



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:33 AM UTC

PDB ID : 5CJX / pdb_00005cjx
Title : Crystal structure of 8ANC195 Fab in complex with BG505 SOSIP.664 HIV-1 Env trimer
Authors : Scharf, L.; Bjorkman, P.J.
Deposited on : 2015-07-15
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:

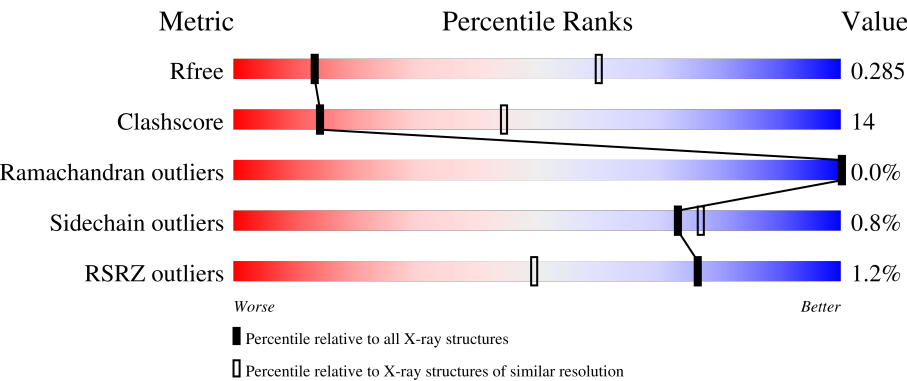
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

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.



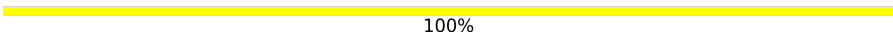
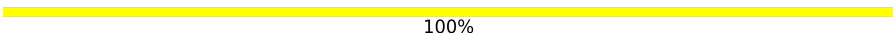
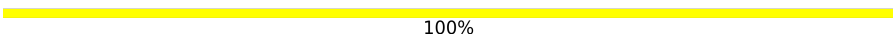
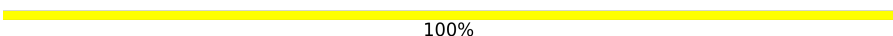
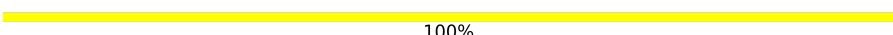
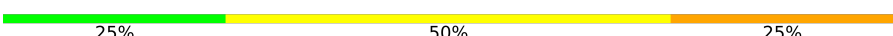
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1521 (3.66-3.50)
Clashscore	190562	1595 (3.66-3.50)
Ramachandran outliers	187476	1551 (3.66-3.50)
Sidechain outliers	187428	1551 (3.66-3.50)
RSRZ outliers	180081	1520 (3.66-3.50)

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	244	
1	D	244	
1	H	244	
2	B	153	
2	J	153	

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Mol	Chain	Length	Quality of chain
2	X	153	
3	C	215	
3	E	215	
3	L	215	
4	G	479	
4	K	479	
4	Y	479	
5	F	5	
5	b	5	
6	I	2	
6	M	2	
6	P	2	
6	Q	2	
6	V	2	
7	N	10	
7	S	10	
7	Z	10	
8	O	6	
8	U	6	
9	R	4	
9	W	4	
10	T	3	
10	a	3	

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
7	MAN	Z	9	-	-	X	-
8	NAG	O	1	-	-	X	-

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 22299 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 8ANC195 G52K5 heavy chain, IG gamma-1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	224	Total	C	N	O	S	0	0	0
			1655	1049	278	323	5			
1	D	224	Total	C	N	O	S	0	0	0
			1614	1022	269	318	5			
1	H	227	Total	C	N	O	S	0	0	0
			1675	1063	282	325	5			

- Molecule 2 is a protein called BG505 Env gp41.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	124	Total	C	N	O	S	0	0	0
			956	602	164	184	6			
2	J	124	Total	C	N	O	S	0	0	0
			957	605	163	183	6			
2	X	121	Total	C	N	O	S	0	0	0
			945	596	165	178	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	559	PRO	ILE	engineered mutation	UNP Q2N0S6
B	605	CYS	THR	engineered mutation	UNP Q2N0S6
J	559	PRO	ILE	engineered mutation	UNP Q2N0S6
J	605	CYS	THR	engineered mutation	UNP Q2N0S6
X	559	PRO	ILE	engineered mutation	UNP Q2N0S6
X	605	CYS	THR	engineered mutation	UNP Q2N0S6

- Molecule 3 is a protein called 8ANC195 G52K5 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	214	Total	C	N	O	S	0	0	0
			1593	996	267	325	5			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	214	Total	C	N	O	S	0	0	0
			1560	976	260	319	5			
3	L	214	Total	C	N	O	S	0	0	0
			1540	962	255	318	5			

- Molecule 4 is a protein called BG505 Env gp120.

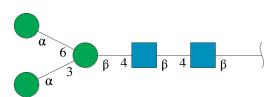
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	398	Total	C	N	O	S	0	0	0
			2857	1807	480	546	24			
4	K	405	Total	C	N	O	S	0	0	0
			2984	1891	502	564	27			
4	Y	397	Total	C	N	O	S	0	0	0
			2947	1864	501	556	26			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	320M	ASN	THR	engineered mutation	UNP Q2N0S6
G	501	CYS	ALA	engineered mutation	UNP Q2N0S6
G	509	ARG	-	expression tag	UNP Q2N0S6
G	510	ARG	-	expression tag	UNP Q2N0S6
G	511	ARG	-	expression tag	UNP Q2N0S6
G	512	ARG	-	expression tag	UNP Q2N0S6
G	513	ARG	-	expression tag	UNP Q2N0S6
K	332	ASN	THR	engineered mutation	UNP Q2N0S6
K	501	CYS	ALA	engineered mutation	UNP Q2N0S6
K	509	ARG	-	expression tag	UNP Q2N0S6
K	510	ARG	-	expression tag	UNP Q2N0S6
K	511	ARG	-	expression tag	UNP Q2N0S6
K	512	ARG	-	expression tag	UNP Q2N0S6
K	513	ARG	-	expression tag	UNP Q2N0S6
Y	320M	ASN	THR	engineered mutation	UNP Q2N0S6
Y	501	CYS	ALA	engineered mutation	UNP Q2N0S6
Y	509	ARG	-	expression tag	UNP Q2N0S6
Y	510	ARG	-	expression tag	UNP Q2N0S6
Y	511	ARG	-	expression tag	UNP Q2N0S6
Y	512	ARG	-	expression tag	UNP Q2N0S6
Y	513	ARG	-	expression tag	UNP Q2N0S6

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-a

cetamido-2-deoxy-beta-D-glucopyranose.



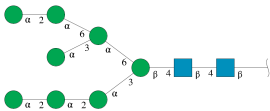
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	F	5	Total	C	N	O	0	0	0
			61	34	2	25			
5	b	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
6	M	2	Total	C	N	O	0	0	0
			28	16	2	10			
6	P	2	Total	C	N	O	0	0	0
			28	16	2	10			
6	Q	2	Total	C	N	O	0	0	0
			28	16	2	10			
6	V	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



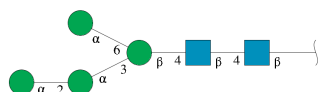
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	N	10	Total	C	N	O	0	0	0
			116	64	2	50			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	S	10	Total	C	N	O	0	0	0
			116	64	2	50			
7	Z	10	Total	C	N	O	0	0	0
			116	64	2	50			

- Molecule 8 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



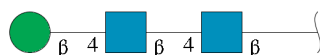
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
8	O	6	Total	C	N	O	0	0	0
			72	40	2	30			
8	U	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 9 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



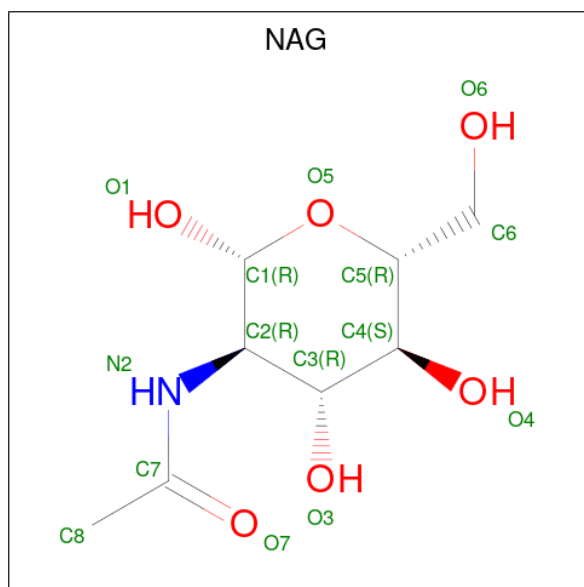
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	R	4	Total	C	N	O	0	0	0
			50	28	2	20			
9	W	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 10 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
10	T	3	Total	C	N	O	0	0	0
			39	22	2	15			
10	a	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 11 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

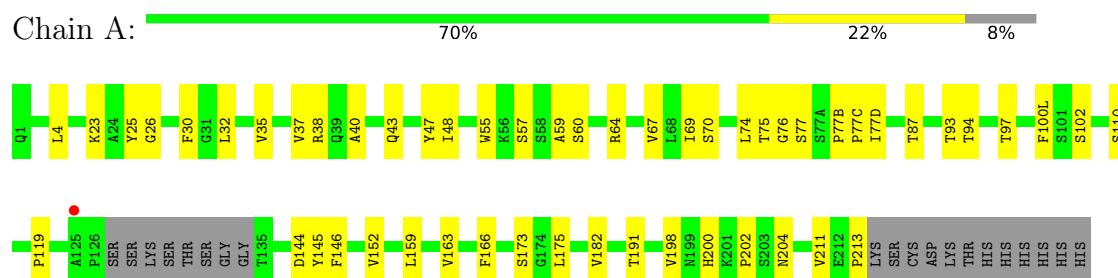


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	A	1	Total	C	N	O	0	0
			14	8	1	5		
11	B	1	Total	C	N	O	0	0
			14	8	1	5		
11	J	1	Total	C	N	O	0	0
			14	8	1	5		
11	K	1	Total	C	N	O	0	0
			14	8	1	5		
11	Y	1	Total	C	N	O	0	0
			14	8	1	5		
11	Y	1	Total	C	N	O	0	0
			14	8	1	5		

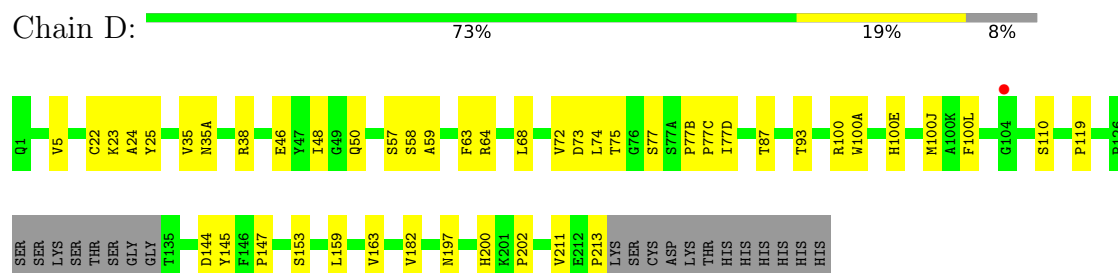
3 Residue-property plots [i](#)

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.

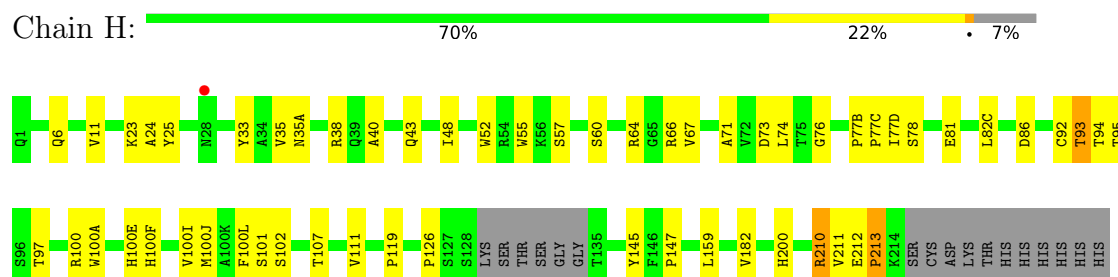
- Molecule 1: 8ANC195 G52K5 heavy chain, IG gamma-1 chain



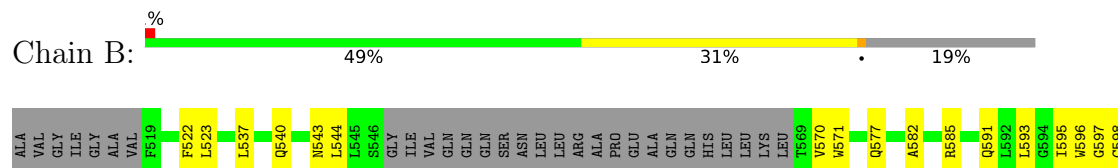
- Molecule 1: 8ANC195 G52K5 heavy chain, IG gamma-1 chain



- Molecule 1: 8ANC195 G52K5 heavy chain, IG gamma-1 chain

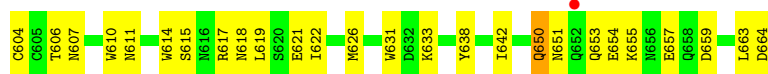
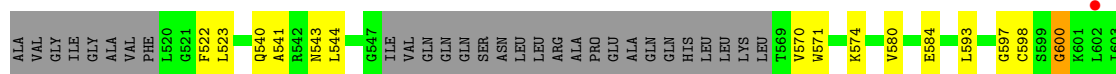


- Molecule 2: BG505 Env gp41

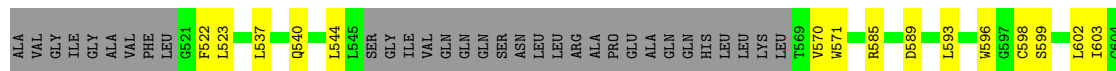




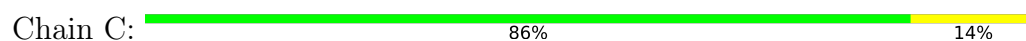
• Molecule 2: BG505 Env gp41



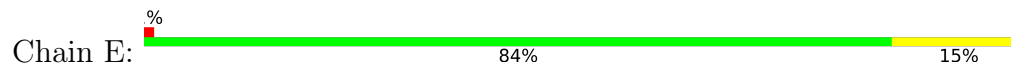
• Molecule 2: BG505 Env gp41



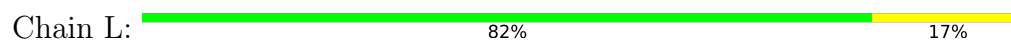
• Molecule 3: 8ANC195 G52K5 light chain

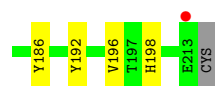


• Molecule 3: 8ANC195 G52K5 light chain

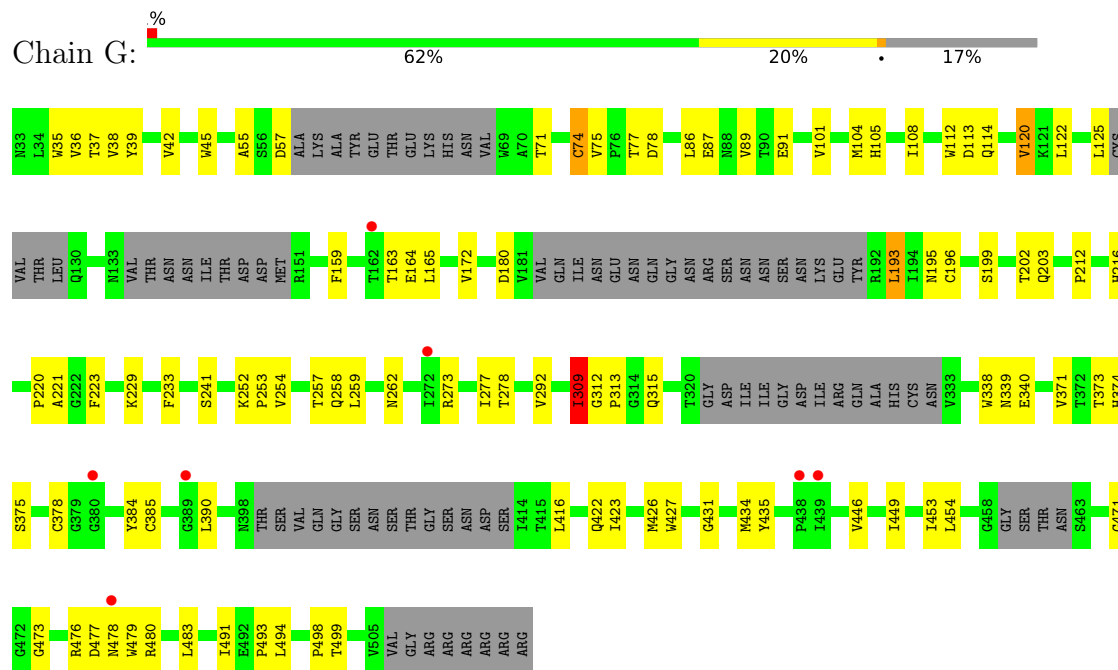


• Molecule 3: 8ANC195 G52K5 light chain

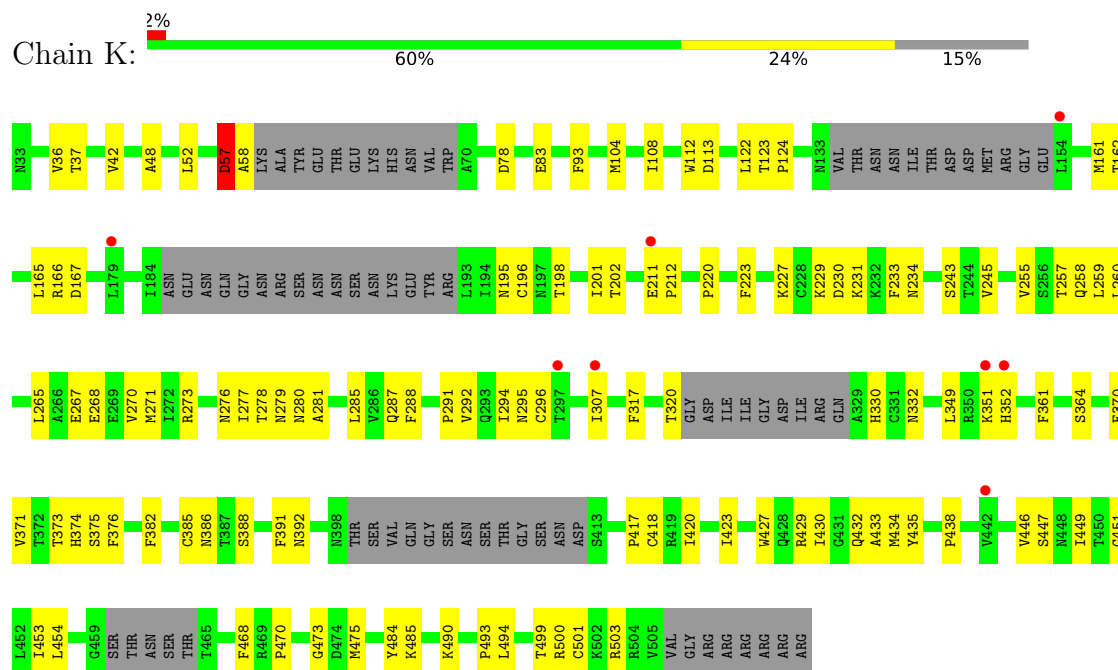




- Molecule 4: BG505 Env gp120

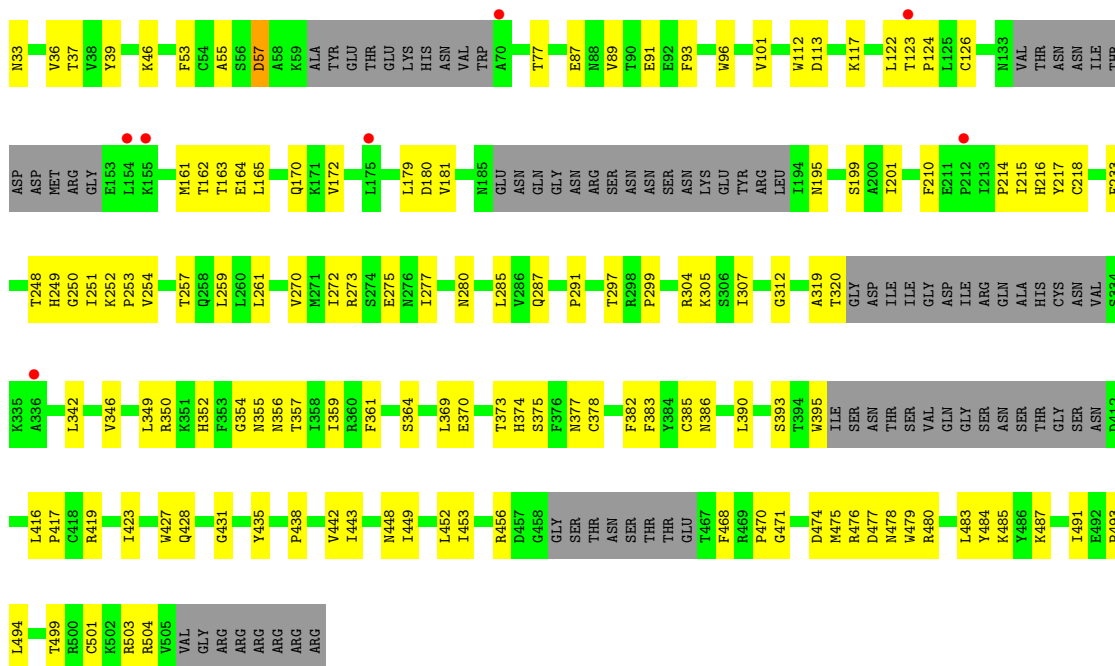


- Molecule 4: BG505 Env gp120

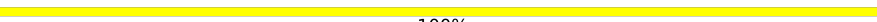


- Molecule 4: BG505 Env gp120



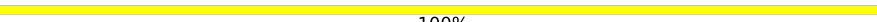


● Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%

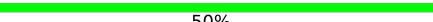


● Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain b:  100%

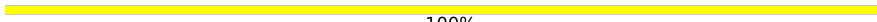


● Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  50%  50%



● Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  100%

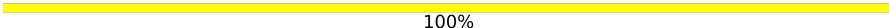
MAG1
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:  50% 50%

