5-C6-O6
2	V	1	NAG	C4-C5-C6-O6
2	U	3	NAG	C4-C5-C6-O6
3	K	1	NAG	C4-C5-C6-O6
2	Z	3	NAG	C8-C7-N2-C2
3	R	3	NAG	O7-C7-N2-C2
3	W	1	NAG	C8-C7-N2-C2
3	W	1	NAG	C3-C2-N2-C7
2	N	1	NAG	C4-C5-C6-O6
3	K	5	NAG	O5-C5-C6-O6
3	S	3	NAG	O5-C5-C6-O6
2	J	1	NAG	C8-C7-N2-C2
2	J	1	NAG	O7-C7-N2-C2
2	N	3	NAG	C8-C7-N2-C2
2	O	3	NAG	C8-C7-N2-C2
2	O	3	NAG	O7-C7-N2-C2
2	Q	3	NAG	C8-C7-N2-C2
3	S	5	NAG	C8-C7-N2-C2
3	S	7	NAG	C8-C7-N2-C2
3	S	7	NAG	O7-C7-N2-C2
3	X	3	NAG	C8-C7-N2-C2
3	X	5	NAG	C8-C7-N2-C2
3	X	5	NAG	O7-C7-N2-C2
3	X	7	NAG	O7-C7-N2-C2
3	W	3	NAG	C4-C5-C6-O6
3	W	5	NAG	C4-C5-C6-O6
2	N	3	NAG	O7-C7-N2-C2
3	W	1	NAG	O7-C7-N2-C2
2	Z	1	NAG	O5-C5-C6-O6
2	a	3	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	a	1	NAG	O5-C5-C6-O6
3	M	7	NAG	C8-C7-N2-C2
3	S	5	NAG	O7-C7-N2-C2
3	W	1	NAG	O5-C5-C6-O6
3	W	3	NAG	O5-C5-C6-O6
3	K	3	NAG	C4-C5-C6-O6
3	W	5	NAG	C8-C7-N2-C2
3	Y	5	NAG	C8-C7-N2-C2
3	S	1	NAG	O5-C5-C6-O6
2	Q	1	NAG	O5-C5-C6-O6
3	X	3	NAG	O7-C7-N2-C2
3	Y	3	NAG	O5-C5-C6-O6
3	S	7	NAG	C4-C5-C6-O6
2	U	3	NAG	O5-C5-C6-O6
2	Q	3	NAG	O7-C7-N2-C2
2	N	1	NAG	C1-C2-N2-C7
3	S	1	NAG	C1-C2-N2-C7
3	S	7	NAG	O5-C5-C6-O6
3	P	3	NAG	O5-C5-C6-O6
2	O	1	NAG	O5-C5-C6-O6
3	M	3	NAG	O5-C5-C6-O6
3	M	7	NAG	O7-C7-N2-C2
3	K	7	NAG	O5-C5-C6-O6
3	W	5	NAG	O5-C5-C6-O6
3	P	7	NAG	C3-C2-N2-C7
3	R	5	NAG	O5-C5-C6-O6
3	W	7	NAG	C4-C5-C6-O6
3	W	5	NAG	O7-C7-N2-C2
3	R	7	NAG	O5-C5-C6-O6
3	Y	5	NAG	O7-C7-N2-C2
3	M	1	NAG	C3-C2-N2-C7
2	a	3	NAG	C1-C2-N2-C7
3	K	7	NAG	C1-C2-N2-C7
2	a	1	NAG	C4-C5-C6-O6
3	P	1	NAG	C1-C2-N2-C7
2	a	4	GCD	C4-C5-C6-O6A
2	a	1	NAG	C8-C7-N2-C2
3	R	1	NAG	C8-C7-N2-C2
2	a	3	NAG	C4-C5-C6-O6
3	Y	1	NAG	O5-C5-C6-O6
2	N	2	BDP	C4-C5-C6-O6A
3	W	4	BDP	C4-C5-C6-O6B

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Mol	Chain	Res	Type	Atoms
3	X	6	BDP	C4-C5-C6-O6A
3	X	6	BDP	C4-C5-C6-O6B
2	N	3	NAG	C3-C2-N2-C7
2	a	3	NAG	C3-C2-N2-C7
3	K	5	NAG	C3-C2-N2-C7
3	K	7	NAG	C3-C2-N2-C7
3	L	3	NAG	C3-C2-N2-C7
3	P	5	NAG	C3-C2-N2-C7
3	Y	5	NAG	C3-C2-N2-C7
2	N	3	NAG	C4-C5-C6-O6
2	a	1	NAG	O7-C7-N2-C2
2	a	1	NAG	C3-C2-N2-C7
3	S	1	NAG	C4-C5-C6-O6
3	K	5	NAG	C1-C2-N2-C7
3	L	3	NAG	C1-C2-N2-C7
3	P	5	NAG	C1-C2-N2-C7
3	Y	5	NAG	C1-C2-N2-C7
3	K	1	NAG	C1-C2-N2-C7
3	Y	1	NAG	C3-C2-N2-C7
3	S	1	NAG	C8-C7-N2-C2
3	W	7	NAG	O5-C5-C6-O6
3	X	1	NAG	O5-C5-C6-O6
2	Z	1	NAG	C3-C2-N2-C7
3	R	1	NAG	O7-C7-N2-C2
3	P	1	NAG	C3-C2-N2-C7
3	S	1	NAG	C3-C2-N2-C7
3	M	4	BDP	O5-C5-C6-O6B
3	Y	4	BDP	O5-C5-C6-O6A
3	Y	4	BDP	O5-C5-C6-O6B
2	N	1	NAG	C3-C2-N2-C7

All (9) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	M	6	BDP	C1-C2-C3-C4-C5-O5
3	Y	6	BDP	C1-C2-C3-C4-C5-O5
2	U	2	BDP	C1-C2-C3-C4-C5-O5
2	V	2	BDP	C1-C2-C3-C4-C5-O5
3	R	6	BDP	C1-C2-C3-C4-C5-O5
3	Y	4	BDP	C1-C2-C3-C4-C5-O5
3	L	4	BDP	C1-C2-C3-C4-C5-O5
3	X	6	BDP	C1-C2-C3-C4-C5-O5

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Mol	Chain	Res	Type	Atoms
2	N	2	BDP	C1-C2-C3-C4-C5-O5

56 monomers are involved in 88 short contacts:

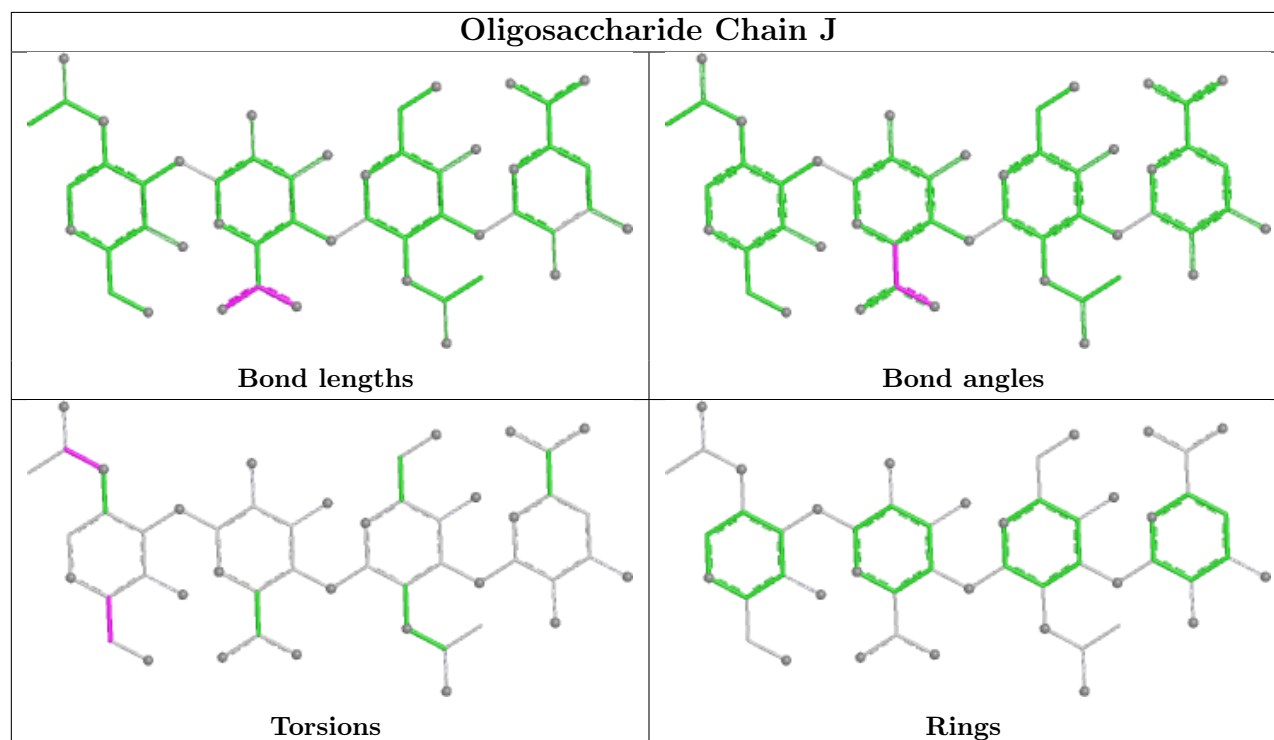
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	X	4	BDP	2	0
2	O	2	BDP	2	0
3	P	8	GCD	1	0
3	Y	3	NAG	3	0
3	R	6	BDP	1	0
3	K	1	NAG	3	0
3	L	6	BDP	1	0
2	U	1	NAG	3	0
3	S	5	NAG	3	0
3	Y	5	NAG	1	0
3	R	1	NAG	3	0
3	P	1	NAG	3	0
3	K	7	NAG	2	0
3	S	4	BDP	2	0
2	T	1	NAG	1	0
3	R	8	GCD	1	0
3	W	1	NAG	4	0
2	O	3	NAG	5	0
2	T	2	BDP	1	0
3	R	7	NAG	1	0
2	Z	2	BDP	1	0
3	P	2	BDP	1	0
2	V	1	NAG	2	0
3	X	3	NAG	1	0
3	X	5	NAG	1	0
3	P	4	BDP	2	0
3	M	1	NAG	7	0
2	U	2	BDP	3	0
3	S	6	BDP	1	0
3	X	6	BDP	1	0
2	N	2	BDP	1	0
3	S	1	NAG	2	0
2	T	4	GCD	2	0
3	K	8	GCD	1	0
3	Y	4	BDP	1	0
2	T	3	NAG	1	0
3	Y	7	NAG	1	0
3	Y	8	GCD	1	0