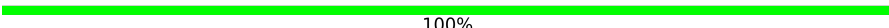
MAG1
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  100%

MAG1
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain V:  100%

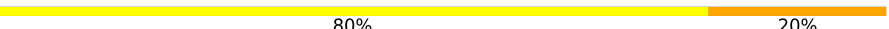
MAG1
MAG2

- Molecule 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  20% 50% 30%

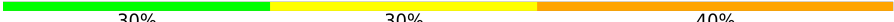
MAG1
MAG2
BNA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

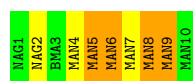
- Molecule 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S:  80% 20%

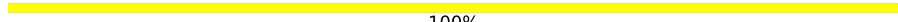
MAG1
MAG2
BNA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

- Molecule 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z:  30% 30% 40%



- Molecule 8: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  100%



- Molecule 8: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain U:  50% 50%

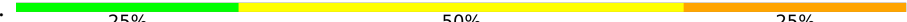


- Molecule 9: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain R:  25% 75%



- Molecule 9: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain W:  25% 50% 25%



- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:  67% 33%



- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain a:  67% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	117.74Å 195.22Å 119.09Å 90.00° 101.60° 90.00°	Depositor
Resolution (Å)	29.73 – 3.58 29.73 – 3.58	Depositor EDS
% Data completeness (in resolution range)	99.0 (29.73-3.58) 98.9 (29.73-3.58)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 3.55Å)	Xtriage
Refinement program	PHENIX (1.10pre_2100: ???)	Depositor
R, R_{free}	0.239 , 0.286 0.239 , 0.285	Depositor DCC
R_{free} test set	2015 reflections (3.29%)	wwPDB-VP
Wilson B-factor (Å ²)	130.2	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 104.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.029 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	22299	wwPDB-VP
Average B, all atoms (Å ²)	149.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.24	1/1700 (0.1%)	0.48	2/2334 (0.1%)
1	D	0.24	1/1656 (0.1%)	0.43	1/2280 (0.0%)
1	H	0.28	2/1720 (0.1%)	0.46	2/2359 (0.1%)
2	B	0.30	0/973	0.53	0/1327
2	J	0.26	0/974	0.52	2/1325 (0.2%)
2	X	0.26	0/962	0.46	1/1309 (0.1%)
3	C	0.14	0/1628	0.36	0/2221
3	E	0.15	0/1594	0.36	0/2181
3	L	0.14	0/1574	0.35	0/2159
4	G	0.28	1/2916 (0.0%)	0.40	0/4002
4	K	0.23	0/3047	0.49	1/4172 (0.0%)
4	Y	0.20	0/3009	0.45	1/4119 (0.0%)
All	All	0.23	5/21753 (0.0%)	0.44	10/29788 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
4	G	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	309	ILE	C-N	10.99	1.49	1.33
1	H	212	GLU	C-N	5.88	1.40	1.33
1	H	213	PRO	N-CD	5.30	1.55	1.47
1	D	213	PRO	N-CD	5.28	1.55	1.47
1	A	213	PRO	N-CD	5.10	1.54	1.47

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	191	THR	N-CA-C	9.60	121.51	111.14
1	H	212	GLU	CA-C-N	-6.70	113.39	120.03
1	H	212	GLU	C-N-CA	-6.70	113.39	120.03
1	A	144	ASP	N-CA-C	6.53	120.66	111.52
2	X	596	TRP	N-CA-C	-6.16	101.36	110.48
2	J	650	GLN	N-CA-C	5.84	120.70	112.93
4	Y	57	ASP	CB-CA-C	-5.73	109.95	116.54
2	J	600	GLY	N-CA-C	-5.55	104.87	111.36
4	K	57	ASP	CB-CA-C	-5.54	110.20	116.63
1	D	144	ASP	N-CA-C	5.29	118.93	111.52

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	G	309	ILE	Mainchain

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1655	0	1577	36	0
1	D	1614	0	1511	48	0
1	H	1675	0	1600	52	0
2	B	956	0	898	71	0
2	J	957	0	915	39	0
2	X	945	0	905	37	0
3	C	1593	0	1489	19	0
3	E	1560	0	1438	23	0
3	L	1540	0	1382	26	0
4	G	2857	0	2565	79	0
4	K	2984	0	2763	108	0
4	Y	2947	0	2734	123	0
5	F	61	0	52	4	0
5	b	61	0	52	2	0
6	I	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	M	28	0	25	2	0
6	P	28	0	25	0	0
6	Q	28	0	25	1	0
6	V	28	0	25	0	0
7	N	116	0	97	7	0
7	S	116	0	96	8	0
7	Z	116	0	96	13	0
8	O	72	0	61	9	0
8	U	72	0	61	5	0
9	R	50	0	43	1	0
9	W	50	0	43	3	0
10	T	39	0	34	1	0
10	a	39	0	34	1	0
11	A	14	0	13	0	0
11	B	14	0	13	0	0
11	J	14	0	13	0	0
11	K	14	0	13	0	0
11	Y	28	0	26	2	0
All	All	22299	0	20649	608	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 14.

All (608) 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)
2:B:645:LEU:O	2:B:649:SER:HB2	1.16	1.32
4:K:231:LYS:HD2	4:K:267:GLU:CD	1.61	1.26
2:B:645:LEU:O	2:B:649:SER:CB	1.88	1.19
4:G:278:THR:HG22	8:O:1:NAG:O6	1.40	1.15
2:B:650:GLN:O	2:B:654:GLU:N	1.77	1.15
2:B:598:CYS:SG	2:B:604:CYS:SG	1.26	1.15
4:Y:216:HIS:CE1	4:Y:250:GLY:HA3	1.81	1.14
2:B:597:GLY:N	2:B:651:ASN:HD21	1.42	1.14
2:J:604:CYS:O	4:K:37:THR:HG23	1.51	1.11
2:B:598:CYS:SG	2:B:604:CYS:CB	2.39	1.10
1:D:59:ALA:HB2	7:Z:9:MAN:H4	1.30	1.09
2:B:601:LYS:O	2:X:655:LYS:NZ	1.92	1.03
4:Y:257:THR:HG21	4:Y:370:GLU:O	1.59	1.02
4:Y:216:HIS:CE1	4:Y:250:GLY:CA	2.43	1.01
2:B:608:VAL:HG22	2:B:649:SER:OG	1.61	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:Y:257:THR:CG2	4:Y:370:GLU:O	2.09	1.00
4:K:386:ASN:OD1	4:K:388:SER:N	2.00	0.93
4:K:231:LYS:CD	4:K:267:GLU:CD	2.44	0.90
2:J:597:GLY:HA3	4:K:503:ARG:HH12	1.39	0.88
2:J:618:ASN:HB3	2:J:621:GLU:HB2	1.55	0.88
2:B:597:GLY:N	2:B:651:ASN:ND2	2.21	0.87
2:B:602:LEU:HB2	2:X:655:LYS:HZ3	1.39	0.87
1:D:59:ALA:HB2	7:Z:9:MAN:C4	2.05	0.86
4:Y:216:HIS:CD2	4:Y:250:GLY:N	2.44	0.86
1:D:38:ARG:HB2	1:D:48:ILE:HD11	1.58	0.85
4:Y:216:HIS:NE2	4:Y:250:GLY:N	2.26	0.83
2:B:597:GLY:CA	2:B:651:ASN:HD21	1.92	0.83
2:X:605:CYS:HA	4:Y:37:THR:HG22	1.61	0.81
4:K:259:LEU:HB2	4:K:374:HIS:CE1	2.15	0.81
1:D:59:ALA:CB	7:Z:9:MAN:H4	2.10	0.81
1:H:210:ARG:HH11	1:H:210:ARG:HG2	1.45	0.81
1:A:74:LEU:HD23	1:A:77(B):PRO:HA	1.63	0.80
2:B:601:LYS:CB	2:B:602:LEU:HD13	2.10	0.80
4:Y:249:HIS:CD2	4:Y:251:ILE:HG12	2.16	0.80
4:G:278:THR:OG1	1:H:74:LEU:HD11	1.82	0.80
8:U:3:BMA:HO4	8:U:5:MAN:HO6	1.30	0.80
1:H:11:VAL:HG21	1:H:147:PRO:HG3	1.64	0.79
2:B:544:LEU:HD21	4:G:493:PRO:HG3	1.66	0.78
4:Y:304:ARG:NH1	4:Y:320:THR:OG1	2.18	0.77
1:D:57:SER:OG	7:Z:8:MAN:H2	1.85	0.77
4:G:477:ASP:OD1	4:G:480:ARG:NH1	2.18	0.76
1:H:38:ARG:HB3	1:H:48:ILE:HD11	1.66	0.76
4:K:231:LYS:HD2	4:K:267:GLU:OE2	1.87	0.75
2:X:593:LEU:HD11	4:Y:494:LEU:HD21	1.68	0.75
2:J:664:ASP:O	4:Y:504:ARG:NH2	2.20	0.75
2:B:650:GLN:O	2:B:654:GLU:CB	2.35	0.75
4:K:233:PHE:O	4:K:273:ARG:NH2	2.18	0.75
4:K:113:ASP:OD1	4:K:429:ARG:NH2	2.21	0.74
11:Y:1041:NAG:O6	10:a:1:NAG:O6	2.05	0.73
4:Y:386:ASN:HB3	4:Y:417:PRO:HG2	1.69	0.73
4:K:231:LYS:HD2	4:K:267:GLU:OE1	1.88	0.73
2:B:601:LYS:C	2:X:655:LYS:HZ1	1.95	0.73
1:H:6:GLN:HE21	1:H:107:THR:HG23	1.52	0.73
4:G:292:VAL:HB	4:G:449:ILE:HB	1.72	0.72
4:G:278:THR:CG2	8:O:1:NAG:O6	2.30	0.72
2:J:593:LEU:HD11	4:K:494:LEU:HD21	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:598:CYS:O	2:J:600:GLY:N	2.20	0.72
4:Y:257:THR:HG23	4:Y:370:GLU:O	1.87	0.72
1:A:57:SER:HG	7:S:9:MAN:HO6	1.37	0.71
2:X:544:LEU:HD21	4:Y:493:PRO:HG3	1.72	0.71
2:B:608:VAL:CG2	2:B:649:SER:OG	2.36	0.71
2:B:602:LEU:HB2	2:X:655:LYS:NZ	2.04	0.71
4:Y:215:ILE:O	4:Y:251:ILE:N	2.23	0.71
4:Y:55:ALA:HB3	4:Y:216:HIS:HB2	1.71	0.71
1:A:159:LEU:HD21	1:A:182:VAL:HG11	1.73	0.71
2:B:603:ILE:C	2:B:603:ILE:HD12	2.16	0.70
2:B:631:TRP:CE2	2:B:635:ILE:HD11	2.26	0.70
2:B:645:LEU:C	2:B:649:SER:HB2	2.13	0.70
4:G:278:THR:CG2	8:O:1:NAG:H5	2.21	0.70
4:K:295:ASN:HA	4:K:446:VAL:HG12	1.72	0.70
1:D:74:LEU:HB3	1:D:77(C):PRO:HD3	1.72	0.70
2:J:610:TRP:HE3	4:K:36:VAL:HG12	1.56	0.70
4:Y:257:THR:HB	4:Y:375:SER:H	1.58	0.69
1:D:38:ARG:NH2	1:D:46:GLU:OE1	2.23	0.69
4:G:385:CYS:HB3	4:G:416:LEU:HD13	1.74	0.69
1:H:66:ARG:NH1	1:H:86:ASP:OD2	2.25	0.69
4:K:361:PHE:C	4:K:391:PHE:O	2.36	0.69
1:A:76:GLY:HA2	4:K:278:THR:HA	1.73	0.68
3:E:186:TYR:O	3:E:192:TYR:OH	2.11	0.68
4:Y:101:VAL:HG11	4:Y:480:ARG:HG2	1.75	0.68
1:D:74:LEU:HD11	4:Y:277:ILE:HB	1.74	0.68
2:J:544:LEU:HD21	4:K:493:PRO:HG3	1.74	0.68
3:L:30(A):GLY:O	3:L:50:ARG:NH1	2.26	0.68
1:H:57:SER:OG	7:N:9:MAN:O6	2.11	0.68
4:Y:304:ARG:NH2	4:Y:435:TYR:OH	2.26	0.68
4:Y:257:THR:HG21	4:Y:373:THR:O	1.94	0.68
4:G:476:ARG:HA	4:G:479:TRP:CD1	2.28	0.68
1:H:119:PRO:HB3	1:H:145:TYR:HB3	1.76	0.68
4:Y:249:HIS:HD2	4:Y:251:ILE:HG12	1.60	0.67
4:K:37:THR:OG1	4:K:499:THR:HG21	1.95	0.67
4:K:231:LYS:CD	4:K:267:GLU:OE2	2.41	0.67
4:K:201:ILE:HD11	4:K:435:TYR:HB2	1.75	0.66
2:B:598:CYS:O	2:B:599:SER:OG	2.12	0.66
4:Y:249:HIS:O	4:Y:251:ILE:HG13	1.95	0.66
1:H:74:LEU:HD13	1:H:77(C):PRO:HD3	1.78	0.66
4:Y:161:MET:SD	4:Y:162:THR:N	2.69	0.66
4:G:163:THR:HG23	4:G:165:LEU:H	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:K:231:LYS:CE	4:K:267:GLU:OE2	2.44	0.66
1:H:52:TRP:NE1	1:H:97:THR:HG21	2.11	0.65
4:G:278:THR:HG22	8:O:1:NAG:C6	2.26	0.65
4:K:292:VAL:HB	4:K:449:ILE:HB	1.79	0.65
4:G:71:THR:O	4:G:74:CYS:HB3	1.97	0.65
2:B:631:TRP:CZ2	2:B:635:ILE:HD11	2.31	0.65
1:H:73:ASP:OD1	1:H:74:LEU:N	2.26	0.65
1:H:94:THR:HB	1:H:102:SER:HB2	1.78	0.65
1:A:23:LYS:HA	1:A:77(D):ILE:HG12	1.79	0.65
3:E:186:TYR:O	3:E:211:ARG:NH2	2.29	0.64
4:G:278:THR:HG21	8:O:1:NAG:H5	1.80	0.64
1:D:100:ARG:NH1	4:Y:91:GLU:OE1	2.30	0.64
2:B:623:TRP:HZ2	4:G:499:THR:HG22	1.61	0.64
1:A:25:TYR:CD1	8:U:2:NAG:H2	2.33	0.64
4:Y:215:ILE:N	4:Y:251:ILE:O	2.28	0.64
4:Y:257:THR:CG2	4:Y:373:THR:O	2.46	0.64
3:L:136:LEU:HD21	3:L:196:VAL:HG21	1.79	0.64
4:K:165:LEU:HD23	4:K:167:ASP:H	1.62	0.63
1:H:25:TYR:CD1	8:O:2:NAG:H2	2.33	0.63
2:X:651:ASN:ND2	2:X:655:LYS:HD2	2.14	0.63
4:Y:216:HIS:ND1	4:Y:250:GLY:CA	2.60	0.63
4:K:231:LYS:CD	4:K:267:GLU:OE1	2.45	0.63
1:A:57:SER:OG	7:S:9:MAN:O6	2.13	0.63
1:D:68:LEU:HD23	7:Z:6:MAN:C6	2.29	0.63
1:H:23:LYS:HA	1:H:77(D):ILE:HG12	1.81	0.63
2:X:610:TRP:CE3	4:Y:36:VAL:HG12	2.34	0.63
4:K:231:LYS:HE3	4:K:267:GLU:OE2	1.99	0.62
2:B:570:VAL:HG12	2:B:571:TRP:H	1.63	0.62
2:X:610:TRP:HE3	4:Y:36:VAL:HG12	1.64	0.62
9:W:1:NAG:H62	9:W:2:NAG:C7	2.29	0.62
4:Y:216:HIS:CG	4:Y:250:GLY:HA2	2.35	0.62
4:Y:112:TRP:CH2	4:Y:210:PHE:HE2	2.18	0.62
4:K:391:PHE:HD2	4:K:470:PRO:HD3	1.64	0.62
5:F:1:NAG:H62	5:F:2:NAG:C7	2.29	0.62
1:H:57:SER:HG	7:N:9:MAN:HO6	1.40	0.61
4:Y:254:VAL:HG11	4:Y:261:LEU:O	2.00	0.61
4:Y:57:ASP:O	4:Y:77:THR:HG22	2.00	0.61
2:X:585:ARG:NH2	4:Y:491:ILE:O	2.34	0.61
4:Y:216:HIS:NE2	4:Y:250:GLY:CA	2.62	0.61
2:B:623:TRP:CZ2	4:G:499:THR:HG22	2.35	0.61
1:D:25:TYR:HD1	5:b:2:NAG:H2	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:159:LEU:HD21	1:D:182:VAL:HG11	1.83	0.61
4:K:273:ARG:HB2	4:K:285:LEU:HG	1.83	0.61
2:J:650:GLN:O	2:J:654:GLU:HB2	2.00	0.61
2:X:650:GLN:O	2:X:654:GLU:N	2.33	0.61
4:Y:91:GLU:OE2	4:Y:487:LYS:NZ	2.32	0.61
2:B:650:GLN:O	2:B:654:GLU:HB3	2.01	0.60
4:K:122:LEU:HB2	4:K:201:ILE:HG23	1.80	0.60
3:E:30:THR:HG21	2:X:616:ASN:H	1.65	0.60
4:K:265:LEU:HD21	4:K:291:PRO:HD3	1.82	0.60
4:K:361:PHE:HB3	4:K:391:PHE:O	2.01	0.60
2:B:591:GLN:NE2	2:J:541:ALA:O	2.36	0.59
4:Y:359:ILE:HD13	4:Y:361:PHE:HE1	1.68	0.59
1:A:211:VAL:HG12	1:A:211:VAL:O	2.01	0.59
4:K:223:PHE:CE1	4:K:490:LYS:HG2	2.37	0.59
1:A:25:TYR:HD1	8:U:2:NAG:H2	1.66	0.59
4:Y:259:LEU:HB2	4:Y:374:HIS:CE1	2.37	0.59
4:Y:475:MET:SD	4:Y:478:ASN:ND2	2.75	0.59
1:D:93:THR:HG21	1:D:100(L):PHE:CD1	2.38	0.59
4:Y:216:HIS:ND1	4:Y:250:GLY:HA2	2.17	0.59
2:B:593:LEU:HG	2:B:599:SER:H	1.66	0.59
4:G:253:PRO:HA	4:G:479:TRP:HZ3	1.68	0.59
4:G:478:ASN:OD1	4:G:479:TRP:N	2.35	0.59
4:K:195:ASN:HB3	4:K:423:ILE:HD12	1.85	0.59
1:H:210:ARG:HH11	1:H:210:ARG:CG	2.16	0.59
4:K:257:THR:HG21	4:K:370:GLU:O	2.03	0.58
4:K:391:PHE:CD2	4:K:470:PRO:HD3	2.37	0.58
4:Y:257:THR:OG1	4:Y:375:SER:OG	2.13	0.58
4:G:259:LEU:HB2	4:G:374:HIS:CE1	2.38	0.58
3:C:30:THR:HG21	2:J:615:SER:HA	1.86	0.58
2:B:602:LEU:HD22	2:X:655:LYS:HZ3	1.67	0.58
2:J:650:GLN:O	2:J:654:GLU:CB	2.51	0.58
4:Y:53:PHE:CZ	4:Y:218:CYS:HB2	2.37	0.58
4:K:233:PHE:C	4:K:271:MET:HE2	2.28	0.58
2:J:597:GLY:N	2:J:651:ASN:OD1	2.37	0.58
2:B:650:GLN:O	2:B:654:GLU:CA	2.50	0.58
4:G:423:ILE:HG22	4:G:435:TYR:HA	1.84	0.58
4:G:309:ILE:HG23	4:G:309:ILE:O	2.04	0.57
2:X:608:VAL:N	2:X:650:GLN:OE1	2.37	0.57
1:A:87:THR:HG23	1:A:110:SER:HA	1.85	0.57
1:H:93:THR:HG21	1:H:100(L):PHE:CD1	2.39	0.57
1:H:25:TYR:CE1	1:H:77(B):PRO:HG3	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:K:276:ASN:HB3	4:K:279:ASN:HB3	1.84	0.57
1:H:6:GLN:OE1	1:H:92:CYS:N	2.37	0.57
2:X:610:TRP:NE1	2:X:614:TRP:O	2.34	0.57
4:Y:428:GLN:O	4:Y:428:GLN:HG3	2.05	0.57
1:A:93:THR:HG21	1:A:100(L):PHE:CD1	2.40	0.57
1:H:126:PRO:HG2	1:H:213:PRO:HB3	1.86	0.57
4:Y:233:PHE:O	4:Y:273:ARG:NH2	2.38	0.57
2:B:523:LEU:HA	2:B:540:GLN:HE21	1.70	0.57
1:H:81:GLU:OE1	7:N:6:MAN:H62	2.05	0.57
1:A:75:THR:HG22	1:A:77:SER:H	1.69	0.56
4:Y:350:ARG:O	4:Y:354:GLY:N	2.36	0.56
4:Y:215:ILE:HD12	4:Y:253:PRO:HG3	1.87	0.56
1:D:73:ASP:HA	1:D:77(C):PRO:HB3	1.86	0.56
3:L:31:ASN:O	3:L:50:ARG:HA	2.05	0.56
1:D:68:LEU:CD2	7:Z:6:MAN:O6	2.53	0.56
4:Y:484:TYR:CE1	4:Y:485:LYS:HG2	2.41	0.56
4:G:212:PRO:HG2	6:M:1:NAG:O7	2.06	0.56
4:G:373:THR:HG21	4:G:384:TYR:HB3	1.86	0.56
4:K:373:THR:OG1	4:K:385:CYS:N	2.38	0.56
4:G:45:TRP:O	1:H:100(A):TRP:HH2	1.88	0.56
2:J:663:LEU:HD12	2:J:664:ASP:HB2	1.88	0.56
2:B:663:LEU:HD11	4:K:500:ARG:HE	1.72	0.55
1:D:5:VAL:O	1:D:23:LYS:N	2.36	0.55
1:H:52:TRP:HE1	1:H:97:THR:HG21	1.69	0.55
2:J:655:LYS:HZ3	2:X:602:LEU:H	1.52	0.55
4:Y:257:THR:O	4:Y:259:LEU:N	2.34	0.55
4:K:392:ASN:OD1	4:K:392:ASN:O	2.24	0.55
4:K:211:GLU:OE1	4:K:211:GLU:N	2.37	0.55
4:Y:305:LYS:N	4:Y:319:ALA:O	2.39	0.55
4:Y:291:PRO:HG3	11:Y:1041:NAG:H82	1.89	0.55
2:J:617:ARG:NH1	2:J:626:MET:SD	2.79	0.55
4:Y:217:TYR:O	4:Y:248:THR:HG23	2.05	0.55
2:B:617:ARG:NH1	2:B:634:GLU:OE2	2.34	0.55
3:E:65:SER:OG	3:E:72:THR:OG1	2.20	0.55
4:G:453:ILE:O	4:G:471:GLY:N	2.39	0.55
3:L:49:TYR:O	3:L:53:ALA:HB3	2.06	0.55
2:X:618:ASN:OD1	2:X:621:GLU:N	2.39	0.55
4:G:426:MET:HG3	4:G:427:TRP:CD2	2.42	0.55
4:K:195:ASN:HD22	4:K:201:ILE:HB	1.71	0.55
4:K:220:PRO:HG2	4:K:223:PHE:CD2	2.42	0.55
4:G:257:THR:O	4:G:259:LEU:N	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:618:ASN:O	2:J:622:ILE:HG13	2.07	0.55
4:K:257:THR:O	4:K:258:GLN:HB2	2.05	0.55
2:X:537:LEU:HD11	4:Y:39:TYR:HD2	1.72	0.55
2:B:614:TRP:HA	2:B:638:TYR:CD2	2.42	0.54
4:G:55:ALA:HA	4:G:75:VAL:O	2.08	0.54
3:C:37:GLN:HB2	3:C:86:PHE:CE1	2.42	0.54
4:G:57:ASP:HA	4:G:77:THR:HG22	1.89	0.54
2:B:585:ARG:NH2	4:G:491:ILE:O	2.41	0.54
1:D:100(A):TRP:O	9:W:2:NAG:H81	2.08	0.54
4:K:257:THR:O	4:K:374:HIS:ND1	2.40	0.54
4:Y:122:LEU:HB2	4:Y:201:ILE:HG23	1.88	0.54
4:G:252:LYS:HB3	4:G:254:VAL:HG23	1.89	0.54
1:A:25:TYR:CD1	1:A:77(B):PRO:HG3	2.43	0.54
1:H:159:LEU:HD21	1:H:182:VAL:HG11	1.89	0.54
3:L:53:ALA:HA	5:F:2:NAG:C6	2.38	0.54
4:G:202:THR:O	4:G:434:MET:HA	2.08	0.53
4:K:104:MET:O	4:K:108:ILE:HG12	2.07	0.53
4:K:48:ALA:HB3	4:K:490:LYS:HG3	1.90	0.53
2:X:585:ARG:NH1	2:X:589:ASP:OD2	2.39	0.53
2:J:653:GLN:O	2:J:657:GLU:HB2	2.09	0.53
2:B:608:VAL:HG22	2:B:649:SER:HG	1.71	0.53
4:K:112:TRP:CE3	4:K:427:TRP:HZ3	2.26	0.53
2:B:603:ILE:HD12	2:B:603:ILE:O	2.08	0.53
4:Y:280:ASN:HB3	4:Y:456:ARG:HB3	1.90	0.53
2:B:537:LEU:HD11	4:G:39:TYR:HD2	1.73	0.53
4:G:164:GLU:OE1	4:G:313:PRO:HA	2.09	0.53
2:B:601:LYS:CB	2:B:602:LEU:CD1	2.85	0.53
2:J:597:GLY:HA3	4:K:503:ARG:NH1	2.18	0.53
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.90	0.53
4:K:257:THR:HB	4:K:375:SER:H	1.73	0.53
2:X:523:LEU:HA	2:X:540:GLN:HE21	1.73	0.53
4:Y:161:MET:N	4:Y:170:GLN:O	2.42	0.53
4:Y:273:ARG:NH1	4:Y:484:TYR:CG	2.77	0.53
1:H:211:VAL:HG12	1:H:211:VAL:O	2.08	0.52
1:A:4:LEU:HD11	1:A:94:THR:HG22	1.90	0.52
4:Y:427:TRP:O	4:Y:428:GLN:C	2.51	0.52
1:D:24:ALA:HB3	1:D:77(C):PRO:O	2.08	0.52
4:K:48:ALA:CB	4:K:490:LYS:HG3	2.40	0.52
4:K:161:MET:SD	4:K:162:THR:N	2.82	0.52
4:Y:172:VAL:HG11	4:Y:307:ILE:HD13	1.92	0.52
4:K:233:PHE:O	4:K:271:MET:HE2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:Y:216:HIS:CD2	4:Y:250:GLY:CA	2.92	0.52
4:Y:377:ASN:HA	4:Y:382:PHE:HA	1.92	0.52
1:A:35:VAL:HA	1:A:93:THR:O	2.10	0.52
1:A:40:ALA:HB3	1:A:43:GLN:HB2	1.91	0.52
1:A:55:TRP:CZ3	7:S:2:NAG:H61	2.44	0.52
1:A:163:VAL:HG22	1:A:182:VAL:HG22	1.91	0.52
2:B:602:LEU:CD1	2:B:602:LEU:N	2.73	0.52
2:X:570:VAL:HG22	2:X:571:TRP:H	1.74	0.52
4:K:223:PHE:CD1	4:K:490:LYS:HG2	2.44	0.52
2:B:601:LYS:O	2:X:655:LYS:CE	2.58	0.52
3:L:186:TYR:HA	3:L:192:TYR:OH	2.10	0.52
4:Y:123:THR:N	4:Y:124:PRO:CD	2.73	0.52
4:Y:216:HIS:CG	4:Y:250:GLY:CA	2.92	0.52
4:Y:364:SER:HB2	4:Y:470:PRO:HG2	1.91	0.52
4:K:260:LEU:HD12	4:K:451:GLY:HA3	1.91	0.52
4:K:296:CYS:SG	4:K:376:PHE:HZ	2.33	0.52
2:B:645:LEU:O	2:B:649:SER:CA	2.57	0.52
4:K:349:LEU:HD21	4:K:468:PHE:CE2	2.45	0.52
2:J:574:LYS:HE3	4:K:52:LEU:O	2.10	0.52
1:D:25:TYR:CD1	1:D:77(B):PRO:HG3	2.45	0.51
3:E:37:GLN:CB	3:E:47:LEU:HD11	2.39	0.51
4:K:295:ASN:N	4:K:332:ASN:O	2.36	0.51
4:K:349:LEU:HD21	4:K:468:PHE:HE2	1.75	0.51
4:K:364:SER:HB2	4:K:470:PRO:HG2	1.92	0.51
3:E:37:GLN:HB2	3:E:86:PHE:CE1	2.45	0.51
4:K:212:PRO:HG2	10:T:1:NAG:O7	2.10	0.51
1:D:100(A):TRP:CZ2	4:Y:46:LYS:HE2	2.45	0.51
3:L:19:VAL:HG12	3:L:75:ILE:HB	1.92	0.51
1:D:100:ARG:O	2:X:633:LYS:NZ	2.39	0.51
2:J:607:ASN:OD1	2:J:607:ASN:N	2.44	0.51
2:B:647:GLU:CG	2:B:648:GLU:N	2.73	0.51
4:Y:304:ARG:HD2	4:Y:320:THR:HA	1.92	0.51
2:B:602:LEU:CB	2:X:655:LYS:HZ3	2.18	0.51
2:J:606:THR:OG1	4:K:36:VAL:O	2.29	0.51
1:A:152:VAL:HG22	1:A:198:VAL:HG22	1.93	0.50
1:H:100(I):VAL:HG12	3:L:91:TYR:HB2	1.93	0.50
4:G:55:ALA:HB3	4:G:216:HIS:HB2	1.93	0.50
4:G:112:TRP:CD2	4:G:427:TRP:HZ3	2.29	0.50
4:Y:499:THR:HG23	4:Y:501:CYS:H	1.75	0.50
1:A:32:LEU:C	1:A:97:THR:HG22	2.36	0.50
1:D:25:TYR:CD1	5:b:2:NAG:H2	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:68:LEU:HD23	7:Z:6:MAN:H62	1.91	0.50
1:D:68:LEU:HD23	7:Z:6:MAN:O6	2.11	0.50
4:K:280:ASN:OD1	4:K:281:ALA:N	2.44	0.50
4:Y:299:PRO:HA	4:Y:442:VAL:HG13	1.93	0.50
1:A:204:ASN:O	1:A:204:ASN:ND2	2.44	0.50
2:B:645:LEU:O	2:B:649:SER:N	2.45	0.50
1:D:35(A):ASN:O	1:D:93:THR:HB	2.12	0.50
2:J:610:TRP:CE3	4:K:36:VAL:HG12	2.43	0.50
1:H:25:TYR:HD1	8:O:2:NAG:H2	1.75	0.50
4:K:122:LEU:O	4:K:123:THR:C	2.55	0.50
2:B:598:CYS:O	2:B:600:GLY:N	2.42	0.50
2:J:638:TYR:O	2:J:642:ILE:HG13	2.11	0.50
1:D:50:GLN:N	1:D:58:SER:OG	2.40	0.50
1:D:64:ARG:HD2	7:Z:9:MAN:O3	2.12	0.50
3:E:189:HIS:NE2	3:E:191:VAL:HG22	2.27	0.50
2:J:570:VAL:HG22	2:J:571:TRP:H	1.77	0.50
4:G:193:LEU:HD12	4:G:193:LEU:H	1.76	0.50
2:J:651:ASN:O	2:J:655:LYS:HG2	2.12	0.50
4:K:277:ILE:HD12	7:S:1:NAG:H81	1.94	0.50
4:Y:383:PHE:CZ	4:Y:385:CYS:SG	3.05	0.50
1:H:74:LEU:CD1	1:H:77(B):PRO:HA	2.42	0.49
3:C:28:SER:HA	3:C:69:THR:HG22	1.94	0.49
2:X:660:LEU:O	2:X:663:LEU:HG	2.12	0.49
4:K:255:VAL:HG13	4:K:475:MET:SD	2.51	0.49
3:L:54:LEU:HD11	5:F:3:BMA:H3	1.94	0.49
3:L:65:SER:OG	3:L:72:THR:OG1	2.29	0.49
4:Y:180:ASP:OD1	4:Y:181:VAL:HG13	2.12	0.49
2:B:610:TRP:HE3	4:G:36:VAL:HG22	1.77	0.49
3:C:166:GLN:HG3	3:C:173:TYR:CZ	2.47	0.49
2:B:614:TRP:O	3:L:30:THR:HG21	2.12	0.49
2:J:522:PHE:CD1	2:J:543:ASN:HB3	2.47	0.49
4:K:391:PHE:CE2	4:K:470:PRO:HB3	2.48	0.49
1:H:74:LEU:HB2	1:H:77(C):PRO:HG3	1.95	0.49
2:X:585:ARG:HH22	4:Y:493:PRO:HD3	1.77	0.49
4:Y:257:THR:HB	4:Y:375:SER:N	2.26	0.49
2:B:596:TRP:C	2:B:651:ASN:ND2	2.71	0.49
2:B:596:TRP:C	2:B:651:ASN:HD21	2.15	0.49
2:B:615:SER:HA	3:L:30:THR:HG21	1.93	0.49
4:G:78:ASP:OD1	4:G:78:ASP:N	2.46	0.49
4:K:123:THR:N	4:K:124:PRO:CD	2.76	0.49
3:C:30:THR:OG1	2:J:615:SER:N	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:147:PRO:O	1:D:200:HIS:NE2	2.37	0.48
4:G:104:MET:O	4:G:108:ILE:HG12	2.13	0.48
4:Y:195:ASN:HB3	4:Y:423:ILE:CD1	2.43	0.48
1:H:210:ARG:HG2	1:H:210:ARG:NH1	2.20	0.48
2:B:595:ILE:HG22	2:B:647:GLU:HB2	1.95	0.48
2:B:598:CYS:C	2:B:600:GLY:H	2.20	0.48
4:G:180:ASP:CG	4:G:422:GLN:H	2.21	0.48
4:Y:96:TRP:CE3	4:Y:275:GLU:HB2	2.49	0.48
4:Y:285:LEU:HD21	4:Y:477:ASP:HB3	1.95	0.48
1:D:163:VAL:HG22	1:D:182:VAL:HG22	1.95	0.48
1:H:126:PRO:HD2	1:H:213:PRO:HA	1.95	0.48
4:Y:217:TYR:N	4:Y:248:THR:OG1	2.41	0.48
1:D:23:LYS:HA	1:D:77(D):ILE:HG22	1.94	0.48
4:G:257:THR:OG1	4:G:375:SER:N	2.43	0.48
2:X:598:CYS:O	2:X:599:SER:HB3	2.13	0.48
4:Y:179:LEU:HD13	4:Y:369:LEU:HD21	1.96	0.48
2:B:645:LEU:O	2:B:649:SER:HB3	2.03	0.48
4:G:120:VAL:HG23	4:G:315:GLN:HB3	1.95	0.48
4:Y:259:LEU:HD13	4:Y:449:ILE:HD13	1.95	0.48
3:E:61:ARG:NH1	3:E:82:ASP:OD2	2.47	0.48
3:E:137:ASN:OD1	3:E:174:SER:HB3	2.13	0.48
2:B:610:TRP:HH2	2:B:635:ILE:HD13	1.79	0.48
1:D:35:VAL:HA	1:D:93:THR:O	2.14	0.48
4:K:270:VAL:HB	4:K:287:GLN:O	2.14	0.48
4:K:429:ARG:CZ	4:K:432:GLN:HG2	2.44	0.48
4:G:101:VAL:HG21	4:G:480:ARG:HG2	1.96	0.48
4:K:93:PHE:HB2	4:K:233:PHE:HZ	1.76	0.48
4:Y:427:TRP:O	4:Y:428:GLN:HG2	2.14	0.48
4:G:278:THR:CG2	8:O:1:NAG:C5	2.92	0.47
1:H:33:TYR:CD2	1:H:94:THR:HG22	2.49	0.47
1:A:146:PHE:HB2	1:A:175:LEU:HD12	1.96	0.47
3:C:35:TRP:CZ3	3:C:88:CYS:HB3	2.49	0.47
1:D:211:VAL:O	1:D:211:VAL:HG12	2.13	0.47
1:H:210:ARG:CG	1:H:210:ARG:NH1	2.75	0.47
4:K:57:ASP:HB3	4:K:58:ALA:H	1.49	0.47
3:E:19:VAL:HG12	3:E:75:ILE:HB	1.96	0.47
4:K:371:VAL:HG22	4:K:473:GLY:N	2.29	0.47
4:G:338:TRP:CZ2	4:G:390:LEU:HB3	2.50	0.47
2:B:650:GLN:C	2:B:654:GLU:HB3	2.39	0.47
1:D:23:LYS:HA	1:D:77(D):ILE:CG2	2.44	0.47
1:D:68:LEU:HD21	7:Z:6:MAN:O6	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:42:VAL:HG22	4:G:493:PRO:O	2.14	0.47
4:K:229:LYS:HE2	4:K:243:SER:OG	2.14	0.47
4:Y:346:VAL:O	4:Y:350:ARG:N	2.44	0.47
3:C:123:GLU:HA	3:C:126:LYS:HZ3	1.80	0.47
1:D:74:LEU:HB3	1:D:77(C):PRO:CD	2.43	0.47
4:K:330:HIS:HA	4:K:417:PRO:HA	1.95	0.47
4:K:386:ASN:OD1	4:K:388:SER:CB	2.63	0.47
4:K:484:TYR:CE2	4:K:485:LYS:HG3	2.50	0.47
1:H:35(A):ASN:O	1:H:93:THR:HB	2.14	0.47
2:J:523:LEU:H	2:J:540:GLN:HG3	1.79	0.47
1:A:166:PHE:CE2	3:C:176:SER:HB3	2.50	0.47
4:G:113:ASP:OD1	4:G:114:GLN:N	2.48	0.47
1:H:60:SER:O	1:H:64:ARG:N	2.47	0.47
1:H:67:VAL:O	7:N:9:MAN:H62	2.15	0.47
4:K:371:VAL:HG22	4:K:473:GLY:H	1.80	0.47
4:Y:259:LEU:HD22	4:Y:452:LEU:CD2	2.45	0.47
4:Y:215:ILE:C	4:Y:251:ILE:H	2.20	0.47
2:B:598:CYS:C	2:B:600:GLY:N	2.73	0.46
2:X:656:ASN:O	2:X:660:LEU:HG	2.15	0.46
1:D:100(E):HIS:HB2	1:D:100(J):MET:HE2	1.97	0.46
1:A:59:ALA:HB3	7:S:9:MAN:H4	1.97	0.46
1:H:82(C):LEU:HB3	1:H:111:VAL:HG21	1.97	0.46
1:H:100(F):HIS:HB2	3:L:32:TRP:CE3	2.50	0.46
2:J:626:MET:HE2	2:J:631:TRP:HA	1.98	0.46
2:B:618:ASN:HB2	2:B:621:GLU:HG3	1.98	0.46
1:H:93:THR:HA	1:H:102:SER:O	2.15	0.46
1:A:30:PHE:CE2	1:A:77(C):PRO:HB2	2.51	0.46
2:B:602:LEU:N	2:B:602:LEU:HD12	2.30	0.46
4:G:229:LYS:O	4:G:241:SER:OG	2.33	0.46
4:K:202:THR:O	4:K:434:MET:HA	2.15	0.46
3:L:11:LEU:HD13	3:L:19:VAL:HG21	1.97	0.46
2:X:522:PHE:CE1	2:X:523:LEU:HG	2.50	0.46
1:D:119:PRO:HB3	1:D:145:TYR:HB3	1.97	0.46
1:H:24:ALA:O	1:H:77(B):PRO:HB2	2.16	0.46
1:A:67:VAL:O	7:S:9:MAN:H62	2.16	0.46
3:C:146:VAL:HG12	3:C:196:VAL:HG22	1.98	0.46
4:K:227:LYS:HA	4:K:485:LYS:O	2.16	0.46
3:C:19:VAL:HG12	3:C:75:ILE:HB	1.97	0.45
1:D:50:GLN:HG2	1:D:58:SER:OG	2.15	0.45
3:L:124:GLN:HG2	3:L:129:THR:O	2.17	0.45
3:L:134:CYS:HB2	3:L:148:TRP:CH2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:144:ALA:HB2	3:L:198:HIS:HD2	1.81	0.45
3:C:187:GLU:CD	3:C:211:ARG:HD2	2.40	0.45
3:E:134:CYS:HB2	3:E:148:TRP:CH2	2.51	0.45
4:G:494:LEU:HD23	4:G:494:LEU:HA	1.85	0.45
2:J:606:THR:HA	4:K:503:ARG:HE	1.82	0.45
4:K:270:VAL:HG12	4:K:288:PHE:HA	1.97	0.45
4:Y:122:LEU:O	4:Y:123:THR:C	2.59	0.45
1:D:100(A):TRP:CH2	4:Y:46:LYS:HE2	2.52	0.45
2:J:580:VAL:O	2:J:584:GLU:HG2	2.16	0.45
3:L:124:GLN:O	3:L:127:SER:OG	2.31	0.45
2:B:598:CYS:O	2:B:599:SER:CB	2.63	0.45
2:B:603:ILE:C	2:B:603:ILE:CD1	2.85	0.45
1:D:22:CYS:O	1:D:77(D):ILE:HB	2.16	0.45
4:Y:164:GLU:HG2	4:Y:312:GLY:HA3	1.98	0.45
4:Y:272:ILE:HD11	4:Y:352:HIS:ND1	2.32	0.45
3:L:53:ALA:HA	5:F:2:NAG:H62	1.98	0.45
4:Y:33:ASN:OD1	4:Y:33:ASN:N	2.50	0.45
1:D:64:ARG:NE	7:Z:9:MAN:O3	2.50	0.45
3:E:37:GLN:HB3	3:E:47:LEU:HD11	1.98	0.45
2:J:598:CYS:C	2:J:600:GLY:N	2.72	0.45
4:Y:113:ASP:O	4:Y:117:LYS:HD3	2.17	0.45
1:A:70:SER:HB2	7:S:3:BMA:O4	2.16	0.45
2:B:540:GLN:N	2:B:540:GLN:OE1	2.49	0.45
4:K:201:ILE:HA	4:K:433:ALA:HB3	1.99	0.45
4:K:257:THR:HB	4:K:375:SER:N	2.32	0.45
3:E:146:VAL:HG12	3:E:196:VAL:HG22	1.99	0.45
4:G:278:THR:HG22	8:O:1:NAG:C5	2.46	0.45
4:G:371:VAL:HG22	4:G:473:GLY:H	1.81	0.45
4:K:37:THR:OG1	4:K:499:THR:CG2	2.64	0.45
4:Y:359:ILE:HD13	4:Y:361:PHE:CE1	2.48	0.45
1:D:48:ILE:HG23	1:D:63:PHE:CD2	2.52	0.44
1:D:75:THR:HG22	1:D:77:SER:H	1.81	0.44
4:G:87:GLU:O	4:G:89:VAL:HG23	2.17	0.44
4:Y:477:ASP:OD1	4:Y:480:ARG:NH1	2.49	0.44
4:G:105:HIS:CG	4:G:476:ARG:HG2	2.52	0.44
2:X:615:SER:OG	2:X:634:GLU:HG2	2.18	0.44
2:B:604:CYS:HB3	4:G:38:VAL:HB	1.99	0.44
2:J:611:ASN:HB3	2:J:614:TRP:CD2	2.52	0.44
4:Y:249:HIS:CD2	4:Y:251:ILE:CG1	2.95	0.44
4:Y:474:ASP:OD1	4:Y:476:ARG:HG3	2.18	0.44
1:D:87:THR:HG23	1:D:110:SER:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:147:PRO:O	1:H:200:HIS:NE2	2.38	0.44
4:Y:494:LEU:HD23	4:Y:494:LEU:HA	1.66	0.44
4:K:278:THR:CG2	8:U:1:NAG:H62	2.48	0.44
4:K:361:PHE:CB	4:K:391:PHE:O	2.66	0.44
4:Y:356:ASN:OD1	4:Y:357:THR:N	2.51	0.44
4:G:163:THR:OG1	4:G:164:GLU:N	2.50	0.44
4:G:476:ARG:HA	4:G:479:TRP:HD1	1.77	0.44
3:L:37:GLN:HB2	3:L:86:PHE:CE1	2.52	0.44
4:Y:304:ARG:NH1	4:Y:438:PRO:HG2	2.32	0.44
2:B:582:ALA:HB1	4:G:221:ALA:HB3	1.99	0.44
4:G:37:THR:HG21	4:G:499:THR:OG1	2.18	0.44
4:G:195:ASN:OD1	4:G:196:CYS:N	2.51	0.44
1:H:40:ALA:HB3	1:H:43:GLN:HB2	1.98	0.44
4:K:374:HIS:HD2	4:K:376:PHE:CE1	2.36	0.44
2:X:646:LEU:HD21	4:Y:36:VAL:HG21	2.00	0.44
4:Y:249:HIS:CD2	4:Y:249:HIS:O	2.70	0.44
4:Y:390:LEU:HG	4:Y:416:LEU:HD21	1.98	0.44
6:Q:1:NAG:H62	6:Q:2:NAG:C7	2.47	0.44
4:G:91:GLU:OE1	1:H:100:ARG:NH2	2.51	0.44
4:G:112:TRP:CE3	4:G:427:TRP:HZ3	2.36	0.44
4:K:499:THR:HG23	4:K:501:CYS:H	1.83	0.44
4:Y:252:LYS:HA	4:Y:253:PRO:HD2	1.90	0.44
1:A:119:PRO:HB3	1:A:145:TYR:HB3	1.99	0.43
3:E:30(A):GLY:HA3	3:E:32:TRP:CZ3	2.53	0.43
3:E:2:ILE:HG12	3:E:93:THR:HG22	1.99	0.43
4:G:233:PHE:O	4:G:273:ARG:NH2	2.44	0.43
1:H:55:TRP:CZ3	7:N:2:NAG:H61	2.53	0.43
1:H:71:ALA:HA	1:H:78:SER:HA	1.98	0.43
3:L:167:ASP:HB3	3:L:170:ASP:OD1	2.17	0.43
4:Y:259:LEU:HD22	4:Y:452:LEU:HD22	1.99	0.43
4:K:374:HIS:CD2	4:K:376:PHE:CE1	3.07	0.43
4:Y:270:VAL:HG13	4:Y:287:GLN:O	2.19	0.43
4:G:35:TRP:O	4:G:498:PRO:HA	2.18	0.43
4:G:277:ILE:CD1	7:N:1:NAG:H81	2.49	0.43
2:X:653:GLN:O	2:X:657:GLU:HB2	2.17	0.43
1:A:38:ARG:HB3	1:A:48:ILE:HD11	2.00	0.43
1:A:94:THR:CG2	1:A:102:SER:HB2	2.48	0.43
4:G:446:VAL:HB	6:M:2:NAG:O7	2.18	0.43
1:H:35:VAL:HA	1:H:93:THR:O	2.17	0.43
4:Y:163:THR:OG1	4:Y:164:GLU:N	2.52	0.43
3:C:195:GLU:N	3:C:195:GLU:OE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:86:LEU:HB3	4:G:89:VAL:HG21	2.00	0.43
4:G:252:LYS:HG3	4:G:262:ASN:O	2.18	0.43
2:X:570:VAL:O	2:X:571:TRP:HB3	2.19	0.43
2:X:605:CYS:O	4:Y:503:ARG:HD3	2.18	0.43
3:C:4:MET:HE3	3:C:23:CYS:SG	2.59	0.43
3:C:31:ASN:O	3:C:50:ARG:HA	2.19	0.43
3:E:189:HIS:CD2	3:E:191:VAL:H	2.37	0.43
1:H:25:TYR:CD1	1:H:77(B):PRO:HG3	2.54	0.43
2:J:522:PHE:CD1	2:J:540:GLN:HA	2.54	0.43
2:B:577:GLN:N	2:B:577:GLN:OE1	2.51	0.43
4:K:351:LYS:O	4:K:352:HIS:C	2.59	0.43
4:Y:249:HIS:CD2	4:Y:249:HIS:C	2.96	0.43
4:Y:373:THR:HG22	4:Y:419:ARG:HH21	1.84	0.43
3:C:58:VAL:HG13	3:C:59:PRO:HD2	2.01	0.43
2:J:659:ASP:HB2	2:X:603:ILE:HD13	2.00	0.43
4:K:223:PHE:HE1	4:K:490:LYS:HG2	1.83	0.43
4:K:294:ILE:HG22	4:K:447:SER:O	2.19	0.43
3:E:11:LEU:HD13	3:E:19:VAL:HG21	2.01	0.42
3:E:37:GLN:NE2	3:E:82:ASP:O	2.50	0.42
4:K:78:ASP:OD1	4:K:78:ASP:N	2.50	0.42
4:K:382:PHE:O	4:K:420:ILE:HA	2.19	0.42
1:D:153:SER:O	1:D:197:ASN:N	2.39	0.42
4:Y:342:LEU:O	4:Y:346:VAL:HG23	2.19	0.42
2:B:522:PHE:CD1	2:B:543:ASN:HB2	2.55	0.42
3:C:32:TRP:HH2	2:J:633:LYS:HG2	1.85	0.42
4:G:277:ILE:HD11	7:N:1:NAG:H81	2.02	0.42
4:K:122:LEU:O	4:K:124:PRO:N	2.53	0.42
4:Y:453:ILE:O	4:Y:471:GLY:N	2.39	0.42
4:G:122:LEU:HD21	4:G:125:LEU:HD23	2.00	0.42
4:Y:172:VAL:HG11	4:Y:307:ILE:HG21	2.02	0.42
2:B:615:SER:O	2:B:616:ASN:HB3	2.20	0.42
3:C:36:TYR:CE1	3:C:46:LEU:HD13	2.54	0.42
4:Y:199:SER:HB3	4:Y:431:GLY:H	1.85	0.42
1:A:94:THR:HG23	1:A:102:SER:HB2	2.01	0.42
4:K:165:LEU:HD23	4:K:166:ARG:N	2.35	0.42
3:E:149:LYS:HD3	3:E:152:ASN:HA	2.01	0.42
1:H:33:TYR:HD2	1:H:94:THR:HG22	1.84	0.42
4:Y:93:PHE:HB2	4:Y:233:PHE:HZ	1.84	0.42
4:Y:297:THR:HG23	4:Y:443:ILE:O	2.19	0.42
9:R:1:NAG:H62	9:R:2:NAG:C7	2.50	0.42
1:A:26:GLY:HA2	8:U:1:NAG:O3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:20:ARG:HG2	3:C:74:THR:HG22	2.02	0.42
3:C:187:GLU:OE2	3:C:211:ARG:NH1	2.52	0.42
4:Y:101:VAL:HG12	4:Y:483:LEU:HD12	2.02	0.42
1:D:72:VAL:HG23	7:Z:5:MAN:H62	2.02	0.41
4:K:42:VAL:HG22	4:K:493:PRO:O	2.20	0.41
3:L:4:MET:HE3	3:L:23:CYS:SG	2.60	0.41
3:E:189:HIS:CD2	3:E:191:VAL:HG22	2.55	0.41
4:G:101:VAL:HG22	4:G:483:LEU:HD12	2.02	0.41
4:G:339:ASN:OD1	4:G:340:GLU:N	2.53	0.41
4:K:112:TRP:CD2	4:K:427:TRP:HZ3	2.38	0.41
4:K:307:ILE:HD11	4:K:317:PHE:HD2	1.84	0.41
4:K:320:THR:OG1	4:K:438:PRO:HG2	2.20	0.41
1:A:200:HIS:CD2	1:A:202:PRO:HD2	2.55	0.41
4:Y:355:ASN:OD1	4:Y:356:ASN:N	2.53	0.41
9:W:1:NAG:H62	9:W:2:NAG:C8	2.50	0.41
4:K:231:LYS:HD3	4:K:267:GLU:OE1	2.19	0.41
3:L:113:PRO:HB3	3:L:139:PHE:HB3	2.02	0.41
4:Y:87:GLU:O	4:Y:89:VAL:HG23	2.20	0.41
1:D:200:HIS:CD2	1:D:202:PRO:HD2	2.55	0.41
4:G:165:LEU:HA	4:Y:126:CYS:O	2.20	0.41
4:K:198:THR:HG22	4:K:430:ILE:HG23	2.01	0.41
4:Y:214:PRO:HA	4:Y:251:ILE:O	2.20	0.41
4:Y:453:ILE:HD11	4:Y:478:ASN:OD1	2.20	0.41
2:B:608:VAL:HG21	2:B:646:LEU:HD23	2.01	0.41
4:K:257:THR:HG22	4:K:258:GLN:HG3	2.03	0.41
3:E:4:MET:HE3	3:E:23:CYS:SG	2.60	0.41
4:K:453:ILE:O	4:K:454:LEU:HD23	2.21	0.41
4:Y:195:ASN:HB3	4:Y:423:ILE:HD13	2.02	0.41
4:G:122:LEU:HD13	4:G:203:GLN:CD	2.45	0.41
4:G:159:PHE:CZ	4:G:172:VAL:HG21	2.56	0.41
2:J:650:GLN:O	2:J:654:GLU:HB3	2.19	0.41
1:A:69:ILE:N	7:S:9:MAN:O6	2.54	0.41
1:D:64:ARG:CD	7:Z:9:MAN:O3	2.69	0.41
4:G:257:THR:O	4:G:258:GLN:HB2	2.21	0.41
4:G:277:ILE:HG22	1:H:76:GLY:HA2	2.02	0.41
4:G:453:ILE:O	4:G:454:LEU:HD23	2.20	0.41
4:K:83:GLU:HA	4:K:245:VAL:HG12	2.02	0.41
4:K:230:ASP:OD1	4:K:233:PHE:HB2	2.21	0.41
4:K:270:VAL:HG12	4:K:288:PHE:C	2.46	0.41
3:L:140:TYR:CD1	3:L:141:PRO:HA	2.55	0.41
4:Y:195:ASN:HD22	4:Y:201:ILE:HB	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:220:PRO:HG2	4:G:223:PHE:CD2	2.56	0.41
4:Y:163:THR:HG23	4:Y:165:LEU:H	1.84	0.41
2:B:599:SER:OG	2:B:600:GLY:N	2.54	0.40
2:B:611:ASN:OD1	2:B:614:TRP:CG	2.74	0.40
2:B:656:ASN:O	2:B:659:ASP:HB3	2.21	0.40
4:G:199:SER:HB3	4:G:431:GLY:H	1.86	0.40
1:H:6:GLN:NE2	1:H:107:THR:HG23	2.26	0.40
2:J:619:LEU:HD23	2:J:619:LEU:HA	1.86	0.40
3:L:46:LEU:HD23	3:L:55:LEU:HD22	2.02	0.40
1:A:37:VAL:HG22	1:A:47:TYR:HA	2.03	0.40
1:H:95:THR:HA	1:H:101:SER:OG	2.21	0.40
4:Y:253:PRO:HA	4:Y:479:TRP:HZ3	1.86	0.40
3:E:33:VAL:HA	3:E:89:GLN:O	2.20	0.40
1:H:100(E):HIS:HA	1:H:100(J):MET:HB3	2.03	0.40
4:Y:393:SER:HB3	4:Y:395:TRP:NE1	2.37	0.40
2:X:651:ASN:CG	2:X:655:LYS:HD2	2.45	0.40
1:A:60:SER:O	1:A:64:ARG:N	2.55	0.40
3:E:148:TRP:CG	3:E:179:LEU:HD22	2.57	0.40
4:Y:349:LEU:HD21	4:Y:468:PHE:CZ	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	220/244 (90%)	212 (96%)	8 (4%)	0	100	100
1	D	220/244 (90%)	210 (96%)	10 (4%)	0	100	100
1	H	223/244 (91%)	214 (96%)	9 (4%)	0	100	100
2	B	120/153 (78%)	107 (89%)	13 (11%)	0	100	100
2	J	120/153 (78%)	112 (93%)	8 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	X	117/153 (76%)	108 (92%)	9 (8%)	0	100	100
3	C	212/215 (99%)	210 (99%)	2 (1%)	0	100	100
3	E	212/215 (99%)	209 (99%)	3 (1%)	0	100	100
3	L	212/215 (99%)	208 (98%)	4 (2%)	0	100	100
4	G	382/479 (80%)	358 (94%)	23 (6%)	1 (0%)	36	65
4	K	391/479 (82%)	372 (95%)	19 (5%)	0	100	100
4	Y	383/479 (80%)	367 (96%)	16 (4%)	0	100	100
All	All	2812/3273 (86%)	2687 (96%)	124 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	G	312	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	184/210 (88%)	183 (100%)	1 (0%)	81	80
1	D	175/210 (83%)	175 (100%)	0	100	100
1	H	185/210 (88%)	183 (99%)	2 (1%)	65	74
2	B	99/129 (77%)	97 (98%)	2 (2%)	48	66
2	J	100/129 (78%)	100 (100%)	0	100	100
2	X	99/129 (77%)	98 (99%)	1 (1%)	68	75
3	C	170/182 (93%)	170 (100%)	0	100	100
3	E	163/182 (90%)	163 (100%)	0	100	100
3	L	157/182 (86%)	156 (99%)	1 (1%)	78	79
4	G	289/427 (68%)	285 (99%)	4 (1%)	59	71
4	K	315/427 (74%)	310 (98%)	5 (2%)	55	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	Y	312/427 (73%)	310 (99%)	2 (1%)	78	79
All	All	2248/2844 (79%)	2230 (99%)	18 (1%)	73	77