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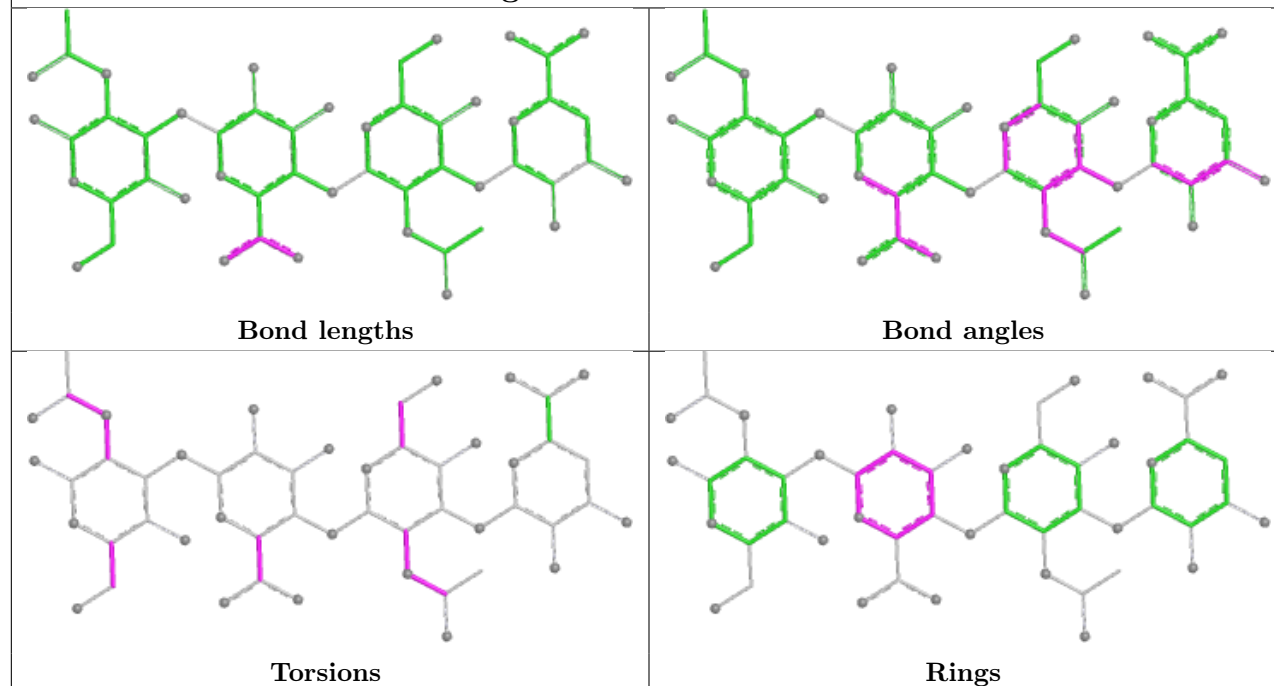
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	V	2	BDP	2	0
3	W	5	NAG	1	0
2	a	2	BDP	1	0
3	P	5	NAG	2	0
2	U	3	NAG	2	0
3	M	7	NAG	1	0
3	X	1	NAG	2	0
3	R	4	BDP	1	0
3	W	6	BDP	1	0
3	P	3	NAG	2	0
3	P	7	NAG	2	0
3	W	2	BDP	2	0
2	Z	3	NAG	1	0
2	N	3	NAG	2	0
3	Y	1	NAG	5	0
3	P	6	BDP	1	0
2	a	1	NAG	1	0
3	R	2	BDP	2	0

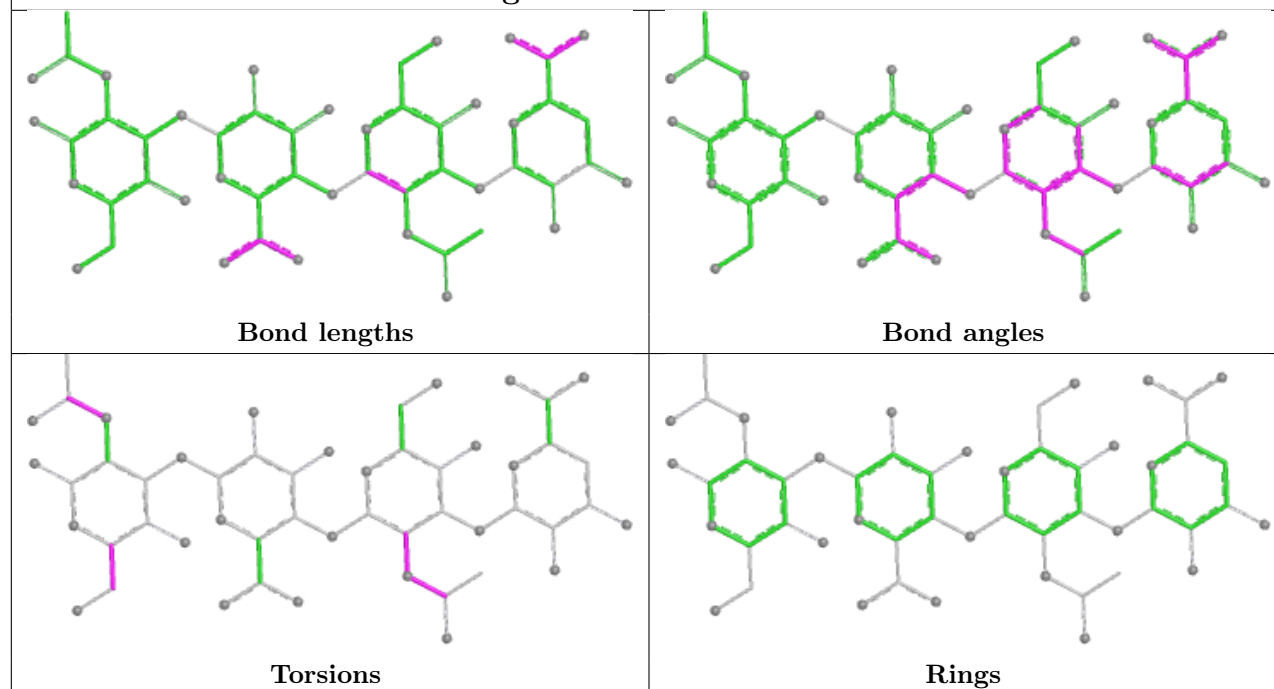
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