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	173	SER
2	B	602	LEU
2	B	664	ASP
4	G	74	CYS
4	G	120	VAL
4	G	193	LEU
4	G	378	CYS
1	H	93	THR
1	H	210	ARG
4	K	57	ASP
4	K	196	CYS
4	K	234	ASN
4	K	268	GLU
4	K	418	CYS
3	L	30	THR
2	X	650	GLN
4	Y	378	CYS
4	Y	448	ASN

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	50	GLN
1	A	171	GLN
2	B	651	ASN
3	C	137	ASN
1	D	171	GLN
3	E	79	GLN
4	G	103	GLN
4	G	302	ASN
4	G	315	GLN
1	H	3	HIS
1	H	164	HIS
4	K	33	ASN
4	K	374	HIS

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Mol	Chain	Res	Type
4	K	398	ASN
4	Y	33	ASN
4	Y	289	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 ⓘ

76 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	$\# Z > 2$
5	NAG	F	1	2,5	14,14,15	0.60	0	17,19,21	0.59	0
5	NAG	F	2	5	14,14,15	0.40	0	17,19,21	0.62	0
5	BMA	F	3	5	11,11,12	0.94	0	15,15,17	0.74	0
5	MAN	F	4	5	11,11,12	0.71	0	15,15,17	1.27	2 (13%)
5	MAN	F	5	5	11,11,12	0.73	0	15,15,17	1.06	2 (13%)
6	NAG	I	1	6,4	14,14,15	0.60	1 (7%)	17,19,21	0.57	0
6	NAG	I	2	6	14,14,15	0.17	0	17,19,21	0.42	0
6	NAG	M	1	6,4	14,14,15	0.48	0	17,19,21	0.39	0
6	NAG	M	2	6	14,14,15	0.15	0	17,19,21	0.39	0
7	NAG	N	1	4,7	14,14,15	0.36	0	17,19,21	0.56	0
7	MAN	N	10	7	11,11,12	0.74	0	15,15,17	0.86	0
7	NAG	N	2	7	14,14,15	0.63	1 (7%)	17,19,21	0.53	0
7	BMA	N	3	7	11,11,12	0.72	0	15,15,17	0.79	0
7	MAN	N	4	7	11,11,12	0.41	0	15,15,17	1.16	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	MAN	N	5	7	11,11,12	0.57	0	15,15,17	1.14	1 (6%)
7	MAN	N	6	7	11,11,12	0.80	0	15,15,17	1.03	2 (13%)
7	MAN	N	7	7	11,11,12	0.57	0	15,15,17	1.15	2 (13%)
7	MAN	N	8	7	11,11,12	0.48	0	15,15,17	1.22	2 (13%)
7	MAN	N	9	7	11,11,12	0.65	0	15,15,17	1.00	2 (13%)
8	NAG	O	1	8,4	14,14,15	0.33	0	17,19,21	0.48	0
8	NAG	O	2	8	14,14,15	0.22	0	17,19,21	0.53	0
8	BMA	O	3	8	11,11,12	0.95	0	15,15,17	0.98	1 (6%)
8	MAN	O	4	8	11,11,12	0.82	1 (9%)	15,15,17	1.11	2 (13%)
8	MAN	O	5	8	11,11,12	0.66	0	15,15,17	1.02	2 (13%)
8	MAN	O	6	8	11,11,12	0.79	0	15,15,17	1.00	2 (13%)
6	NAG	P	1	1,6	14,14,15	0.69	1 (7%)	17,19,21	0.55	0
6	NAG	P	2	6	14,14,15	0.16	0	17,19,21	0.51	0
6	NAG	Q	1	6,2	14,14,15	0.42	0	17,19,21	0.70	0
6	NAG	Q	2	6	14,14,15	0.20	0	17,19,21	0.57	0
9	NAG	R	1	9,2	14,14,15	0.44	0	17,19,21	0.39	0
9	NAG	R	2	9	14,14,15	0.21	0	17,19,21	0.58	0
9	BMA	R	3	9	11,11,12	0.76	0	15,15,17	0.64	0
9	MAN	R	4	9	11,11,12	0.55	0	15,15,17	0.98	1 (6%)
7	NAG	S	1	4,7	14,14,15	1.05	1 (7%)	17,19,21	1.02	2 (11%)
7	MAN	S	10	7	11,11,12	0.85	1 (9%)	15,15,17	0.82	1 (6%)
7	NAG	S	2	7	14,14,15	0.79	1 (7%)	17,19,21	0.51	0
7	BMA	S	3	7	11,11,12	0.81	0	15,15,17	0.92	0
7	MAN	S	4	7	11,11,12	0.51	0	15,15,17	0.97	2 (13%)
7	MAN	S	5	7	11,11,12	0.70	0	15,15,17	0.89	1 (6%)
7	MAN	S	6	7	11,11,12	0.56	0	15,15,17	1.05	2 (13%)
7	MAN	S	7	7	11,11,12	0.97	1 (9%)	15,15,17	0.85	1 (6%)
7	MAN	S	8	7	11,11,12	0.61	0	15,15,17	1.04	2 (13%)
7	MAN	S	9	7	11,11,12	0.76	0	15,15,17	0.88	0
10	NAG	T	1	4,10	14,14,15	0.67	1 (7%)	17,19,21	0.48	0
10	NAG	T	2	10	14,14,15	0.20	0	17,19,21	0.38	0
10	BMA	T	3	10	11,11,12	0.56	0	15,15,17	0.66	0
8	NAG	U	1	8,4	14,14,15	0.41	0	17,19,21	0.75	1 (5%)
8	NAG	U	2	8	14,14,15	0.38	0	17,19,21	0.50	0
8	BMA	U	3	8	11,11,12	0.98	1 (9%)	15,15,17	0.84	0
8	MAN	U	4	8	11,11,12	0.86	1 (9%)	15,15,17	1.11	2 (13%)
8	MAN	U	5	8	11,11,12	0.88	0	15,15,17	0.90	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	MAN	U	6	8	11,11,12	0.74	0	15,15,17	1.04	2 (13%)
6	NAG	V	1	6,2	14,14,15	0.55	0	17,19,21	0.69	0
6	NAG	V	2	6	14,14,15	0.23	0	17,19,21	0.37	0
9	NAG	W	1	9,2	14,14,15	0.74	1 (7%)	17,19,21	0.76	1 (5%)
9	NAG	W	2	9	14,14,15	0.25	0	17,19,21	0.58	0
9	BMA	W	3	9	11,11,12	0.83	0	15,15,17	0.74	0
9	MAN	W	4	9	11,11,12	0.76	0	15,15,17	1.16	2 (13%)
7	NAG	Z	1	4,7	14,14,15	0.37	0	17,19,21	0.55	0
7	MAN	Z	10	7	11,11,12	0.73	0	15,15,17	0.86	0
7	NAG	Z	2	7	14,14,15	0.64	1 (7%)	17,19,21	0.53	0
7	BMA	Z	3	7	11,11,12	0.71	0	15,15,17	0.79	0
7	MAN	Z	4	7	11,11,12	0.41	0	15,15,17	1.16	1 (6%)
7	MAN	Z	5	7	11,11,12	0.57	0	15,15,17	1.14	1 (6%)
7	MAN	Z	6	7	11,11,12	0.80	0	15,15,17	1.03	2 (13%)
7	MAN	Z	7	7	11,11,12	0.57	0	15,15,17	1.15	2 (13%)
7	MAN	Z	8	7	11,11,12	0.48	0	15,15,17	1.22	2 (13%)
7	MAN	Z	9	7	11,11,12	0.66	0	15,15,17	1.00	2 (13%)
10	NAG	a	1	4,10	14,14,15	0.36	0	17,19,21	0.69	1 (5%)
10	NAG	a	2	10	14,14,15	0.29	0	17,19,21	0.41	0
10	BMA	a	3	10	11,11,12	0.55	0	15,15,17	0.65	0
5	NAG	b	1	4,5	14,14,15	0.39	0	17,19,21	0.89	1 (5%)
5	NAG	b	2	5	14,14,15	0.22	0	17,19,21	0.58	0
5	BMA	b	3	5	11,11,12	0.86	0	15,15,17	0.91	1 (6%)
5	MAN	b	4	5	11,11,12	0.78	0	15,15,17	1.22	2 (13%)
5	MAN	b	5	5	11,11,12	0.78	0	15,15,17	0.98	2 (13%)

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
5	NAG	F	1	2,5	-	0/6/23/26	0/1/1/1
5	NAG	F	2	5	-	2/6/23/26	0/1/1/1
5	BMA	F	3	5	-	2/2/19/22	0/1/1/1
5	MAN	F	4	5	-	1/2/19/22	0/1/1/1
5	MAN	F	5	5	-	0/2/19/22	0/1/1/1
6	NAG	I	1	6,4	-	0/6/23/26	0/1/1/1
6	NAG	I	2	6	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	M	1	6,4	-	0/6/23/26	0/1/1/1
6	NAG	M	2	6	-	2/6/23/26	0/1/1/1
7	NAG	N	1	4,7	-	2/6/23/26	0/1/1/1
7	MAN	N	10	7	-	0/2/19/22	0/1/1/1
7	NAG	N	2	7	-	2/6/23/26	0/1/1/1
7	BMA	N	3	7	-	0/2/19/22	0/1/1/1
7	MAN	N	4	7	-	0/2/19/22	0/1/1/1
7	MAN	N	5	7	-	2/2/19/22	0/1/1/1
7	MAN	N	6	7	-	0/2/19/22	0/1/1/1
7	MAN	N	7	7	-	0/2/19/22	0/1/1/1
7	MAN	N	8	7	-	2/2/19/22	0/1/1/1
7	MAN	N	9	7	-	2/2/19/22	0/1/1/1
8	NAG	O	1	8,4	-	0/6/23/26	0/1/1/1
8	NAG	O	2	8	-	1/6/23/26	0/1/1/1
8	BMA	O	3	8	-	0/2/19/22	0/1/1/1
8	MAN	O	4	8	-	1/2/19/22	0/1/1/1
8	MAN	O	5	8	-	0/2/19/22	0/1/1/1
8	MAN	O	6	8	-	0/2/19/22	0/1/1/1
6	NAG	P	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	P	2	6	-	2/6/23/26	0/1/1/1
6	NAG	Q	1	6,2	-	0/6/23/26	0/1/1/1
6	NAG	Q	2	6	-	2/6/23/26	0/1/1/1
9	NAG	R	1	9,2	-	0/6/23/26	0/1/1/1
9	NAG	R	2	9	-	2/6/23/26	0/1/1/1
9	BMA	R	3	9	-	0/2/19/22	0/1/1/1
9	MAN	R	4	9	-	1/2/19/22	0/1/1/1
7	NAG	S	1	4,7	-	2/6/23/26	0/1/1/1
7	MAN	S	10	7	-	0/2/19/22	0/1/1/1
7	NAG	S	2	7	-	0/6/23/26	0/1/1/1
7	BMA	S	3	7	-	0/2/19/22	0/1/1/1
7	MAN	S	4	7	-	0/2/19/22	0/1/1/1
7	MAN	S	5	7	-	2/2/19/22	0/1/1/1
7	MAN	S	6	7	-	0/2/19/22	0/1/1/1
7	MAN	S	7	7	-	0/2/19/22	0/1/1/1
7	MAN	S	8	7	-	2/2/19/22	0/1/1/1
7	MAN	S	9	7	-	0/2/19/22	0/1/1/1
10	NAG	T	1	4,10	-	2/6/23/26	0/1/1/1
10	NAG	T	2	10	-	2/6/23/26	0/1/1/1
10	BMA	T	3	10	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	U	1	8,4	-	2/6/23/26	0/1/1/1
8	NAG	U	2	8	-	2/6/23/26	0/1/1/1
8	BMA	U	3	8	-	0/2/19/22	0/1/1/1
8	MAN	U	4	8	-	1/2/19/22	0/1/1/1
8	MAN	U	5	8	-	0/2/19/22	0/1/1/1
8	MAN	U	6	8	-	0/2/19/22	0/1/1/1
6	NAG	V	1	6,2	-	0/6/23/26	0/1/1/1
6	NAG	V	2	6	-	2/6/23/26	0/1/1/1
9	NAG	W	1	9,2	-	0/6/23/26	0/1/1/1
9	NAG	W	2	9	-	1/6/23/26	0/1/1/1
9	BMA	W	3	9	-	1/2/19/22	0/1/1/1
9	MAN	W	4	9	-	2/2/19/22	0/1/1/1
7	NAG	Z	1	4,7	-	2/6/23/26	0/1/1/1
7	MAN	Z	10	7	-	0/2/19/22	0/1/1/1
7	NAG	Z	2	7	-	2/6/23/26	0/1/1/1
7	BMA	Z	3	7	-	0/2/19/22	0/1/1/1
7	MAN	Z	4	7	-	0/2/19/22	0/1/1/1
7	MAN	Z	5	7	-	2/2/19/22	0/1/1/1
7	MAN	Z	6	7	-	0/2/19/22	0/1/1/1
7	MAN	Z	7	7	-	0/2/19/22	0/1/1/1
7	MAN	Z	8	7	-	2/2/19/22	0/1/1/1
7	MAN	Z	9	7	-	2/2/19/22	0/1/1/1
10	NAG	a	1	4,10	-	2/6/23/26	0/1/1/1
10	NAG	a	2	10	-	2/6/23/26	0/1/1/1
10	BMA	a	3	10	-	0/2/19/22	0/1/1/1
5	NAG	b	1	4,5	-	0/6/23/26	0/1/1/1
5	NAG	b	2	5	-	2/6/23/26	0/1/1/1
5	BMA	b	3	5	-	0/2/19/22	0/1/1/1
5	MAN	b	4	5	-	2/2/19/22	0/1/1/1
5	MAN	b	5	5	-	0/2/19/22	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	S	1	NAG	O5-C1	-3.87	1.37	1.43
7	S	2	NAG	O5-C1	-2.91	1.38	1.43
9	W	1	NAG	O5-C1	-2.63	1.39	1.43
8	U	4	MAN	C1-C2	2.36	1.57	1.52
7	Z	2	NAG	O5-C1	-2.34	1.39	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	P	1	NAG	O5-C1	-2.32	1.39	1.43
7	N	2	NAG	O5-C1	-2.30	1.39	1.43
10	T	1	NAG	C1-C2	2.30	1.55	1.52
8	U	3	BMA	C4-C3	2.24	1.58	1.52
8	O	4	MAN	C1-C2	2.15	1.57	1.52
7	S	7	MAN	O5-C1	-2.11	1.40	1.43
6	I	1	NAG	O5-C1	-2.10	1.40	1.43
7	S	10	MAN	O5-C1	-2.04	1.40	1.43

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	4	MAN	C1-O5-C5	4.05	117.61	112.19
7	N	4	MAN	C1-O5-C5	3.85	117.35	112.19
7	Z	4	MAN	C1-O5-C5	3.84	117.33	112.19
5	b	4	MAN	C1-O5-C5	3.61	117.02	112.19
9	W	4	MAN	C1-O5-C5	3.50	116.88	112.19
8	O	4	MAN	C1-O5-C5	3.30	116.61	112.19
7	Z	7	MAN	C1-O5-C5	3.12	116.37	112.19
7	N	7	MAN	C1-O5-C5	3.11	116.35	112.19
8	U	6	MAN	C1-O5-C5	3.00	116.20	112.19
8	U	4	MAN	C1-O5-C5	2.87	116.03	112.19
5	F	5	MAN	C1-O5-C5	2.82	115.97	112.19
9	R	4	MAN	C1-O5-C5	2.64	115.72	112.19
7	N	5	MAN	C1-O5-C5	2.63	115.72	112.19
8	O	5	MAN	C1-O5-C5	2.63	115.71	112.19
7	Z	5	MAN	C1-O5-C5	2.62	115.70	112.19
7	S	4	MAN	C1-O5-C5	2.59	115.66	112.19
5	b	5	MAN	C1-O5-C5	2.56	115.62	112.19
7	N	8	MAN	C1-O5-C5	2.55	115.61	112.19
8	O	6	MAN	C1-O5-C5	2.55	115.60	112.19
7	Z	8	MAN	C1-O5-C5	2.54	115.59	112.19
7	S	6	MAN	C1-O5-C5	2.53	115.58	112.19
7	Z	6	MAN	C1-O5-C5	2.49	115.53	112.19
9	W	1	NAG	C1-O5-C5	2.48	115.52	112.19
7	N	6	MAN	C1-O5-C5	2.47	115.50	112.19
8	U	4	MAN	O2-C2-C3	-2.45	105.07	110.15
5	b	4	MAN	O2-C2-C3	-2.43	105.11	110.15
7	S	1	NAG	C1-O5-C5	-2.39	108.98	112.19
7	S	8	MAN	O2-C2-C3	-2.37	105.24	110.15
7	N	9	MAN	O2-C2-C3	-2.34	105.29	110.15
7	Z	9	MAN	O2-C2-C3	-2.33	105.32	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	N	9	MAN	C1-O5-C5	2.31	115.28	112.19
8	U	1	NAG	C1-O5-C5	2.29	115.26	112.19
7	Z	9	MAN	C1-O5-C5	2.29	115.26	112.19
10	a	1	NAG	C1-O5-C5	2.29	115.25	112.19
8	O	5	MAN	O2-C2-C3	-2.27	105.44	110.15
8	O	6	MAN	O2-C2-C3	-2.25	105.48	110.15
8	O	4	MAN	O2-C2-C3	-2.22	105.56	110.15
7	N	8	MAN	O2-C2-C3	-2.19	105.61	110.15
7	Z	8	MAN	O2-C2-C3	-2.19	105.61	110.15
7	S	8	MAN	C1-O5-C5	2.19	115.12	112.19
7	Z	6	MAN	O2-C2-C3	-2.18	105.63	110.15
7	N	6	MAN	O2-C2-C3	-2.17	105.66	110.15
7	S	10	MAN	O2-C2-C3	-2.16	105.67	110.15
8	U	6	MAN	O2-C2-C3	-2.16	105.67	110.15
7	S	5	MAN	O2-C2-C3	-2.16	105.67	110.15
8	O	3	BMA	C1-C2-C3	2.16	112.78	109.64
7	Z	7	MAN	O2-C2-C3	-2.15	105.71	110.15
7	N	7	MAN	O2-C2-C3	-2.14	105.72	110.15
5	F	5	MAN	O2-C2-C3	-2.14	105.72	110.15
5	F	4	MAN	O2-C2-C3	-2.11	105.79	110.15
7	S	1	NAG	C3-C4-C5	2.10	114.05	110.23
5	b	3	BMA	C1-O5-C5	2.10	115.00	112.19
8	U	5	MAN	O2-C2-C3	-2.09	105.83	110.15
7	S	4	MAN	O2-C2-C3	-2.08	105.85	110.15
5	b	5	MAN	O2-C2-C3	-2.07	105.87	110.15
7	S	6	MAN	O2-C2-C3	-2.05	105.91	110.15
7	S	7	MAN	O2-C2-C3	-2.04	105.93	110.15
9	W	4	MAN	O2-C2-C3	-2.03	105.94	110.15
5	b	1	NAG	C3-C4-C5	2.01	113.88	110.23

There are no chirality outliers.

All (67) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	N	1	NAG	O5-C5-C6-O6
7	Z	1	NAG	O5-C5-C6-O6
7	N	8	MAN	O5-C5-C6-O6
7	Z	8	MAN	O5-C5-C6-O6
10	a	2	NAG	O5-C5-C6-O6
10	T	2	NAG	O5-C5-C6-O6
7	N	1	NAG	C4-C5-C6-O6
7	Z	1	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	N	8	MAN	C4-C5-C6-O6
7	Z	8	MAN	C4-C5-C6-O6
8	U	2	NAG	O5-C5-C6-O6
7	S	8	MAN	O5-C5-C6-O6
7	S	5	MAN	O5-C5-C6-O6
7	S	5	MAN	C4-C5-C6-O6
10	T	1	NAG	C4-C5-C6-O6
7	S	1	NAG	O5-C5-C6-O6
10	T	2	NAG	C4-C5-C6-O6
10	a	2	NAG	C4-C5-C6-O6
7	N	5	MAN	O5-C5-C6-O6
7	Z	5	MAN	O5-C5-C6-O6
8	U	2	NAG	C4-C5-C6-O6
7	N	9	MAN	O5-C5-C6-O6
7	Z	9	MAN	O5-C5-C6-O6
7	S	8	MAN	C4-C5-C6-O6
10	T	1	NAG	O5-C5-C6-O6
5	F	2	NAG	O5-C5-C6-O6
6	I	2	NAG	O5-C5-C6-O6
5	b	2	NAG	O5-C5-C6-O6
6	M	2	NAG	O5-C5-C6-O6
8	U	1	NAG	O5-C5-C6-O6
7	Z	5	MAN	C4-C5-C6-O6
7	N	5	MAN	C4-C5-C6-O6
9	W	4	MAN	O5-C5-C6-O6
6	P	2	NAG	O5-C5-C6-O6
8	U	1	NAG	C4-C5-C6-O6
6	V	2	NAG	O5-C5-C6-O6
6	Q	2	NAG	O5-C5-C6-O6
5	b	4	MAN	O5-C5-C6-O6
7	S	1	NAG	C4-C5-C6-O6
7	Z	9	MAN	C4-C5-C6-O6
7	N	9	MAN	C4-C5-C6-O6
8	O	4	MAN	O5-C5-C6-O6
9	R	4	MAN	O5-C5-C6-O6
8	O	2	NAG	O5-C5-C6-O6
8	U	4	MAN	O5-C5-C6-O6
6	I	2	NAG	C4-C5-C6-O6
10	a	1	NAG	C4-C5-C6-O6
5	F	4	MAN	O5-C5-C6-O6
7	N	2	NAG	C4-C5-C6-O6
7	Z	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
6	M	2	NAG	C4-C5-C6-O6
5	F	2	NAG	C4-C5-C6-O6
5	b	2	NAG	C4-C5-C6-O6
9	W	2	NAG	O5-C5-C6-O6
6	P	2	NAG	C4-C5-C6-O6
10	a	1	NAG	O5-C5-C6-O6
6	V	2	NAG	C4-C5-C6-O6
7	N	2	NAG	O5-C5-C6-O6
7	Z	2	NAG	O5-C5-C6-O6
5	F	3	BMA	C4-C5-C6-O6
6	Q	2	NAG	C4-C5-C6-O6
9	R	2	NAG	O5-C5-C6-O6
5	F	3	BMA	O5-C5-C6-O6
9	W	4	MAN	C4-C5-C6-O6
9	R	2	NAG	C1-C2-N2-C7
9	W	3	BMA	C4-C5-C6-O6
5	b	4	MAN	C4-C5-C6-O6

There are no ring outliers.

32 monomers are involved in 57 short contacts:

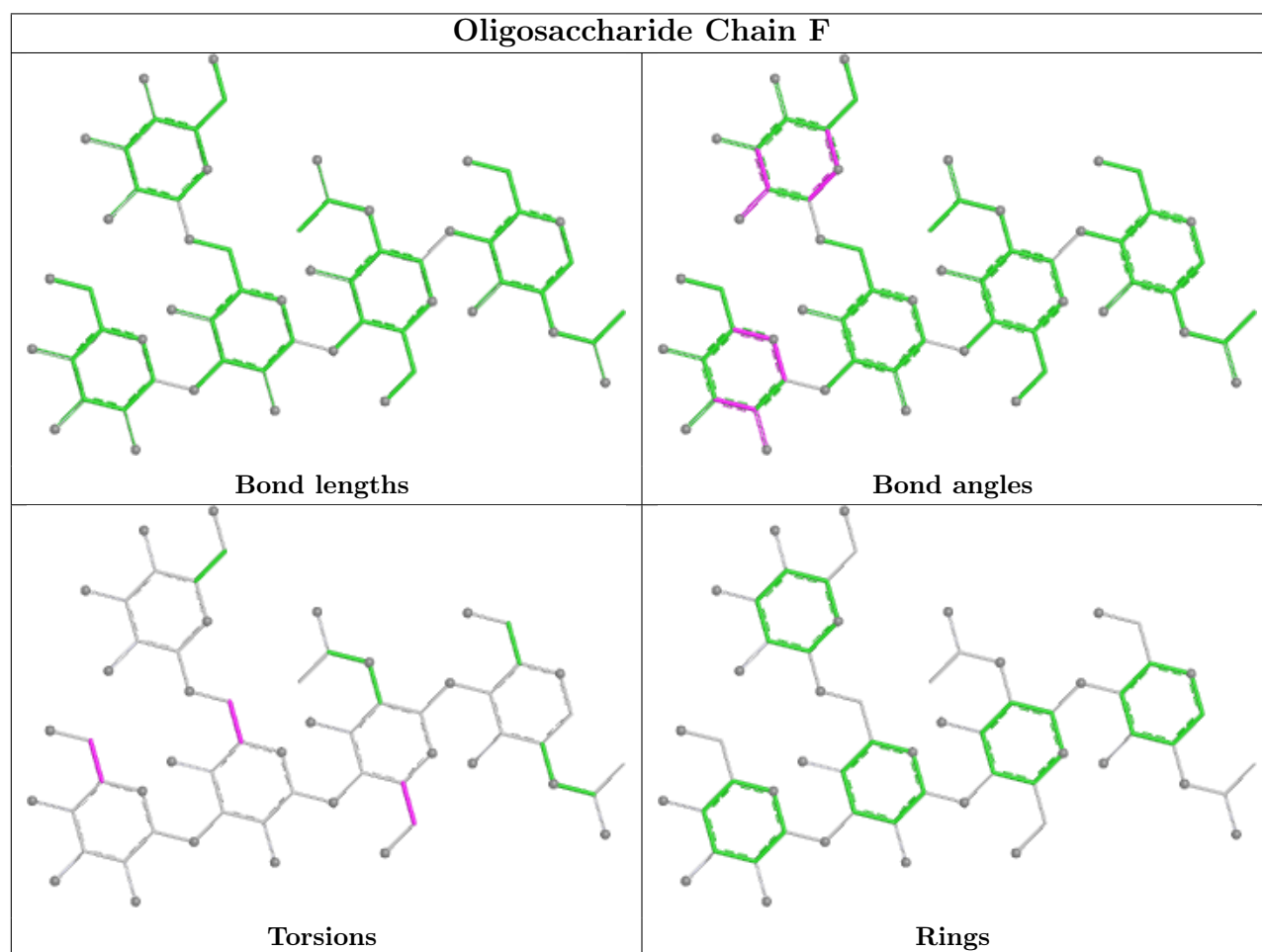
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	Q	2	NAG	1	0
5	b	2	NAG	2	0
8	U	5	MAN	1	0
7	N	9	MAN	3	0
10	a	1	NAG	1	0
6	M	2	NAG	1	0
7	N	6	MAN	1	0
9	W	1	NAG	2	0
7	S	2	NAG	1	0
8	U	1	NAG	2	0
10	T	1	NAG	1	0
5	F	3	BMA	1	0
9	W	2	NAG	3	0
9	R	2	NAG	1	0
7	Z	5	MAN	1	0
8	O	1	NAG	7	0
8	U	3	BMA	1	0
7	S	1	NAG	1	0
8	U	2	NAG	2	0
7	S	3	BMA	1	0