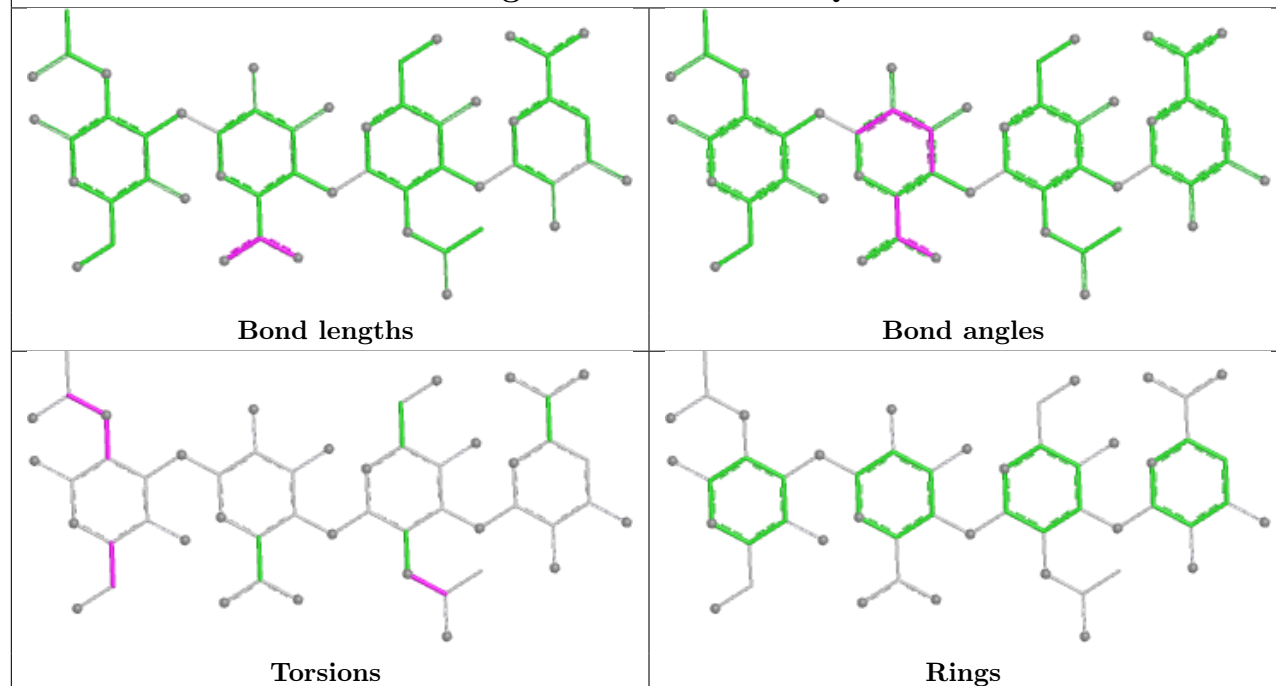
Oligosaccharide Chain N



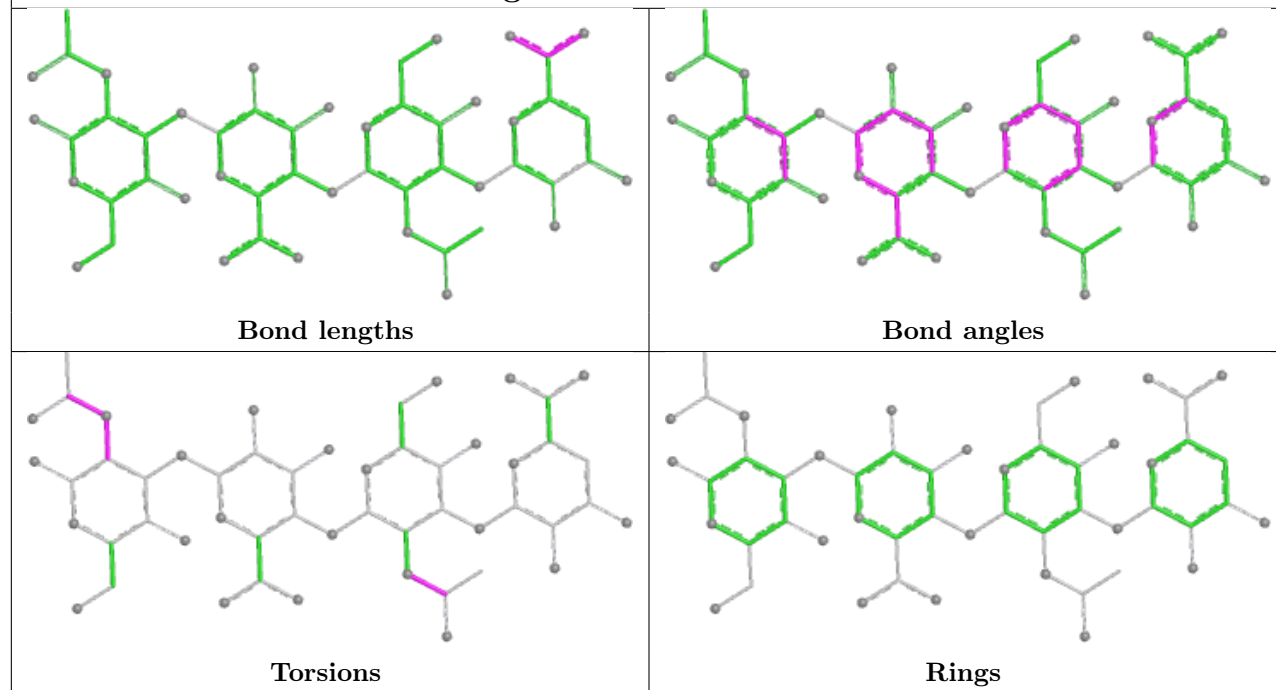
Oligosaccharide Chain O



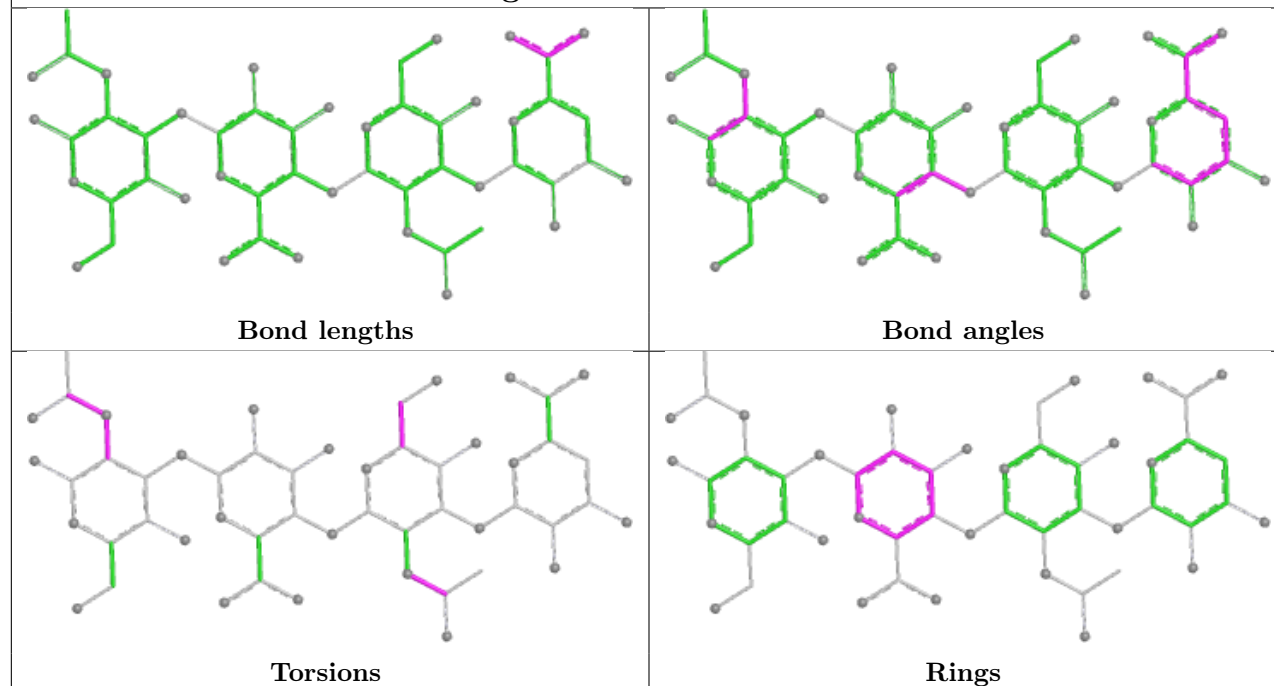
Oligosaccharide Chain Q



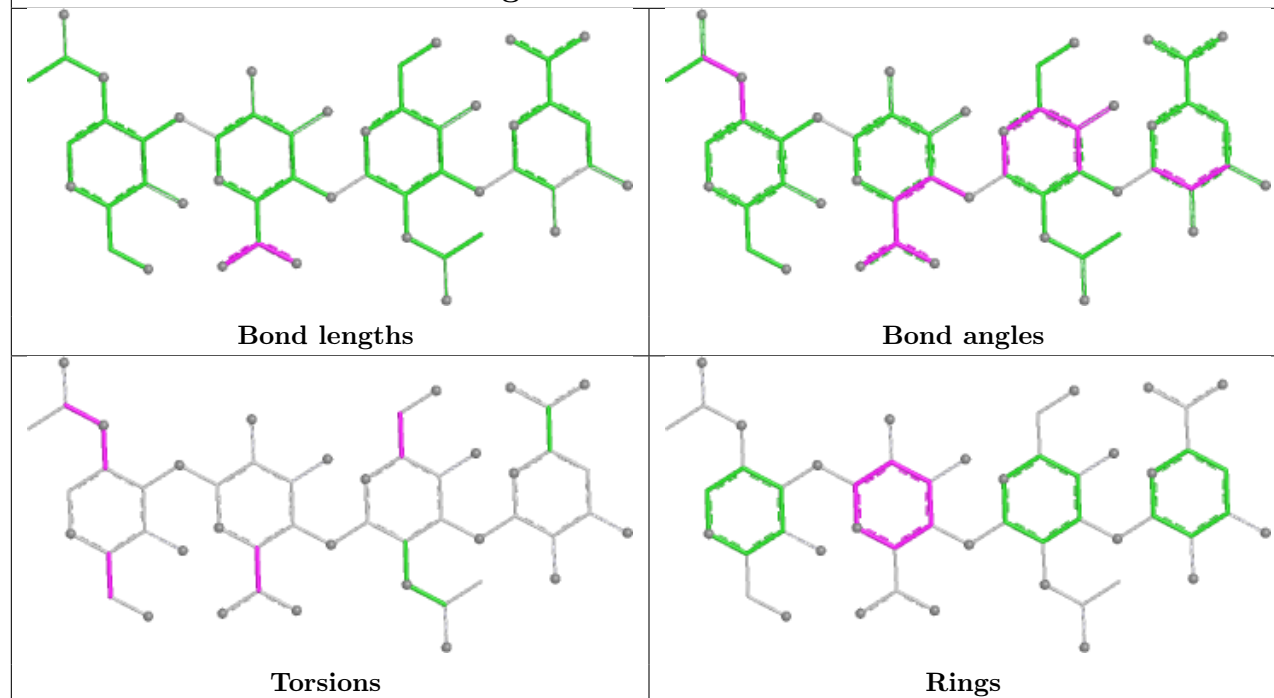
Oligosaccharide Chain T

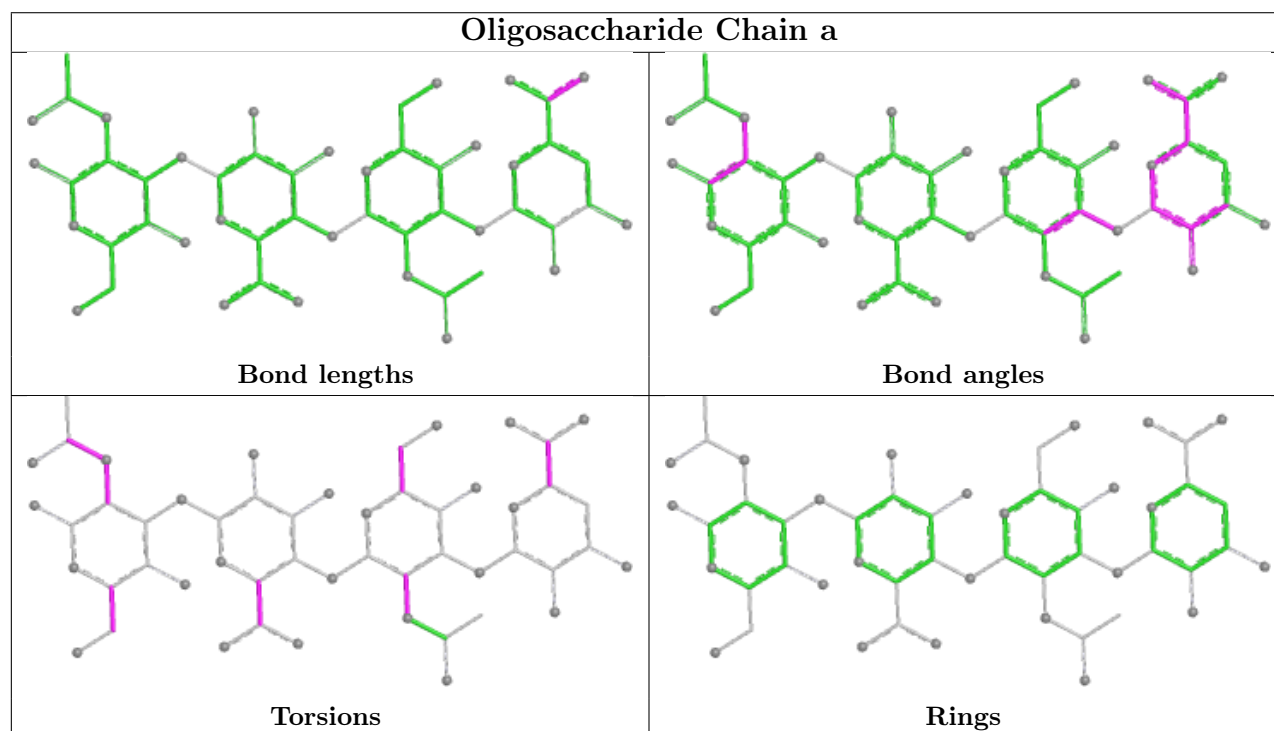
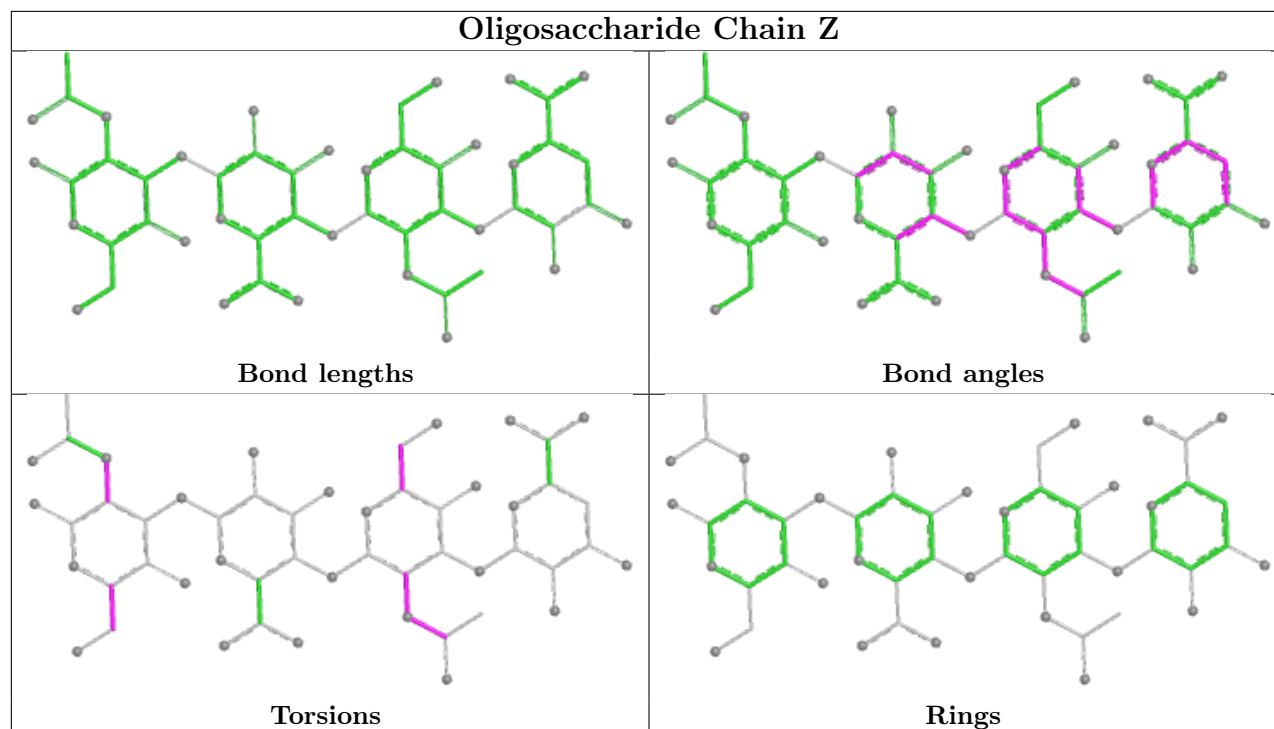