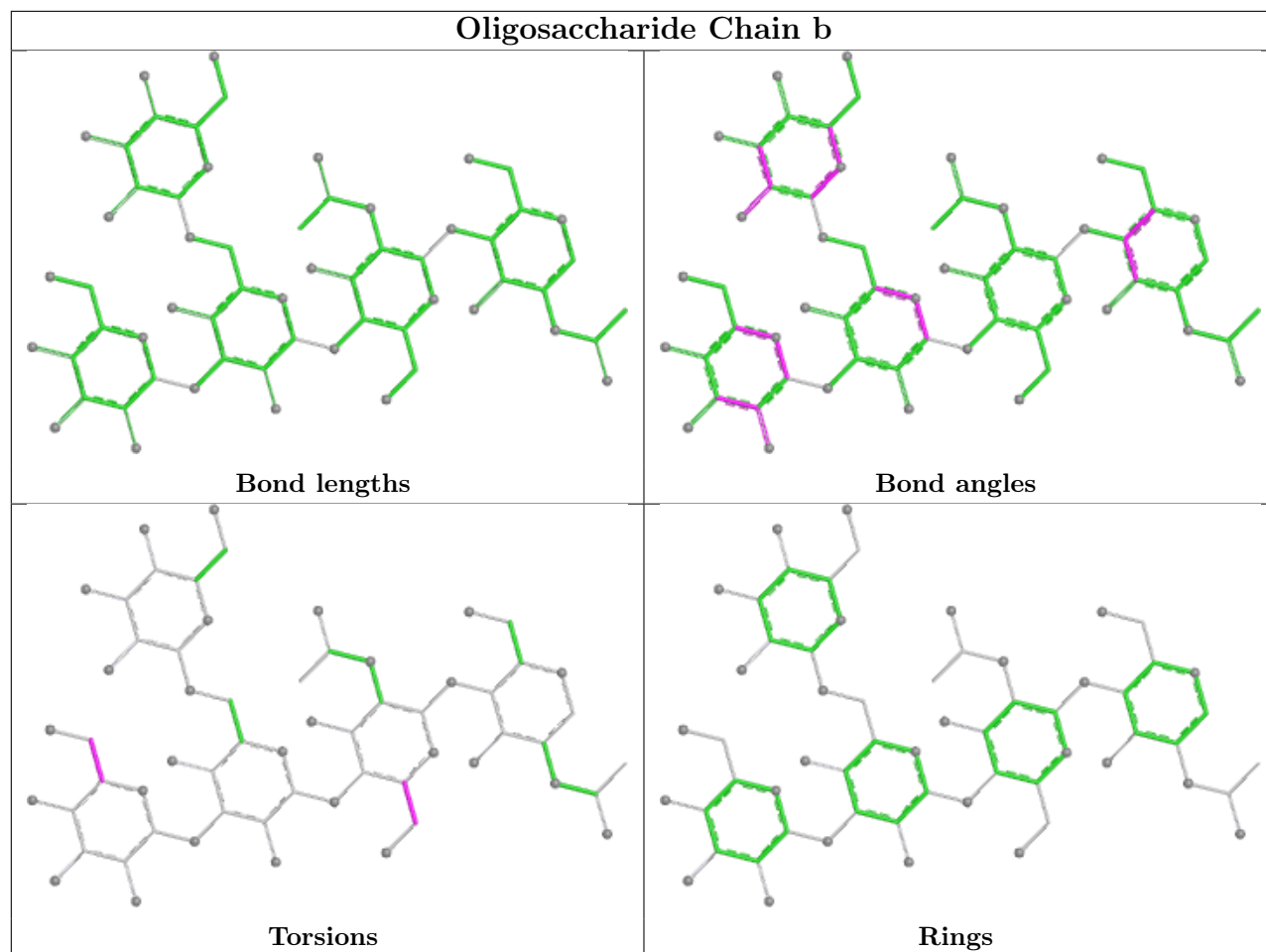
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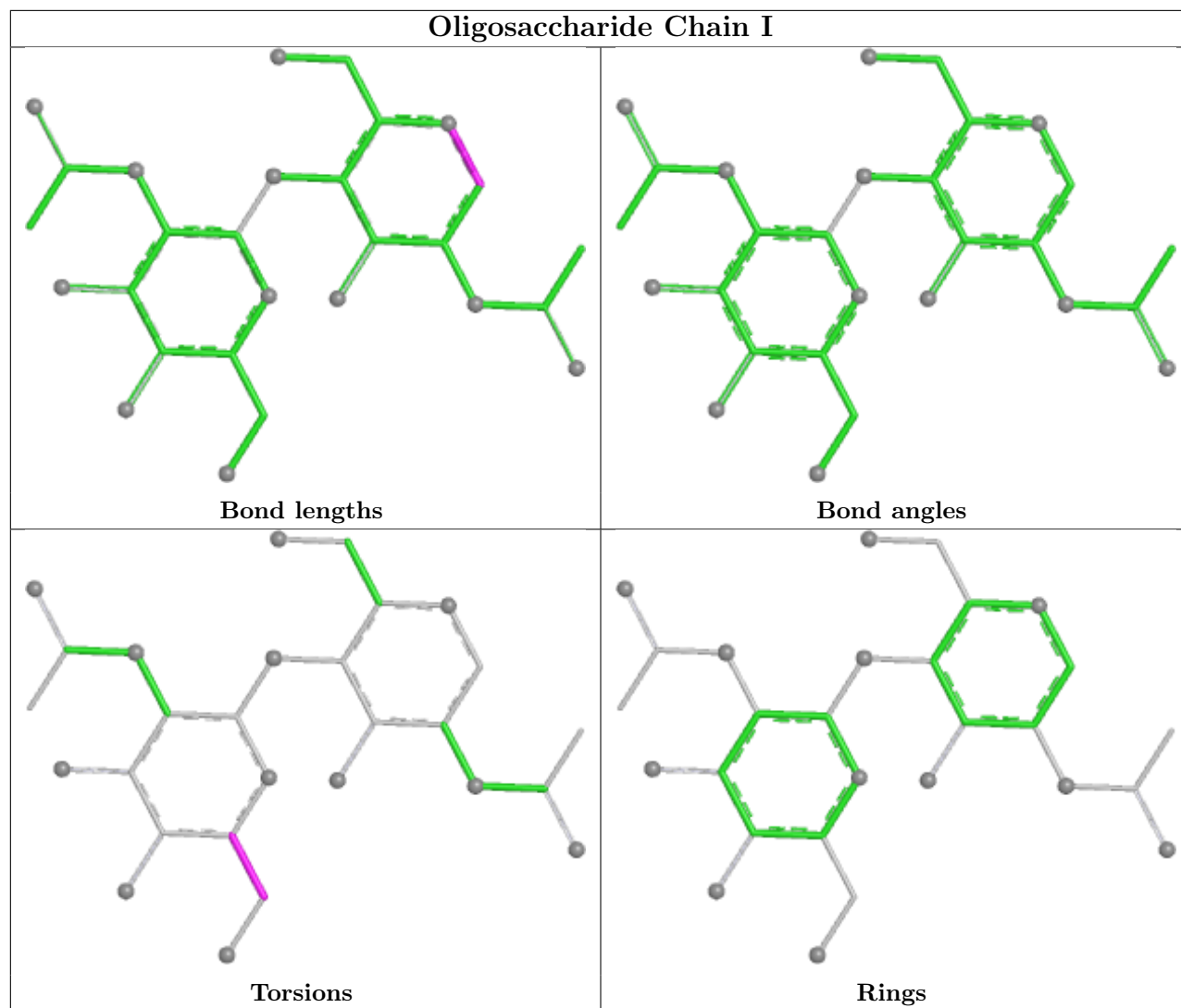
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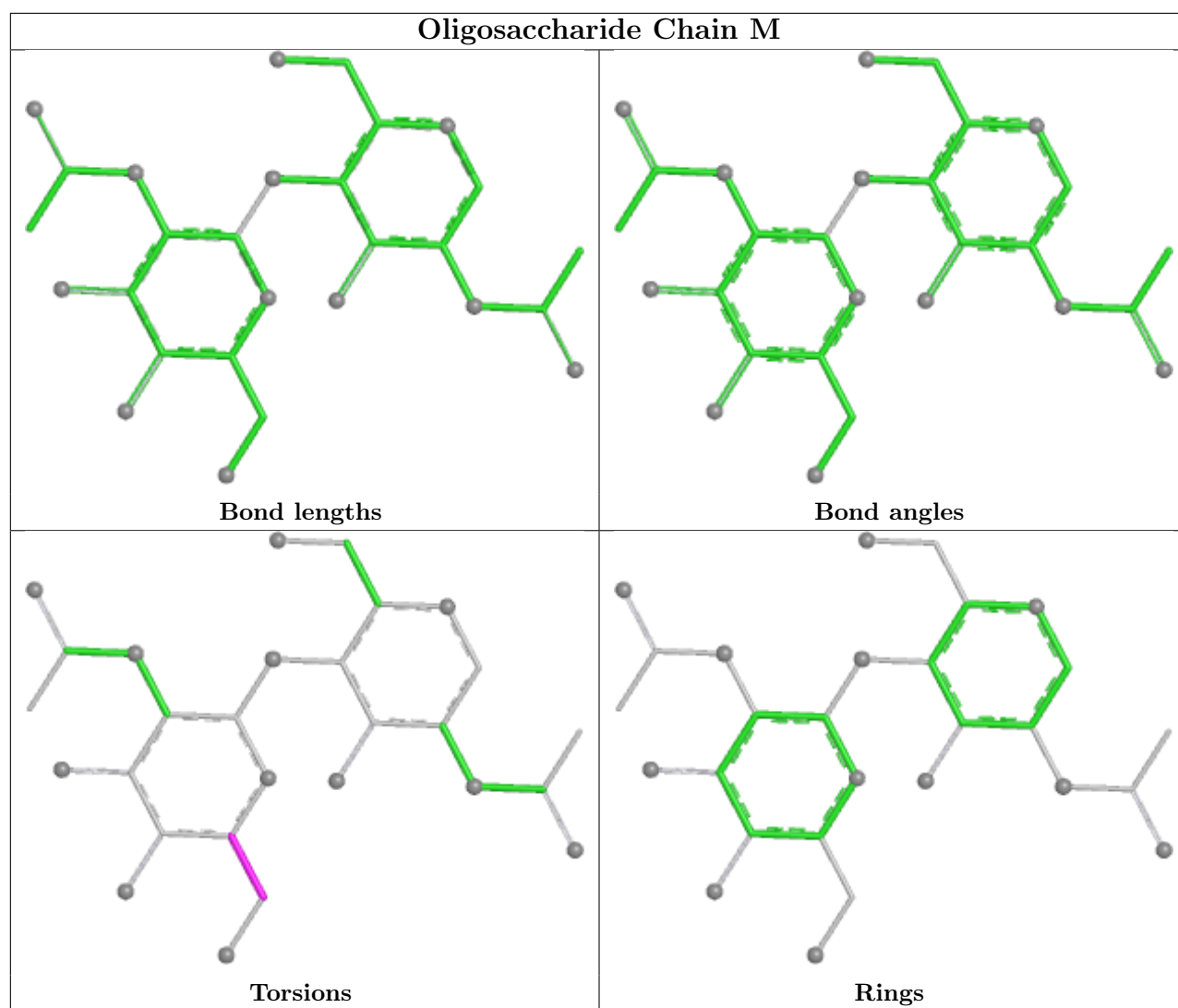
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	S	9	MAN	5	0
7	Z	6	MAN	5	0
6	Q	1	NAG	1	0
7	Z	8	MAN	1	0
7	Z	9	MAN	6	0
7	N	1	NAG	2	0
9	R	1	NAG	1	0
5	F	2	NAG	3	0
6	M	1	NAG	1	0
7	N	2	NAG	1	0
5	F	1	NAG	1	0
8	O	2	NAG	2	0

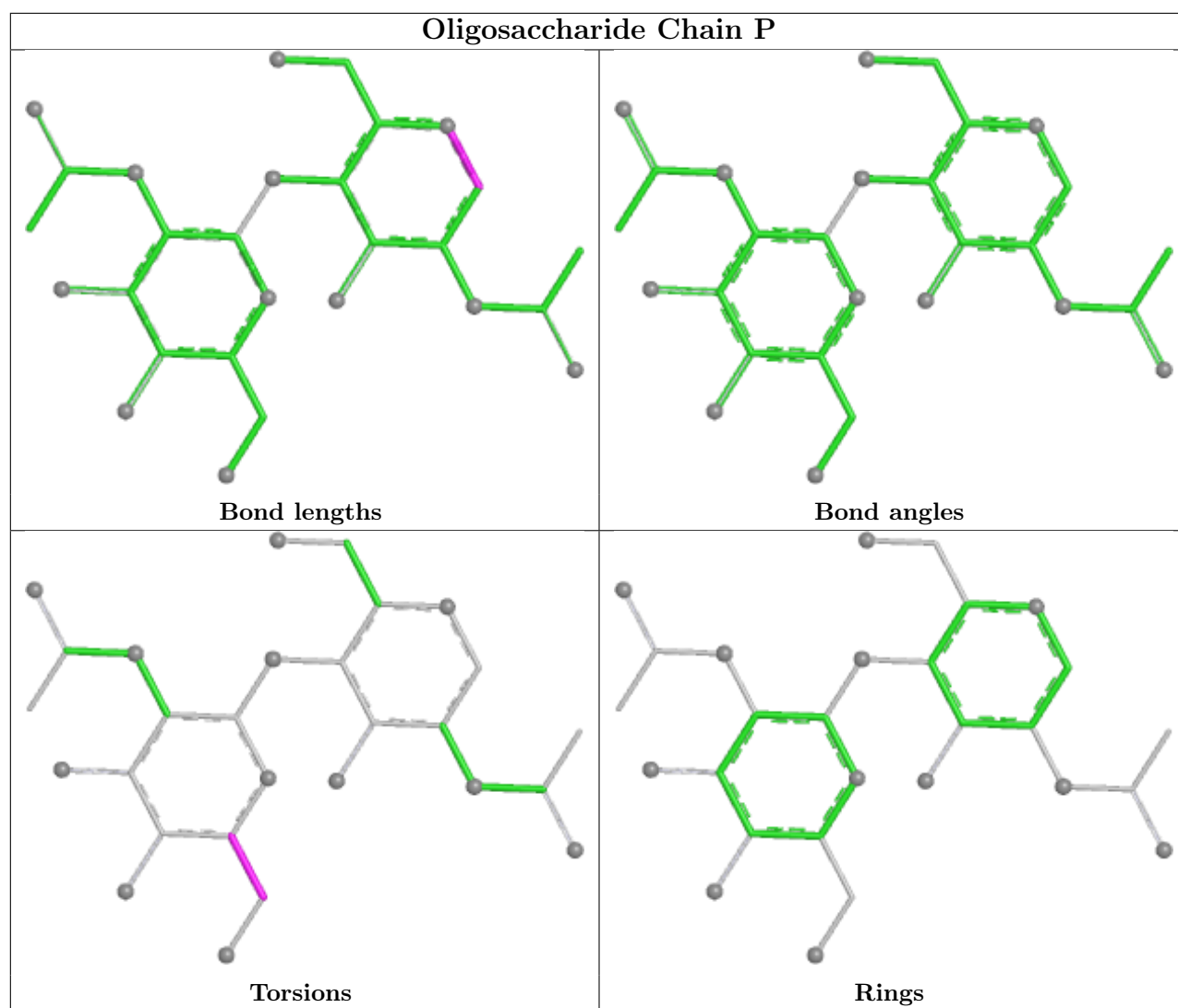
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