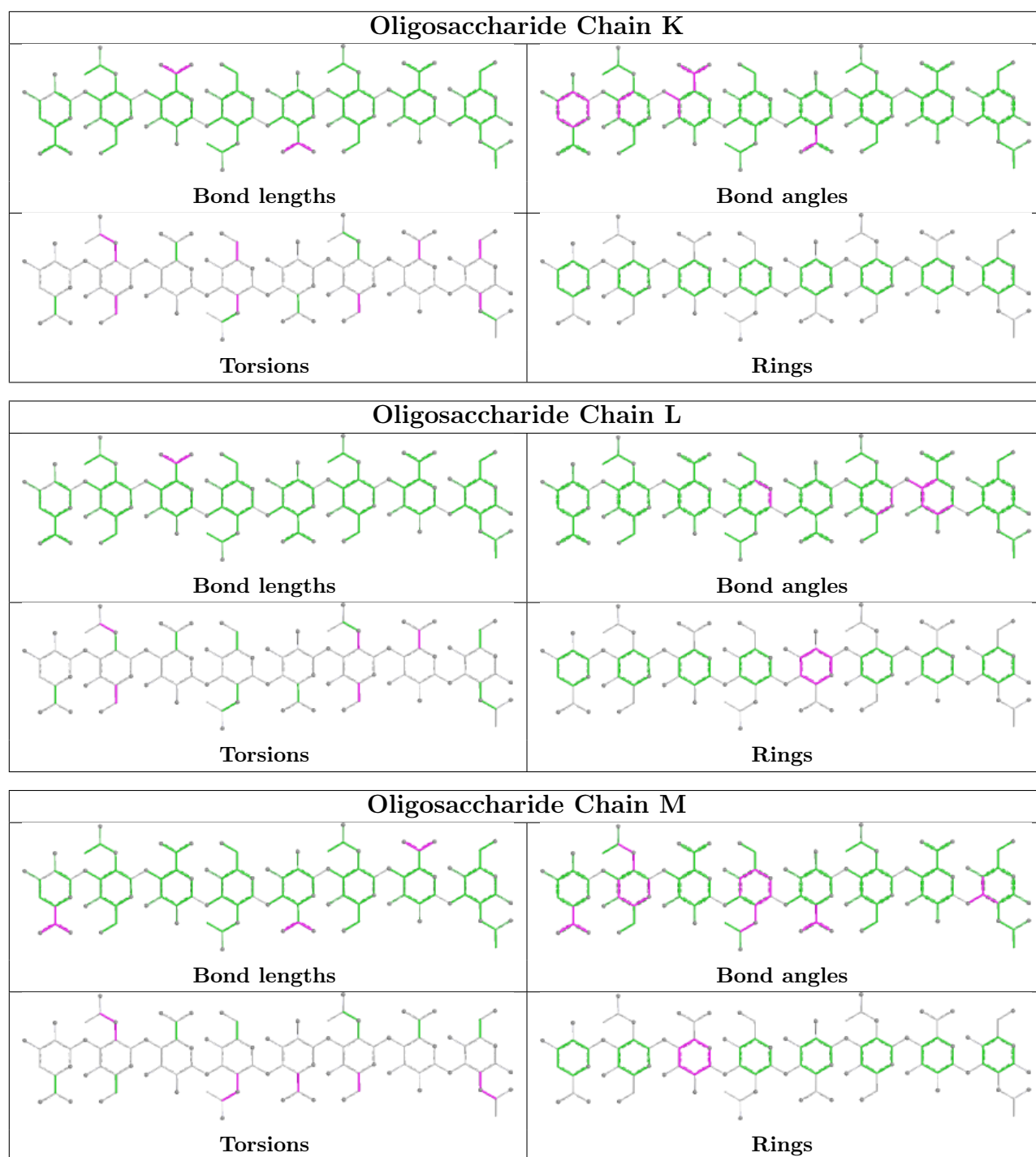
Oligosaccharide Chain U

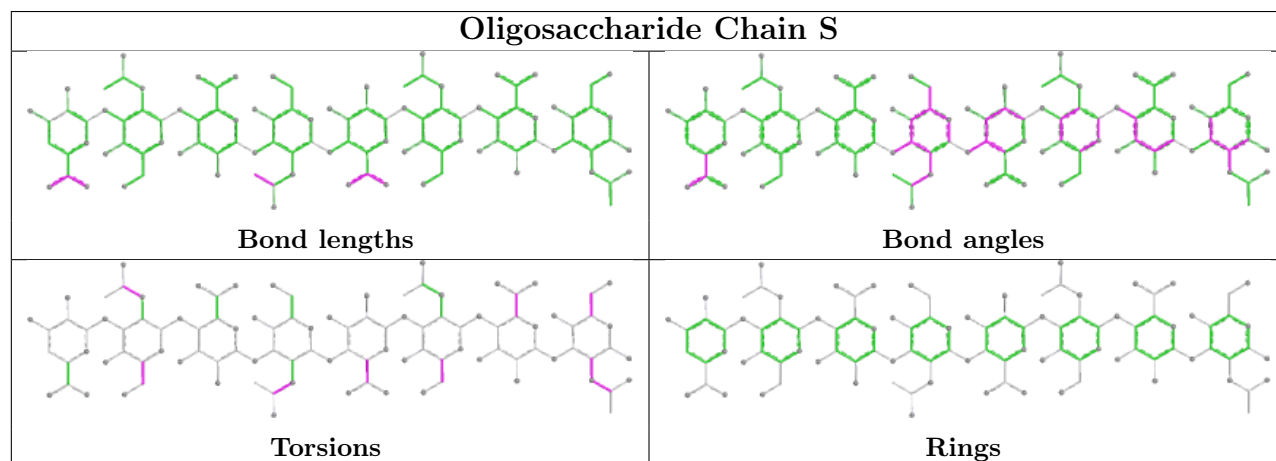
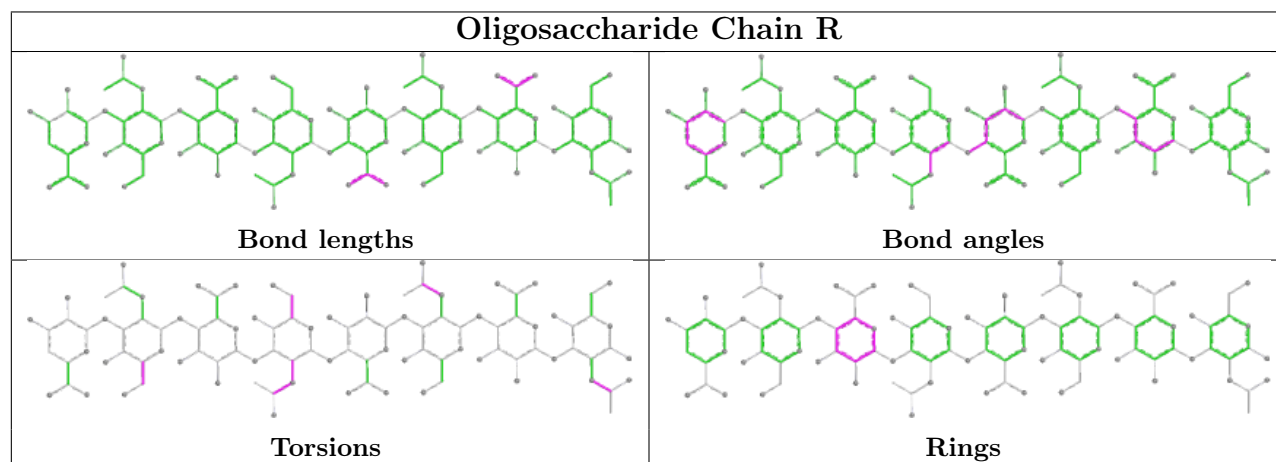
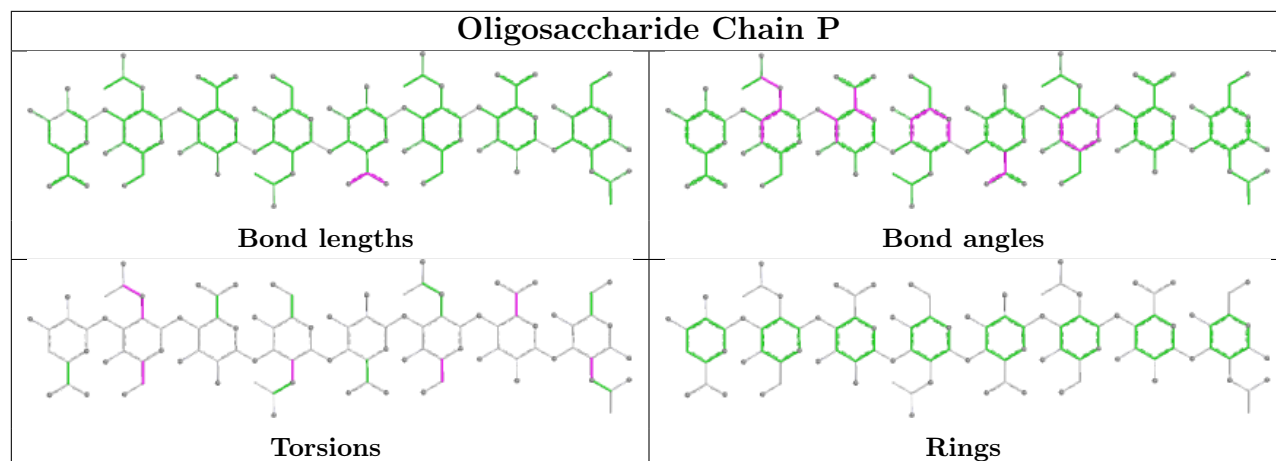


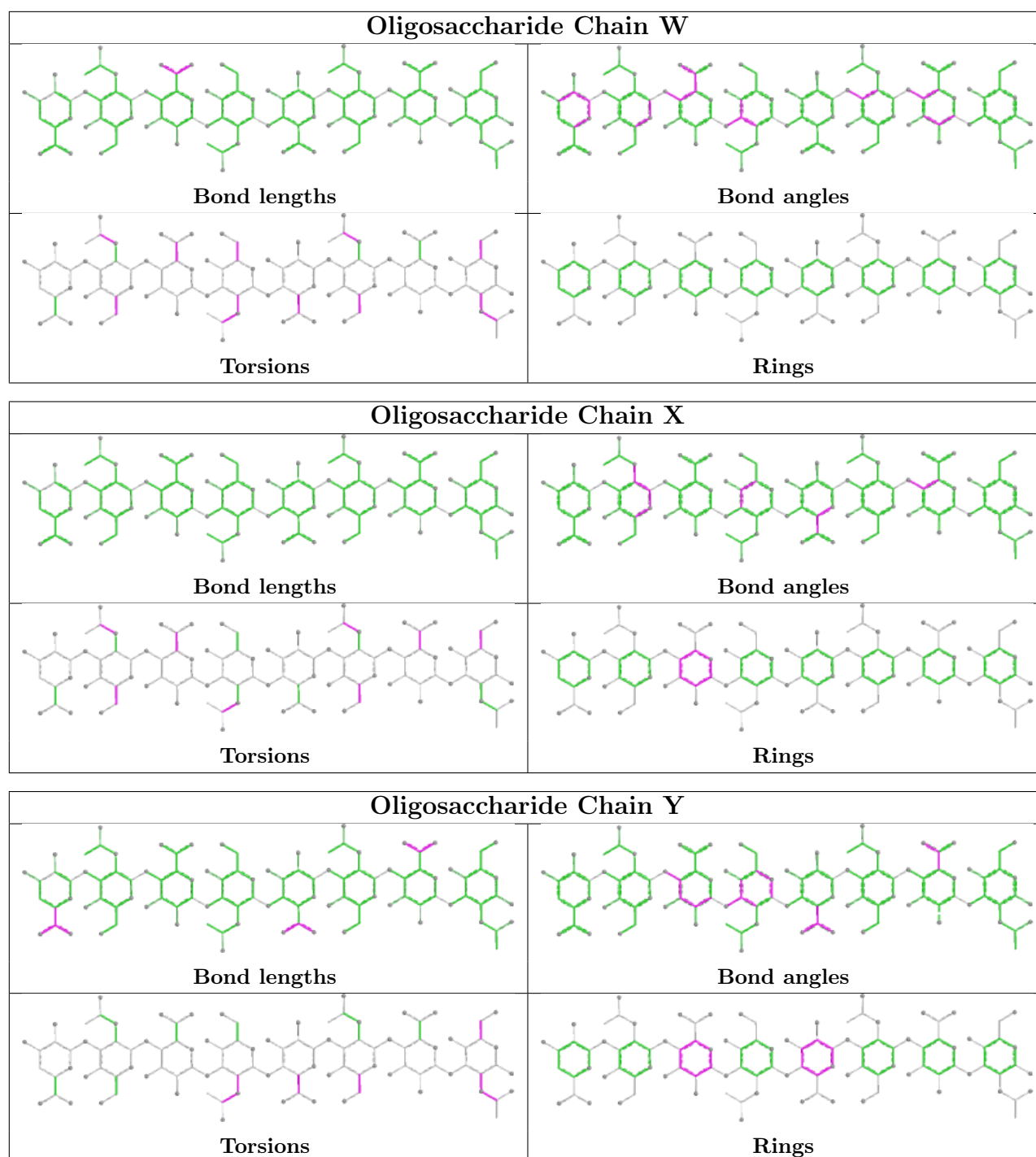
Oligosaccharide Chain V











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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	361/379 (95%)	-0.22	0 100 100	58, 100, 156, 180	0
1	B	361/379 (95%)	-0.23	2 (0%) 85 63	44, 98, 154, 201	0
1	C	361/379 (95%)	-0.20	0 100 100	50, 98, 157, 223	0
1	D	349/379 (92%)	-0.16	1 (0%) 90 74	57, 110, 164, 216	0
1	E	357/379 (94%)	-0.14	1 (0%) 90 74	40, 113, 165, 216	0
1	F	366/379 (96%)	-0.10	1 (0%) 90 74	41, 113, 175, 229	0
1	G	363/379 (95%)	-0.24	0 100 100	53, 103, 152, 205	0
1	H	361/379 (95%)	-0.25	0 100 100	31, 100, 155, 199	0
1	I	362/379 (95%)	-0.23	0 100 100	33, 101, 150, 207	0
All	All	3241/3411 (95%)	-0.20	5 (0%) 91 79	31, 104, 161, 229	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	351	ALA	3.7
1	B	256	GLY	2.8
1	D	42	PHE	2.3
1	B	33	PHE	2.3
1	E	33	PHE	2.2

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

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

6.3 Carbohydrates ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	N	1	15/15	0.45	0.10	116,119,120,122	0
2	GCD	V	4	11/12	0.46	0.14	103,103,104,105	0
3	GCD	X	8	11/12	0.46	0.11	105,107,107,107	0
2	GCD	N	4	11/12	0.48	0.14	92,100,102,103	0
2	GCD	U	4	11/12	0.50	0.09	106,106,107,107	0
2	GCD	O	4	11/12	0.51	0.10	110,110,111,112	0
2	NAG	T	1	15/15	0.53	0.12	91,94,96,96	0
3	GCD	W	8	11/12	0.54	0.09	97,98,100,101	0
2	NAG	O	1	15/15	0.57	0.12	104,105,107,109	0
3	NAG	Y	5	14/15	0.57	0.14	83,91,94,97	0
2	NAG	U	3	14/15	0.59	0.10	107,108,109,110	0
3	GCD	R	8	11/12	0.60	0.10	108,110,111,111	0
2	GCD	T	4	11/12	0.60	0.10	95,97,98,99	0
2	NAG	T	3	14/15	0.61	0.15	92,96,97,97	0
3	NAG	K	7	14/15	0.62	0.11	97,98,99,99	0
3	BDP	L	2	12/13	0.63	0.10	85,86,88,88	0
3	BDP	R	2	12/13	0.63	0.11	96,97,98,99	0
3	BDP	R	4	12/13	0.63	0.12	99,102,105,105	0
3	BDP	R	6	12/13	0.63	0.09	108,110,110,110	0
3	GCD	Y	8	11/12	0.63	0.08	104,104,105,105	0
3	NAG	R	5	14/15	0.64	0.12	103,106,108,109	0
2	NAG	J	1	14/15	0.64	0.11	84,85,87,87	0
3	NAG	X	5	14/15	0.64	0.12	96,100,101,102	0
3	GCD	S	8	11/12	0.65	0.09	99,100,102,102	0
3	NAG	M	7	14/15	0.65	0.09	100,101,102,104	0
3	NAG	M	5	14/15	0.67	0.13	91,94,96,98	0
3	BDP	X	4	12/13	0.67	0.09	96,99,101,101	0
2	GCD	Z	4	11/12	0.67	0.12	103,103,105,105	0
2	BDP	N	2	12/13	0.68	0.10	120,124,124,125	0
2	NAG	Z	1	15/15	0.68	0.09	107,108,108,109	0
3	BDP	X	6	12/13	0.68	0.10	104,105,106,107	0
3	BDP	M	6	12/13	0.69	0.09	100,101,101,102	0
2	NAG	a	1	15/15	-	-	94,96,96,97	0
2	BDP	a	2	12/13	-	-	94,96,97,97	0
2	NAG	a	3	14/15	-	-	93,94,95,95	0
2	GCD	a	4	11/12	-	-	91,93,95,95	0
3	BDP	M	2	12/13	0.69	0.12	90,92,94,94	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	BDP	P	4	12/13	0.69	0.09	91,92,92,93	0
2	GCD	Q	4	11/12	0.69	0.13	86,86,87,87	0
3	NAG	R	1	15/15	0.70	0.12	97,98,100,101	0
3	GCD	K	8	11/12	0.70	0.08	95,97,98,98	0
3	NAG	X	7	14/15	0.70	0.10	102,105,106,107	0
3	NAG	R	3	14/15	0.70	0.15	93,95,96,98	0
3	GCD	M	8	11/12	0.70	0.07	105,105,106,106	0
3	NAG	M	3	14/15	0.70	0.12	90,92,93,93	0
2	BDP	U	2	12/13	0.71	0.09	112,113,113,114	0
3	NAG	R	7	14/15	0.71	0.09	110,111,112,112	0
3	NAG	L	3	14/15	0.72	0.12	85,87,88,89	0
3	NAG	L	5	14/15	0.72	0.14	85,88,91,92	0
3	BDP	K	4	12/13	0.73	0.09	91,93,93,93	0
2	BDP	T	2	12/13	0.74	0.09	96,96,97,97	0
3	BDP	M	4	12/13	0.74	0.08	93,95,95,95	0
3	NAG	L	7	14/15	0.74	0.08	91,94,95,95	0
2	NAG	N	3	14/15	0.74	0.10	109,117,119,120	0
3	NAG	P	7	14/15	0.75	0.09	95,96,97,97	0
3	BDP	W	4	12/13	0.76	0.11	82,86,88,88	0
2	BDP	V	2	12/13	0.76	0.12	104,105,106,106	0
3	BDP	K	6	12/13	0.76	0.08	95,97,97,98	0
2	NAG	V	1	14/15	0.77	0.10	101,103,104,105	0
3	NAG	S	5	14/15	0.77	0.14	84,87,90,90	0
3	GCD	L	8	11/12	0.77	0.07	95,96,97,97	0
2	NAG	V	3	14/15	0.77	0.12	96,103,105,105	0
3	BDP	P	2	12/13	0.78	0.08	83,87,91,91	0
3	NAG	W	7	14/15	0.78	0.09	95,96,97,98	0
2	NAG	U	1	15/15	0.79	0.08	107,113,113,114	0
3	NAG	W	3	14/15	0.79	0.13	77,85,86,87	0
3	NAG	M	1	15/15	0.80	0.10	76,89,91,91	0
2	NAG	O	3	14/15	0.80	0.11	112,114,115,115	0
3	BDP	Y	6	12/13	0.81	0.09	100,102,103,103	0
3	NAG	Y	7	14/15	0.81	0.08	102,104,105,106	0
3	BDP	W	6	12/13	0.81	0.10	92,94,95,95	0
3	NAG	X	3	14/15	0.82	0.09	96,97,98,98	0
3	BDP	Y	4	12/13	0.82	0.10	86,89,92,92	0
3	BDP	L	4	12/13	0.82	0.07	87,88,89,90	0
2	NAG	Q	1	15/15	0.82	0.07	86,86,90,91	0
3	BDP	W	2	12/13	0.82	0.10	85,86,91,91	0
3	BDP	L	6	12/13	0.82	0.07	88,90,91,92	0
3	BDP	P	6	12/13	0.83	0.08	95,96,97,97	0
3	NAG	S	7	14/15	0.83	0.07	94,95,96,98	0

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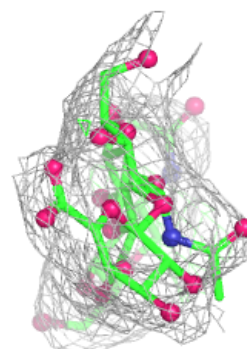
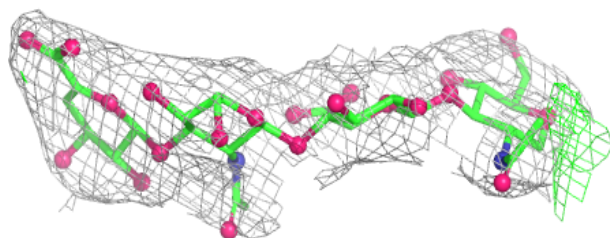
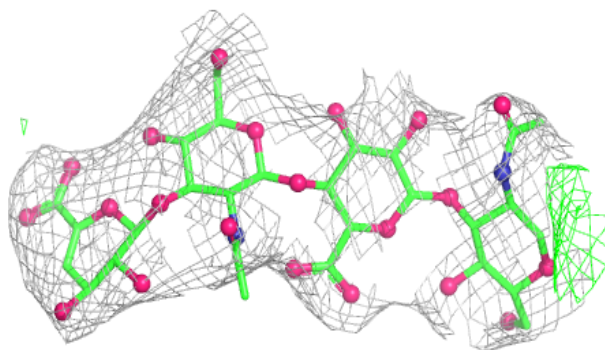
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	BDP	Y	2	12/13	0.83	0.09	81,82,86,87	0
2	BDP	Z	2	12/13	0.83	0.08	108,109,109,110	0
2	GCD	J	4	11/12	0.83	0.07	76,78,80,81	0
3	BDP	K	2	12/13	0.83	0.06	92,93,93,93	0
2	NAG	Q	3	14/15	0.83	0.11	83,86,87,88	0
3	NAG	W	5	14/15	0.83	0.11	87,89,90,91	0
3	NAG	Y	3	14/15	0.84	0.11	79,80,82,85	0
3	NAG	S	3	14/15	0.84	0.10	77,77,78,79	0
3	NAG	P	1	15/15	0.84	0.09	90,90,94,95	0
2	NAG	Z	3	14/15	0.85	0.10	102,106,107,108	0
3	NAG	X	1	15/15	0.85	0.07	97,98,98,99	0
3	GCD	P	8	11/12	0.85	0.05	92,93,94,94	0
2	BDP	Q	2	12/13	0.85	0.08	85,86,87,87	0
3	BDP	S	6	12/13	0.85	0.06	91,92,93,94	0
3	NAG	S	1	15/15	0.86	0.09	79,80,82,83	0
2	BDP	O	2	12/13	0.86	0.09	111,113,113,114	0
3	BDP	X	2	12/13	0.86	0.07	97,98,98,98	0
3	NAG	K	1	15/15	0.86	0.08	93,94,97,97	0
3	NAG	L	1	15/15	0.87	0.08	80,81,83,84	0
3	NAG	W	1	15/15	0.87	0.09	80,83,85,85	0
3	BDP	S	4	12/13	0.88	0.08	77,79,82,83	0
3	NAG	K	3	14/15	0.88	0.10	86,90,91,91	0
3	BDP	S	2	12/13	0.88	0.07	75,78,82,82	0
3	NAG	P	3	14/15	0.88	0.08	85,87,88,90	0
3	NAG	P	5	14/15	0.89	0.10	93,94,95,95	0
3	NAG	Y	1	15/15	0.90	0.08	83,84,86,86	0
2	NAG	J	3	14/15	0.90	0.08	68,74,76,76	0
2	BDP	J	2	12/13	0.91	0.06	77,82,83,83	0
3	NAG	K	5	14/15	0.92	0.06	92,93,94,95	0

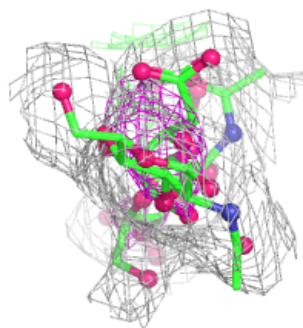
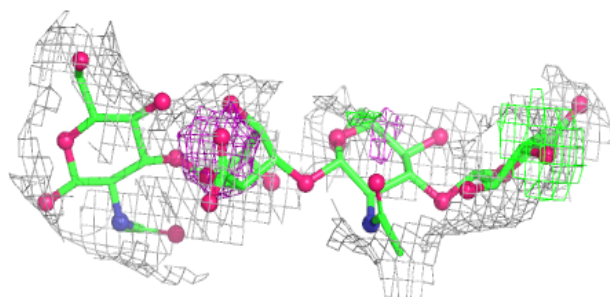
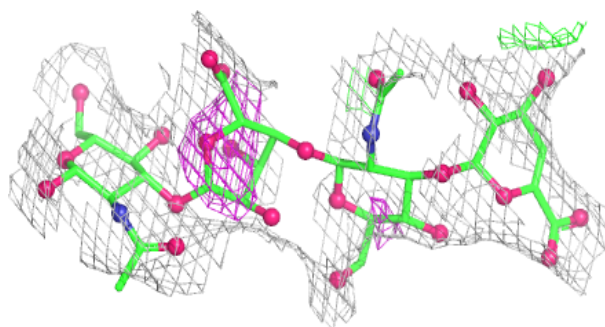
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain J:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