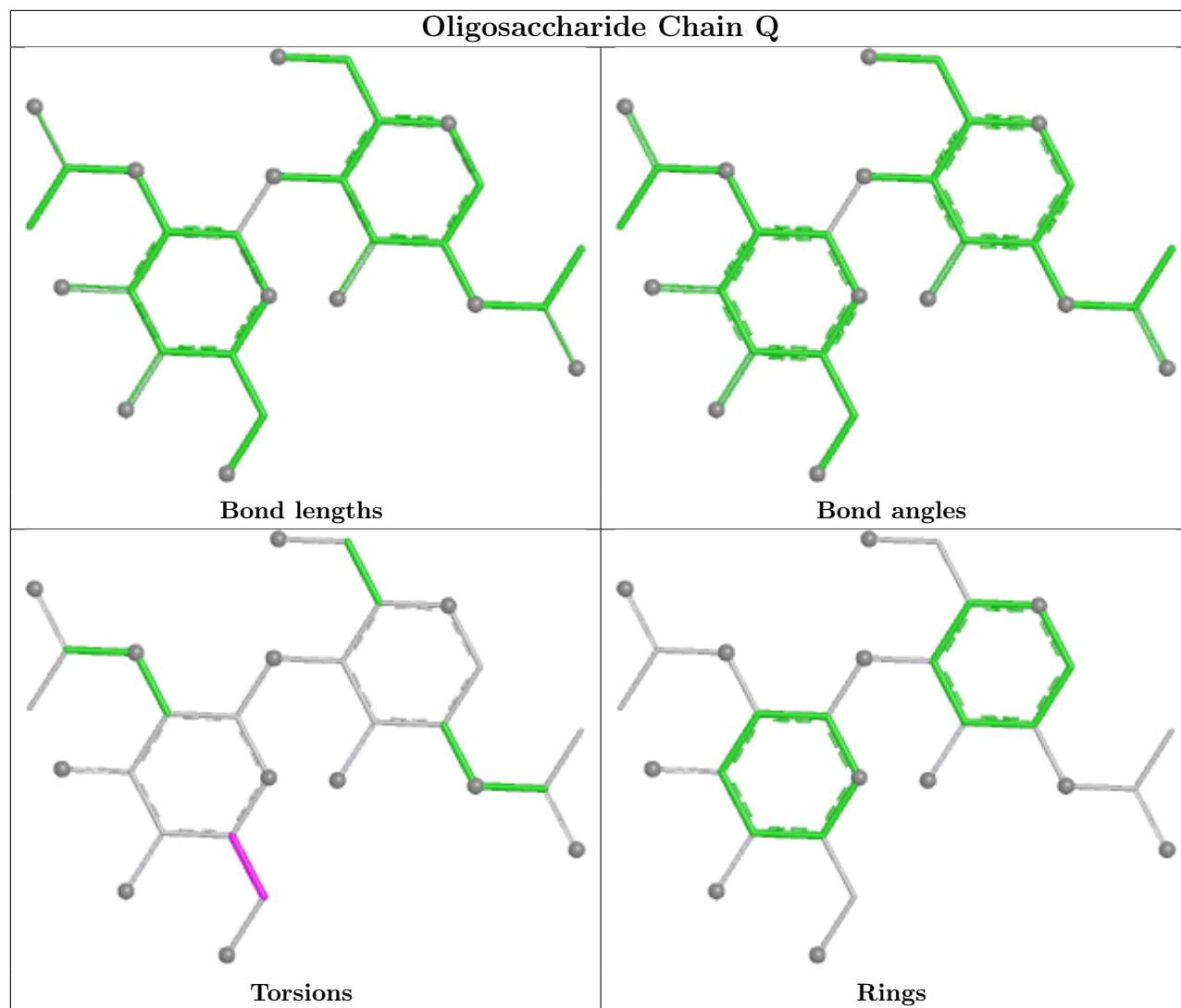


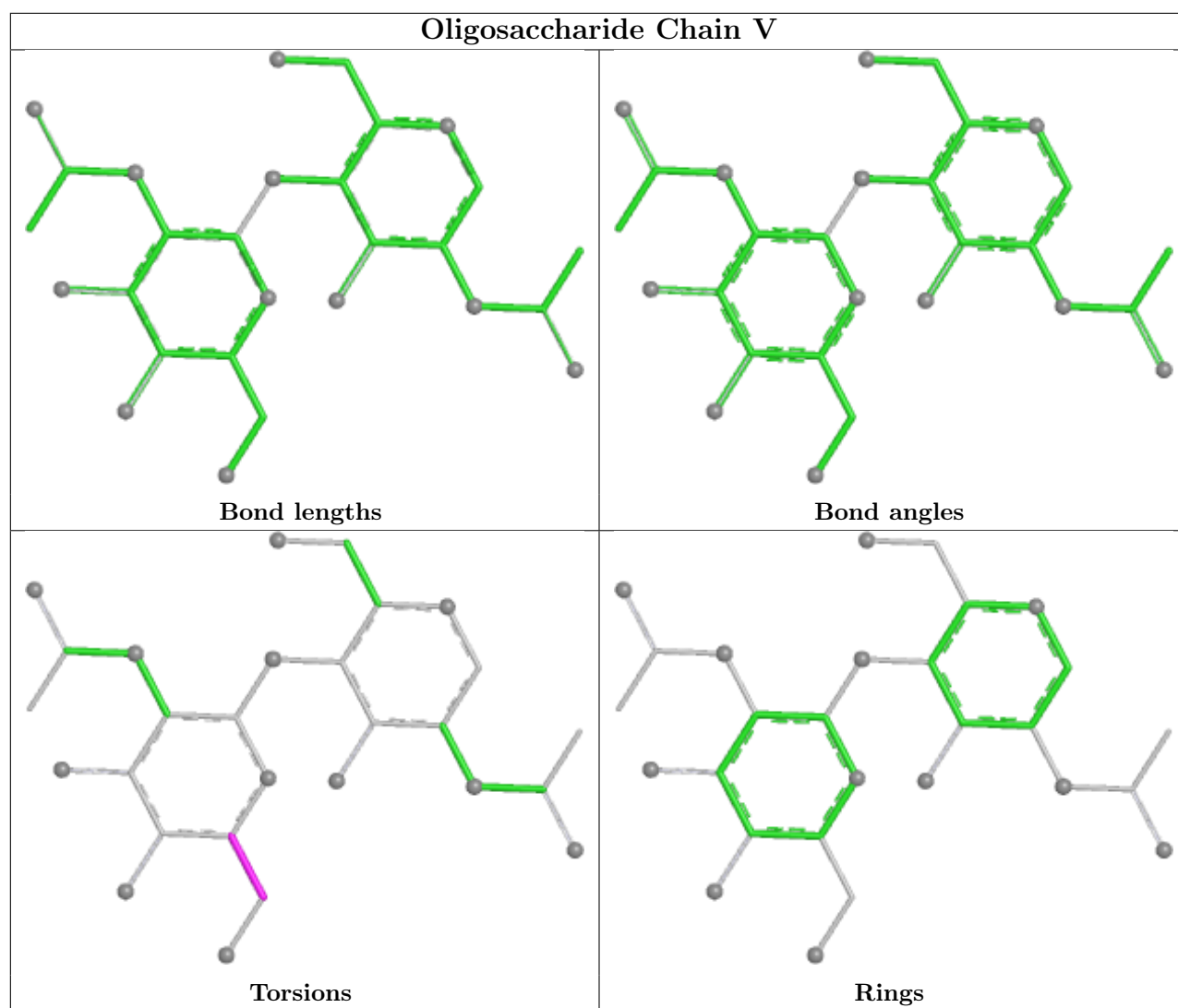


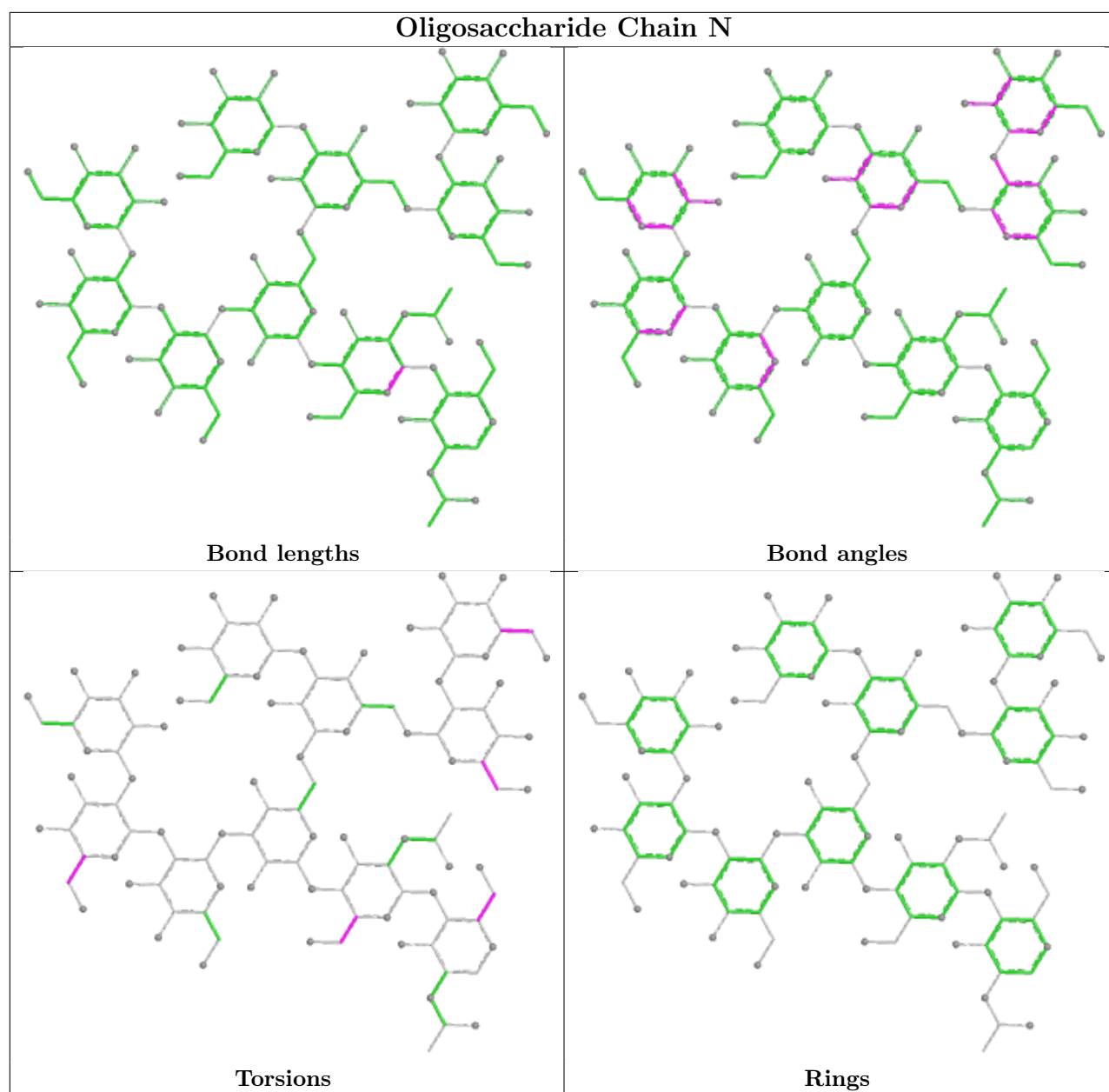


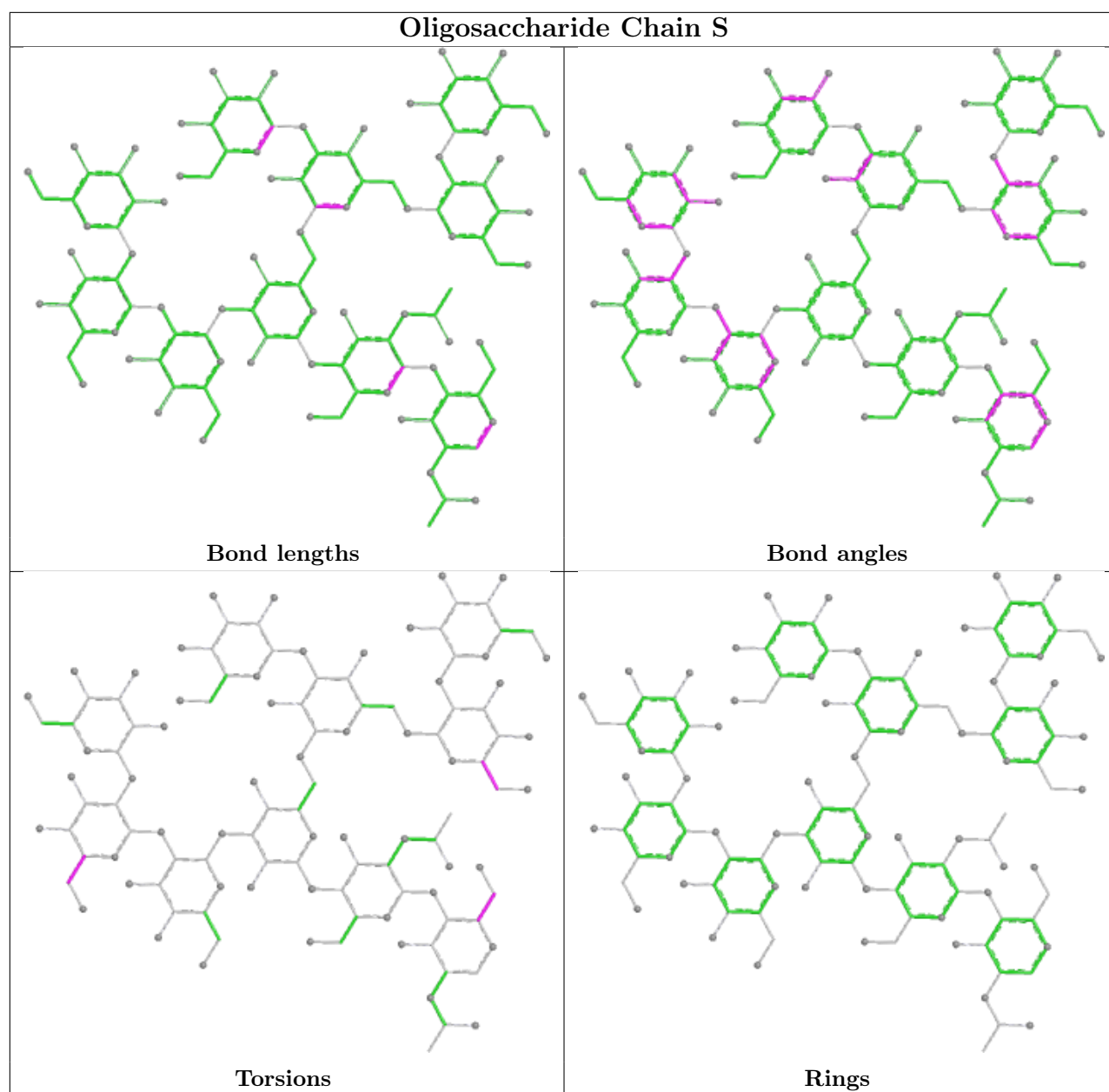


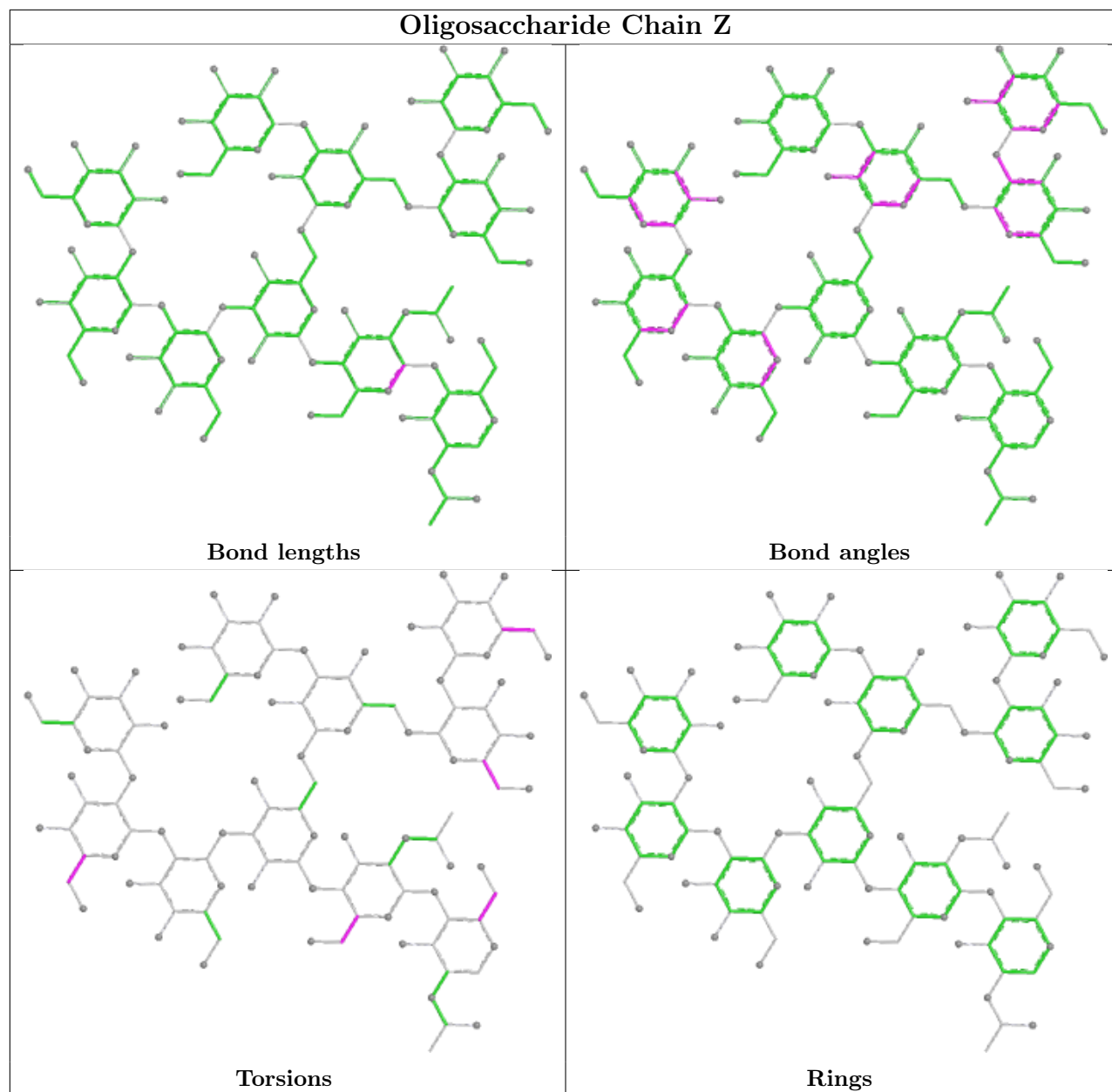


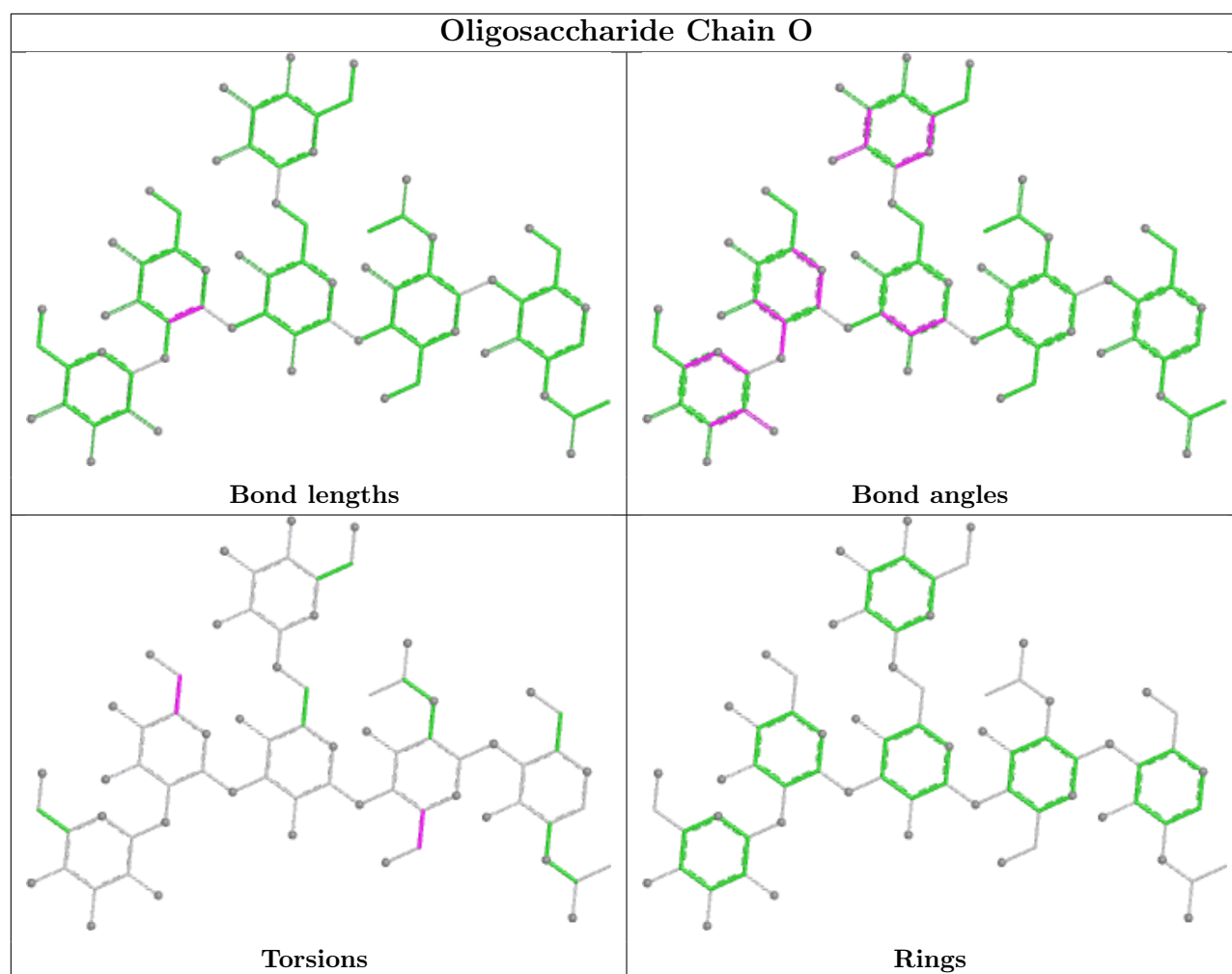




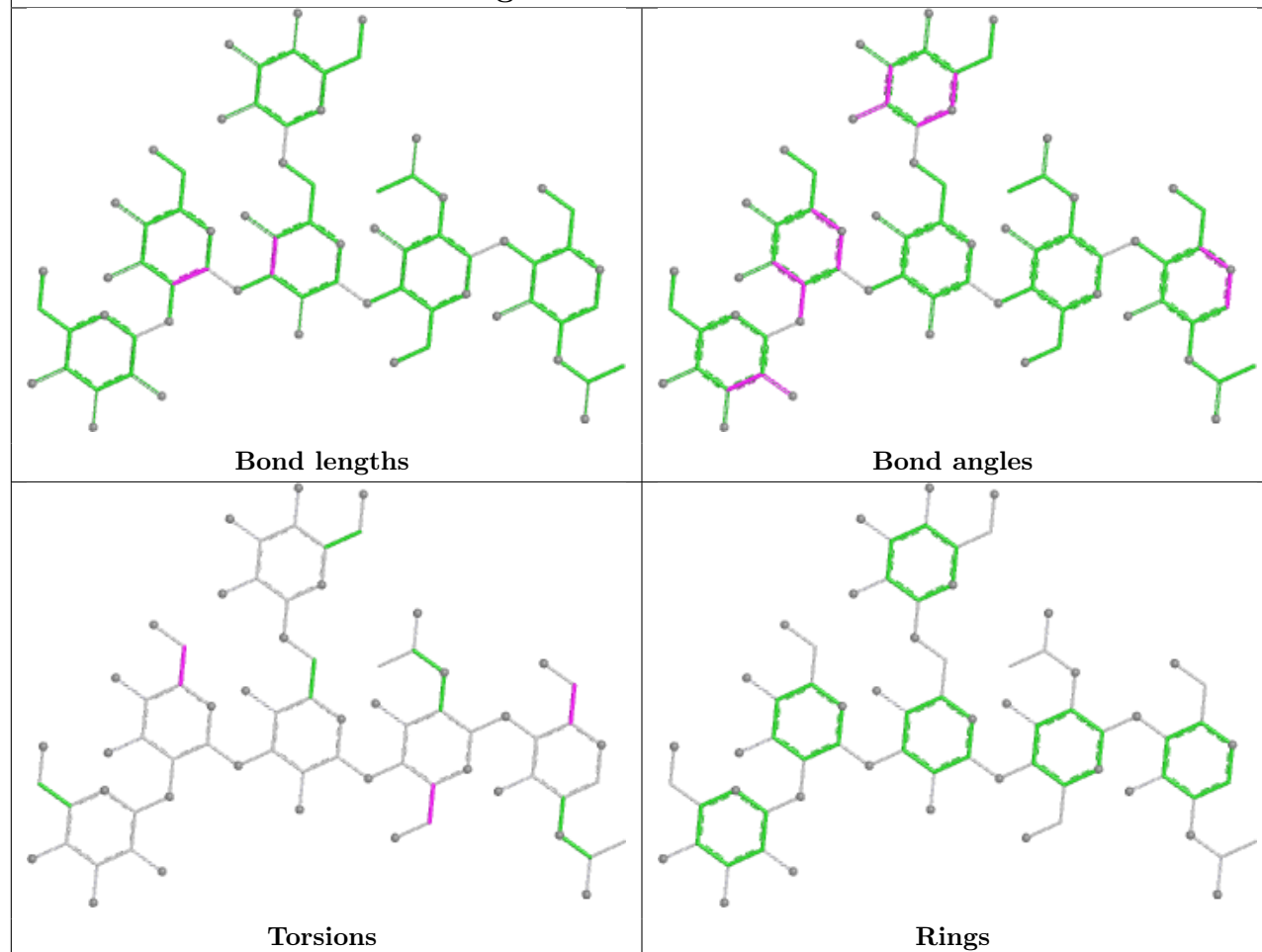




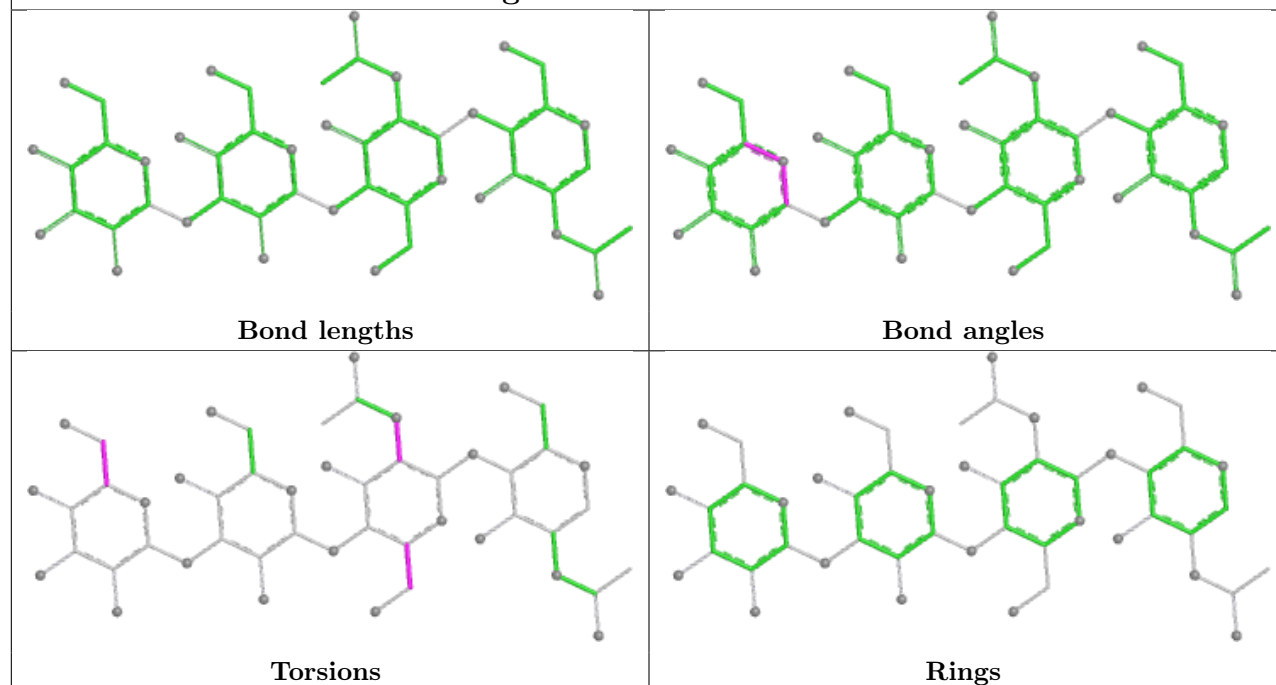




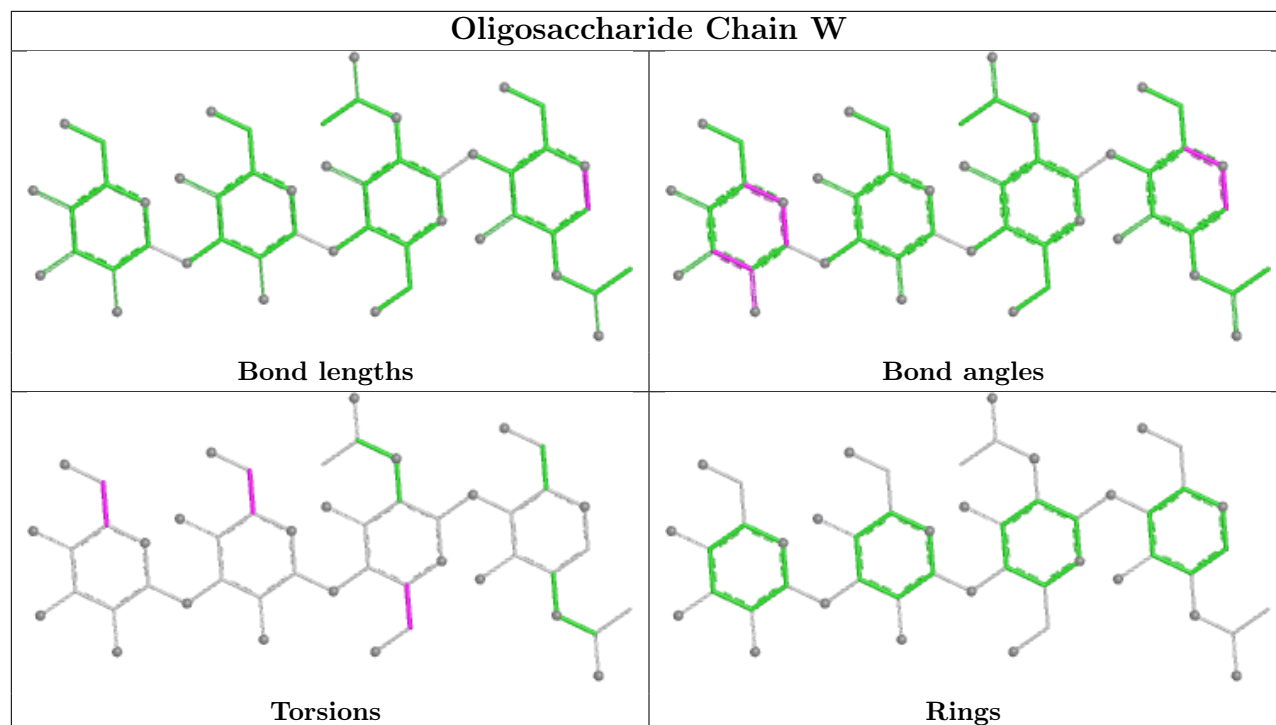
Oligosaccharide Chain U



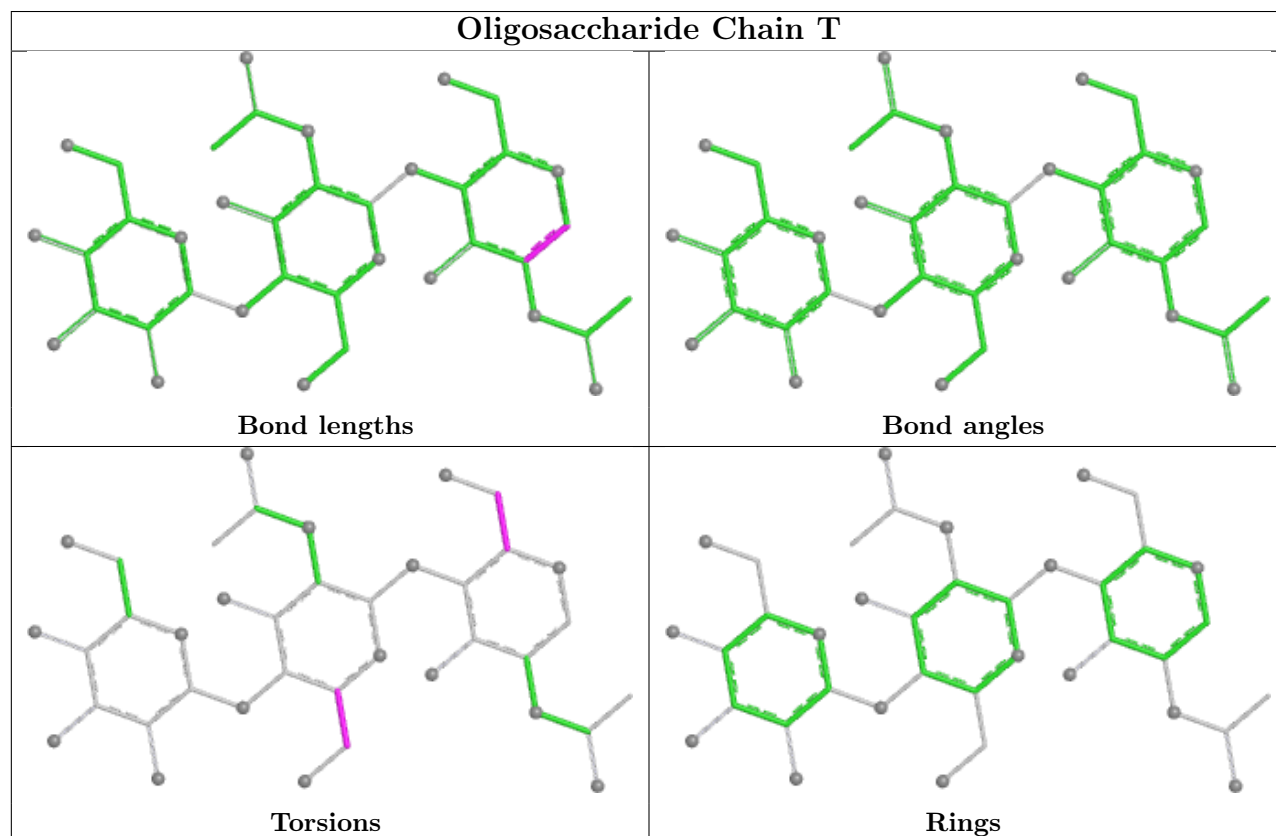
Oligosaccharide Chain R

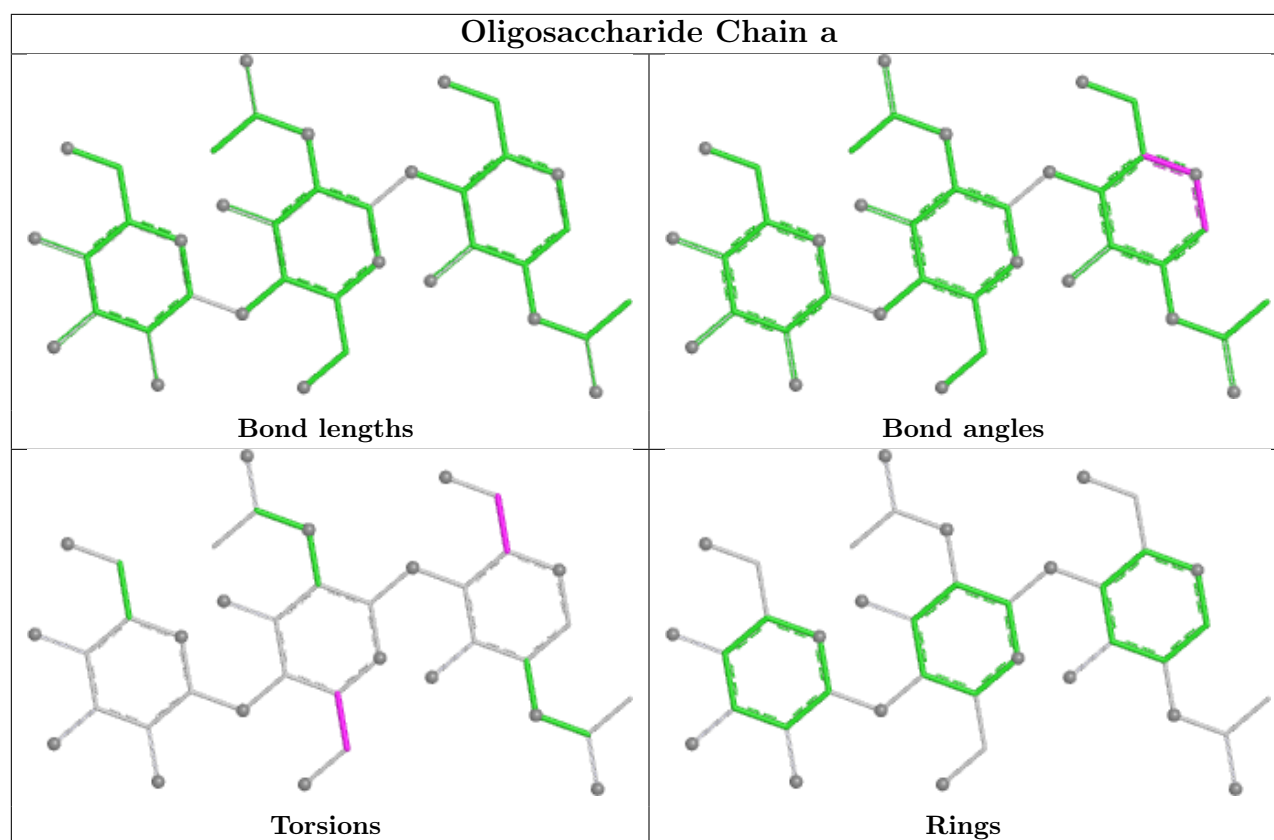


Oligosaccharide Chain W



Oligosaccharide Chain T





5.6 Ligand geometry [i](#)

6 ligands 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	$\# Z > 2$
11	NAG	J	1042	2	14,14,15	0.61	0	17,19,21	0.75	1 (5%)
11	NAG	A	1000	-	14,14,15	0.52	0	17,19,21	0.50	0
11	NAG	Y	1041	4	14,14,15	1.01	1 (7%)	17,19,21	0.82	1 (5%)
11	NAG	K	1040	4	14,14,15	0.61	0	17,19,21	0.48	0
11	NAG	B	1040	2	14,14,15	0.30	0	17,19,21	0.55	0
11	NAG	Y	1040	4	14,14,15	0.54	0	17,19,21	0.63	1 (5%)

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
11	NAG	J	1042	2	-	1/6/23/26	0/1/1/1
11	NAG	A	1000	-	-	1/6/23/26	0/1/1/1
11	NAG	Y	1041	4	-	1/6/23/26	0/1/1/1
11	NAG	K	1040	4	-	1/6/23/26	0/1/1/1
11	NAG	B	1040	2	-	1/6/23/26	0/1/1/1
11	NAG	Y	1040	4	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	Y	1041	NAG	C1-C2	3.21	1.56	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	Y	1041	NAG	C1-O5-C5	2.69	115.79	112.19
11	J	1042	NAG	C1-O5-C5	2.42	115.43	112.19
11	Y	1040	NAG	C1-O5-C5	2.02	114.89	112.19

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	K	1040	NAG	C3-C2-N2-C7
11	A	1000	NAG	C4-C5-C6-O6
11	B	1040	NAG	C3-C2-N2-C7
11	J	1042	NAG	C3-C2-N2-C7
11	Y	1041	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	Y	1041	NAG	2	0

5.7 Other polymers

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	224/244 (91%)	-0.20	1 (0%) 88 70	94, 133, 173, 216	0
1	D	224/244 (91%)	-0.13	1 (0%) 88 70	102, 155, 201, 272	0
1	H	227/244 (93%)	-0.23	1 (0%) 88 70	71, 121, 164, 195	0
2	B	124/153 (81%)	-0.00	2 (1%) 70 42	76, 115, 180, 236	0
2	J	124/153 (81%)	-0.01	2 (1%) 70 42	67, 113, 178, 255	0
2	X	121/153 (79%)	0.05	0 100 100	74, 119, 177, 325	0
3	C	214/215 (99%)	-0.23	1 (0%) 87 66	91, 144, 203, 230	0
3	E	214/215 (99%)	-0.18	3 (1%) 73 45	99, 167, 215, 277	0
3	L	214/215 (99%)	-0.39	1 (0%) 87 66	80, 122, 209, 246	0
4	G	398/479 (83%)	-0.01	7 (1%) 67 40	71, 167, 231, 277	0
4	K	405/479 (84%)	0.04	8 (1%) 65 37	63, 167, 248, 316	0
4	Y	397/479 (82%)	0.04	7 (1%) 67 40	66, 176, 239, 315	0
All	All	2886/3273 (88%)	-0.09	34 (1%) 76 49	63, 147, 222, 325	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	104	GLY	4.1
4	K	351	LYS	3.9
4	Y	154	LEU	3.9
4	K	352	HIS	3.3
4	K	297	THR	3.1
4	K	211	GLU	3.1
2	B	616	ASN	2.9
4	K	307	ILE	2.8
3	L	213	GLU	2.7
4	Y	70	ALA	2.6
3	E	31	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
4	G	438	PRO	2.6
4	Y	175	LEU	2.6
4	G	439	ILE	2.5
4	K	154	LEU	2.5
3	E	173	TYR	2.4
4	Y	212	PRO	2.3
3	C	12	SER	2.3
2	B	664	ASP	2.3
2	J	602	LEU	2.3
4	G	389	GLY	2.3
1	H	28	ASN	2.3
3	E	96	GLY	2.3
1	A	125	ALA	2.2
4	Y	155	LYS	2.2
4	G	380	GLY	2.2
2	J	652	GLN	2.2
4	G	478	ASN	2.2
4	K	179	LEU	2.1
4	Y	336	ALA	2.1
4	K	442	VAL	2.1
4	Y	123	THR	2.1
4	G	272	ILE	2.0
4	G	162	THR	2.0

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 [i](#)

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($\text{\AA}^2$)	Q<0.9
5	NAG	F	1	14/15	-	-	69,88,120,138	0
5	NAG	F	2	14/15	-	-	106,124,138,153	0
5	BMA	F	3	11/12	-	-	135,148,156,157	11
5	MAN	F	4	11/12	-	-	119,131,141,143	11
5	MAN	F	5	11/12	-	-	126,135,143,153	11

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	b	1	14/15	-	-	153,171,178,184	0
5	NAG	b	2	14/15	-	-	155,168,187,195	0
5	BMA	b	3	11/12	-	-	152,158,170,178	0
5	MAN	b	4	11/12	-	-	170,174,183,187	0