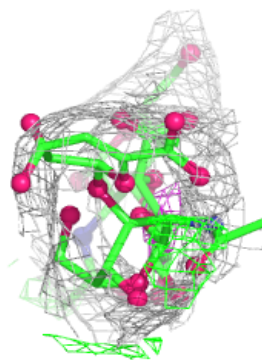
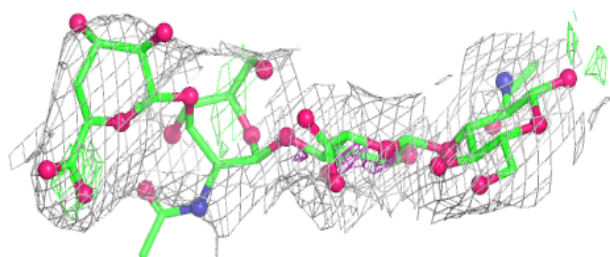
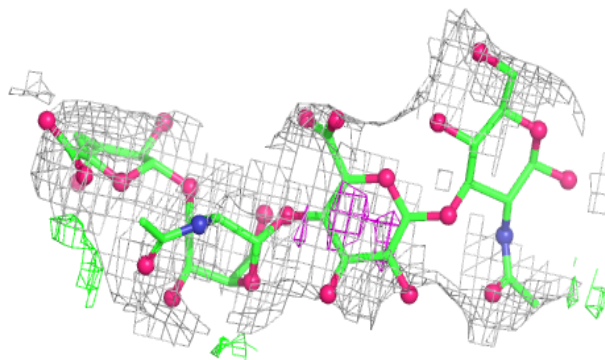
**Electron density around Chain N:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

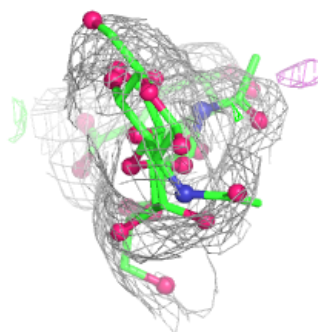
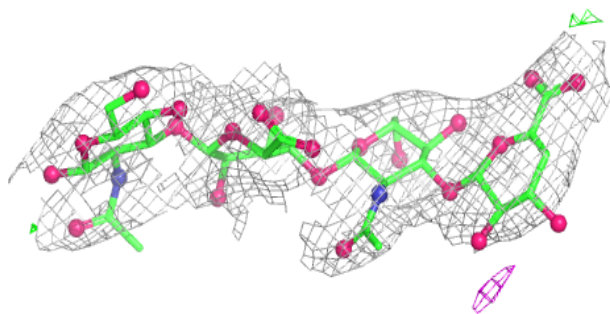
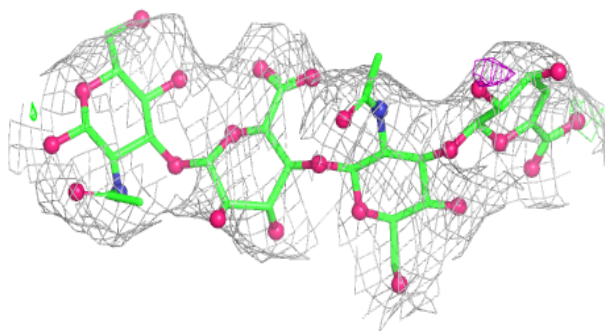


Electron density around Chain O:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

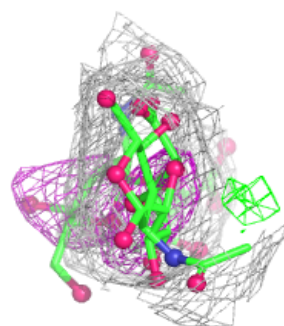
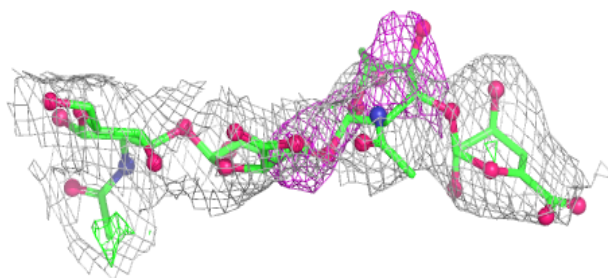
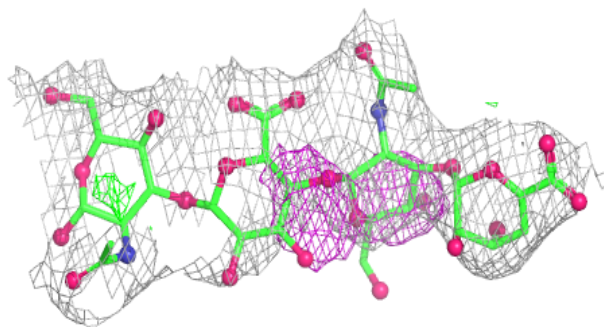
**Electron density around Chain Q:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

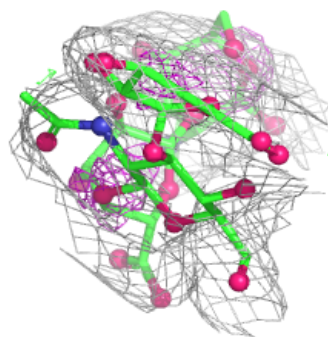
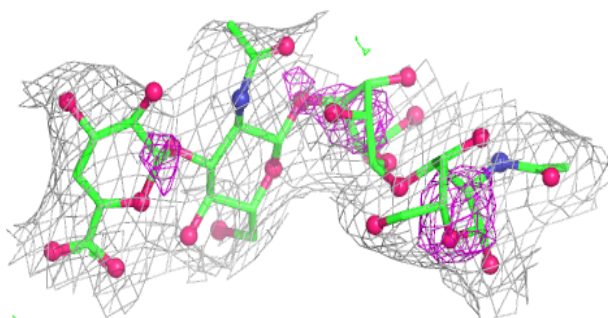
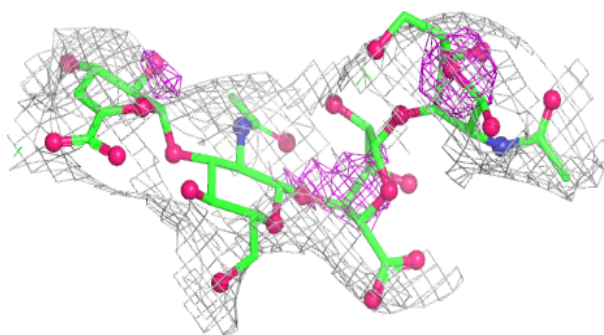


Electron density around Chain T:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

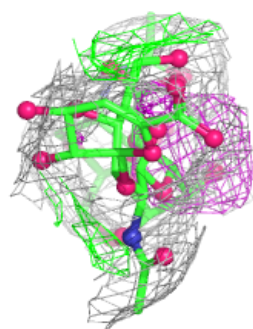
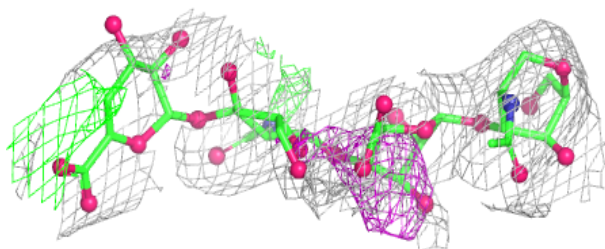
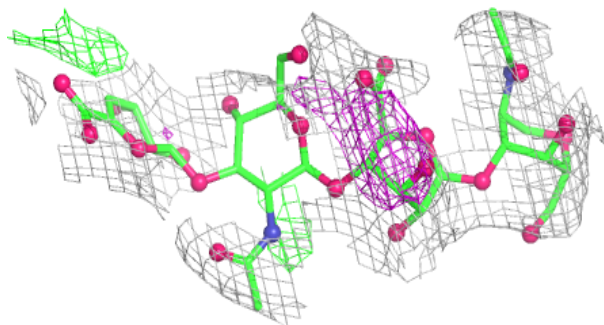
**Electron density around Chain U:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

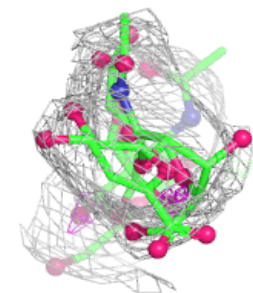
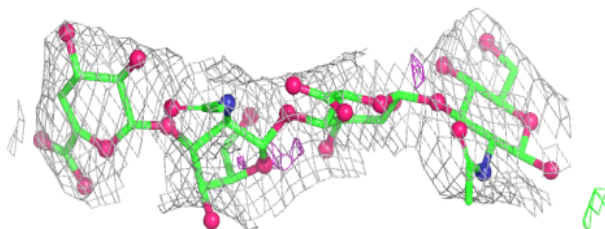
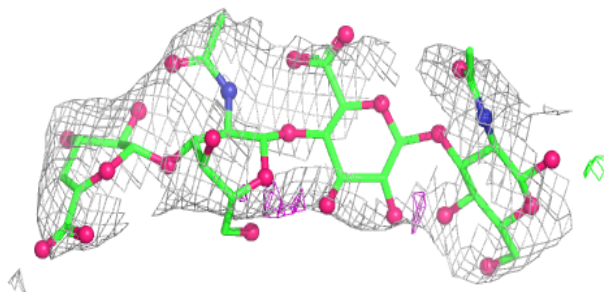


Electron density around Chain V:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

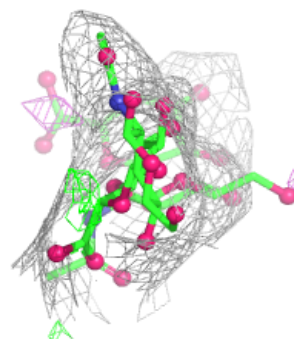
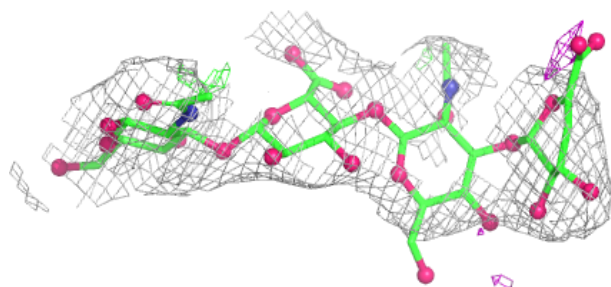
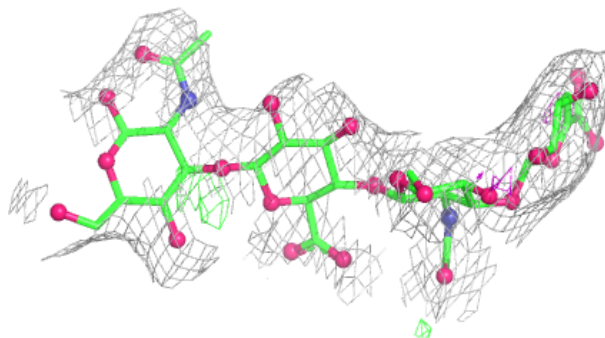
**Electron density around Chain Z:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

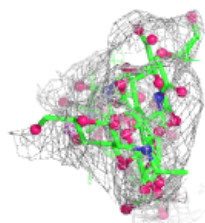
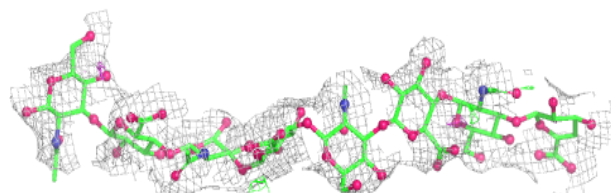
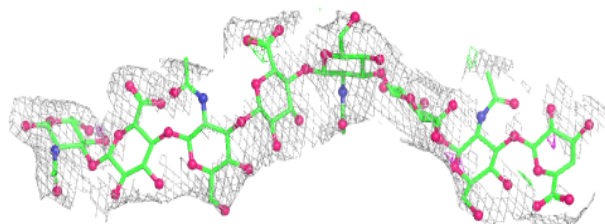


Electron density around Chain a:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

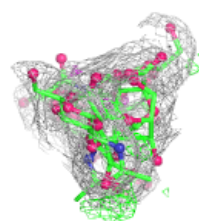
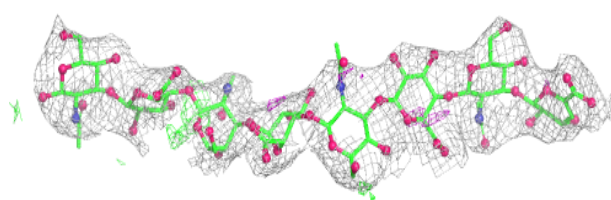
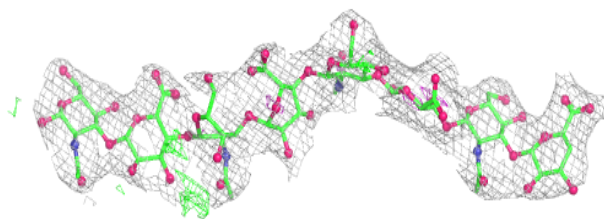
**Electron density around Chain K:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

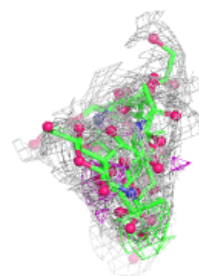
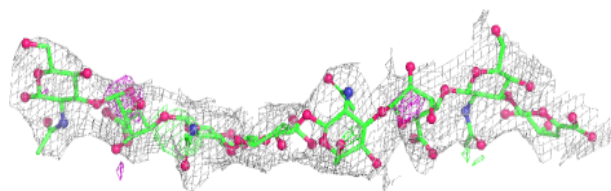
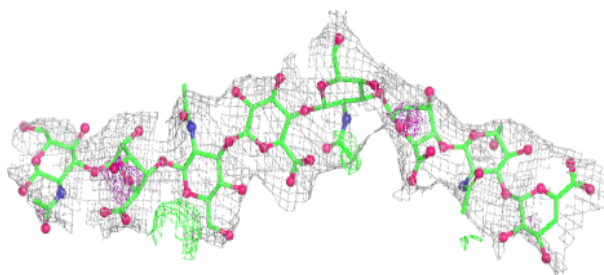


Electron density around Chain L:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

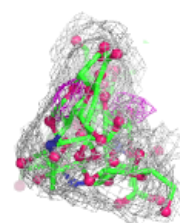
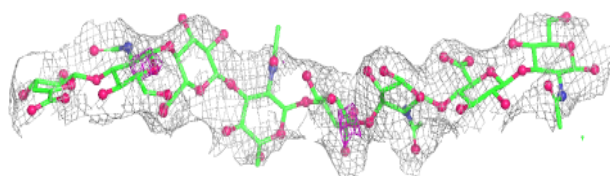
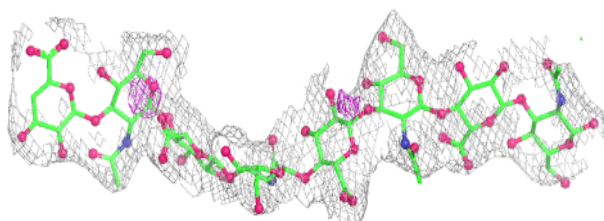
**Electron density around Chain M:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

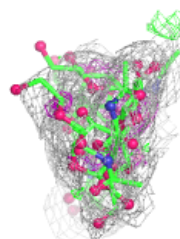
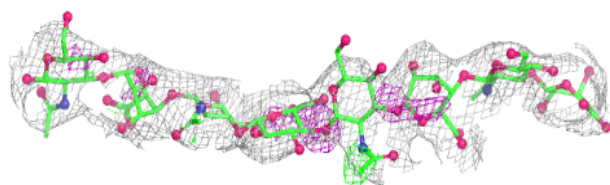
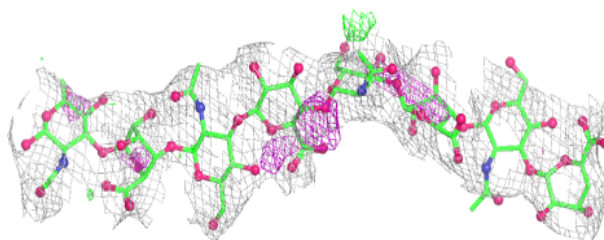


Electron density around Chain P:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

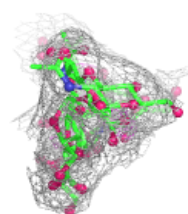
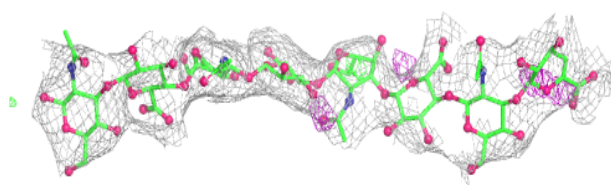
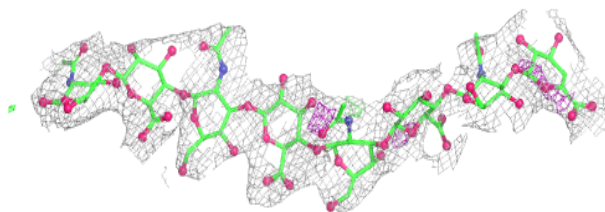
**Electron density around Chain R:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

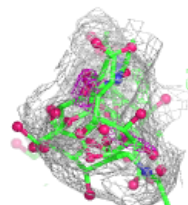
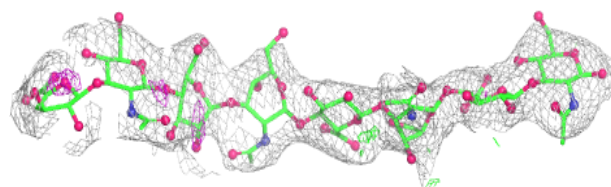
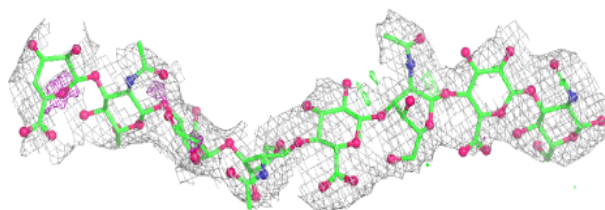


Electron density around Chain S:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

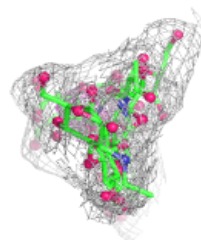
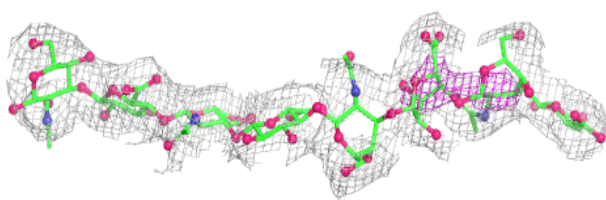
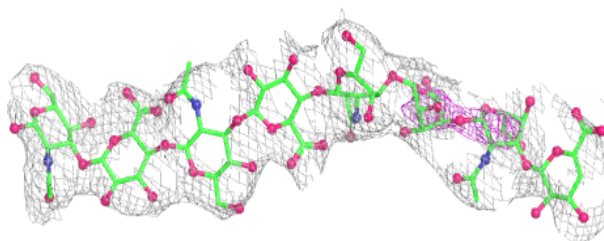
**Electron density around Chain W:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

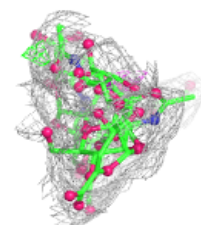
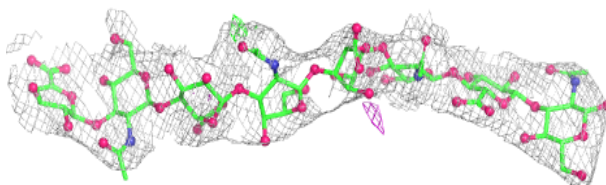
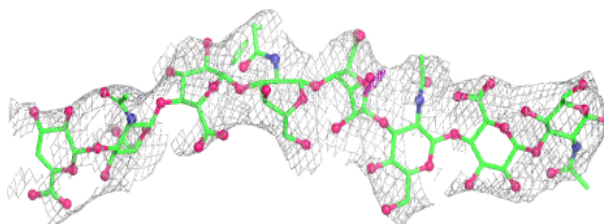


Electron density around Chain X:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around Chain Y:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.