5	MAN	b	5	11/12	-	-	168,173,179,183	0
6	NAG	I	1	14/15	-	-	125,137,150,154	14
6	NAG	I	2	14/15	-	-	157,164,178,178	14
7	NAG	S	2	14/15	0.14	0.13	58,63,79,90	14
9	BMA	W	3	11/12	0.44	0.19	147,166,181,198	0
9	MAN	W	4	11/12	0.47	0.11	205,211,221,222	0
7	MAN	Z	10	11/12	0.51	0.20	121,125,133,138	11
7	NAG	N	2	14/15	0.59	0.08	82,95,126,133	0
10	BMA	T	3	11/12	0.59	0.11	147,151,156,158	11
9	NAG	R	2	14/15	0.63	0.17	93,103,111,125	14
7	NAG	S	1	14/15	0.66	0.10	46,63,97,107	14
9	NAG	R	1	14/15	0.71	0.16	56,77,98,109	14
8	NAG	O	2	14/15	0.73	0.09	118,136,148,155	0
7	BMA	N	3	11/12	-	-	103,116,128,142	0
7	MAN	N	4	11/12	-	-	124,128,138,144	0
7	MAN	N	5	11/12	-	-	116,122,139,143	0
7	MAN	N	6	11/12	-	-	126,127,144,162	0
7	MAN	N	7	11/12	-	-	115,129,156,157	0
7	MAN	N	8	11/12	-	-	114,133,145,153	0
7	MAN	N	9	11/12	-	-	104,106,118,121	11
7	MAN	N	10	11/12	-	-	157,162,178,181	0
7	MAN	Z	4	11/12	0.74	0.14	117,120,124,131	11
7	MAN	Z	8	11/12	0.76	0.14	110,123,137,141	11
7	BMA	S	3	11/12	-	-	93,110,127,142	0
7	MAN	S	4	11/12	-	-	101,105,114,129	11
7	MAN	S	5	11/12	-	-	75,91,95,104	11
7	MAN	S	6	11/12	-	-	80,85,98,107	11
7	MAN	S	7	11/12	-	-	117,130,153,157	0
7	MAN	S	8	11/12	-	-	123,128,153,162	0
7	MAN	S	9	11/12	-	-	103,124,132,145	0
7	MAN	S	10	11/12	-	-	148,156,162,166	0
7	MAN	Z	7	11/12	0.78	0.17	109,115,125,126	11
6	NAG	Q	2	14/15	0.79	0.09	133,170,179,180	14
10	NAG	T	2	14/15	0.84	0.13	114,128,145,149	14
9	NAG	W	2	14/15	0.84	0.12	80,98,116,132	14
9	NAG	W	1	14/15	0.85	0.09	70,86,110,126	14
6	NAG	Q	1	14/15	0.85	0.14	155,164,174,179	0
6	NAG	V	1	14/15	0.85	0.14	124,140,154,155	0

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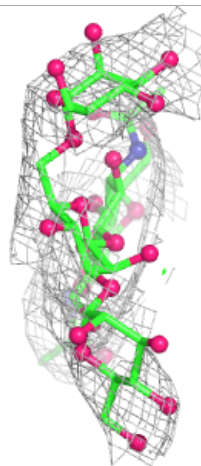
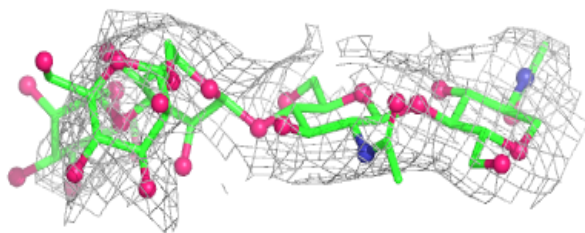
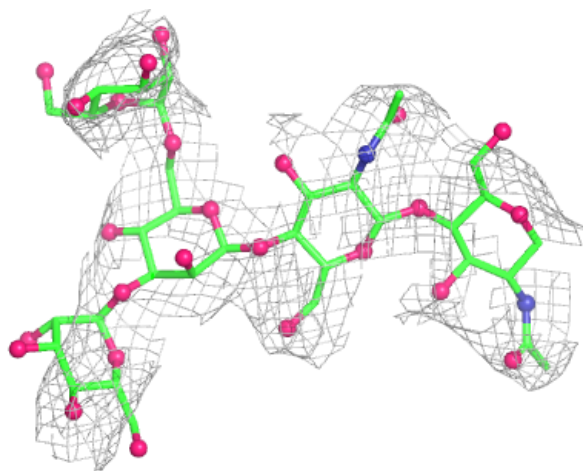
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	NAG	N	1	14/15	0.86	0.09	49,72,100,112	14
7	MAN	Z	9	11/12	0.86	0.09	122,126,136,150	0
6	NAG	M	2	14/15	0.88	0.09	155,163,180,181	0
7	BMA	Z	3	11/12	0.88	0.13	95,106,115,125	11
7	MAN	Z	6	11/12	0.89	0.15	115,121,131,133	11
8	BMA	O	3	11/12	-	-	132,136,155,163	0
8	MAN	O	4	11/12	-	-	171,189,201,208	0
8	MAN	O	5	11/12	-	-	182,193,197,198	0
8	MAN	O	6	11/12	-	-	164,169,181,185	0
7	MAN	Z	5	11/12	0.89	0.11	117,125,141,149	11
10	NAG	T	1	14/15	0.89	0.18	113,135,148,161	0
8	MAN	U	4	11/12	0.89	0.08	175,177,194,206	0
8	MAN	U	6	11/12	0.89	0.13	168,174,187,187	0
8	MAN	U	5	11/12	0.90	0.13	172,209,213,214	0
7	NAG	Z	2	14/15	0.92	0.16	73,88,99,102	14
8	NAG	O	1	14/15	0.92	0.10	117,135,150,161	14
6	NAG	V	2	14/15	0.92	0.11	158,170,180,183	0
9	BMA	R	3	11/12	-	-	139,150,171,186	11
9	MAN	R	4	11/12	-	-	184,196,201,202	0
8	NAG	U	2	14/15	0.92	0.11	134,146,162,176	14
6	NAG	P	2	14/15	0.93	0.08	165,177,184,185	14
7	NAG	Z	1	14/15	0.93	0.14	94,113,131,135	14
8	BMA	U	3	11/12	0.93	0.09	152,163,172,186	0
6	NAG	P	1	14/15	0.95	0.10	96,125,147,148	14
6	NAG	M	1	14/15	0.95	0.07	132,145,171,173	0
8	NAG	U	1	14/15	0.96	0.12	115,132,139,142	14
10	NAG	a	1	14/15	-	-	104,118,150,154	14
10	NAG	a	2	14/15	-	-	145,169,180,182	14
10	BMA	a	3	11/12	-	-	184,194,202,210	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.

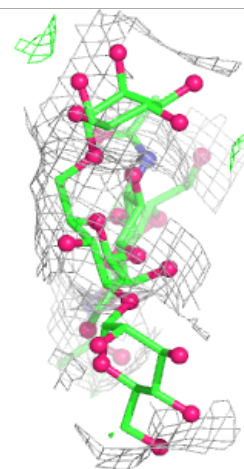
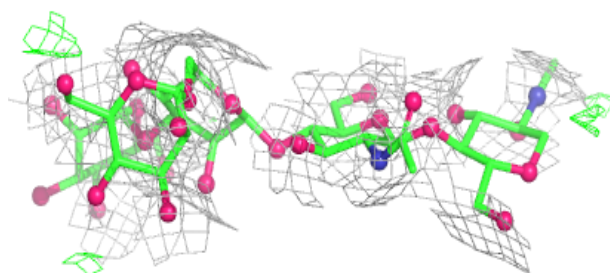
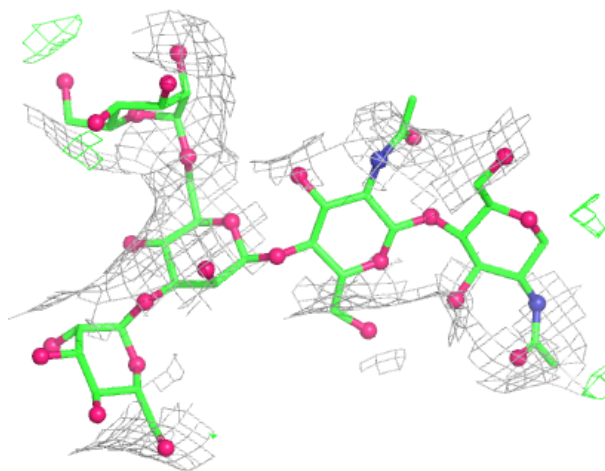
Electron density around Chain F:

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



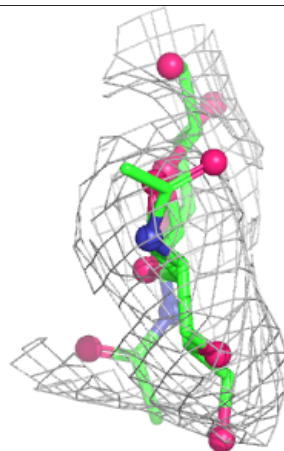
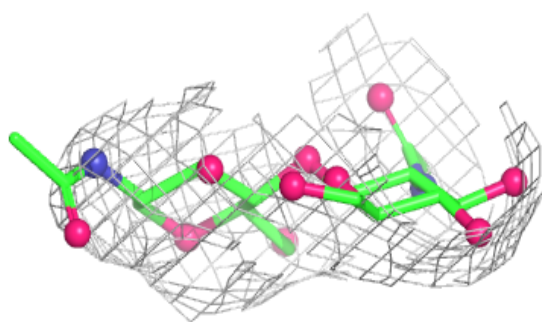
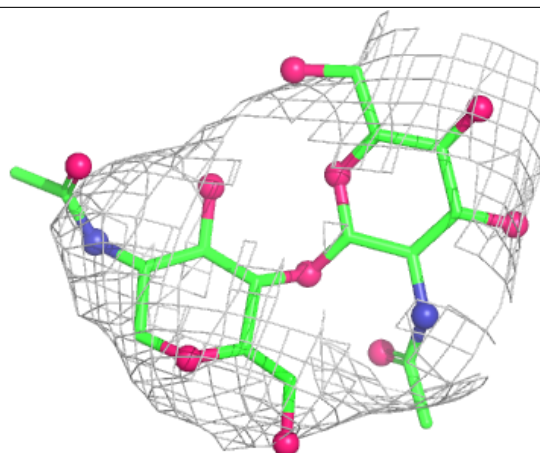
Electron density around Chain b:

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



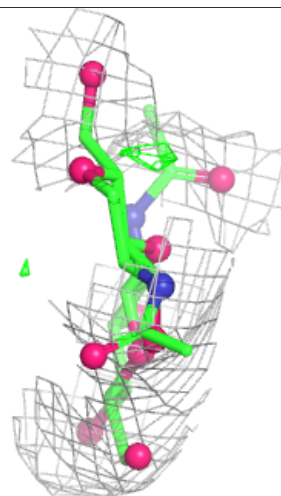
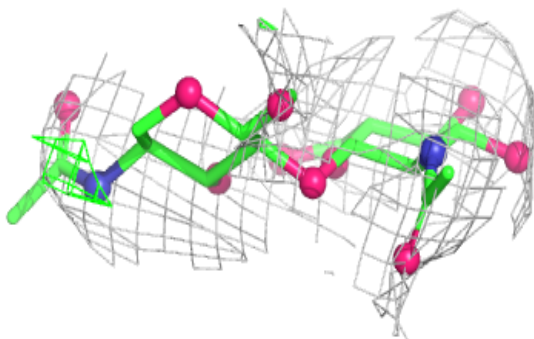
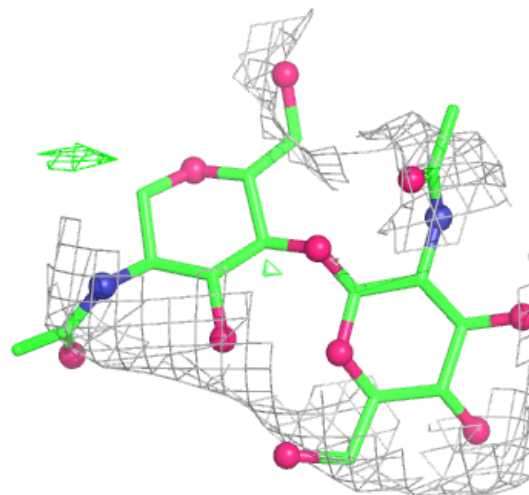
Electron density around Chain I:

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



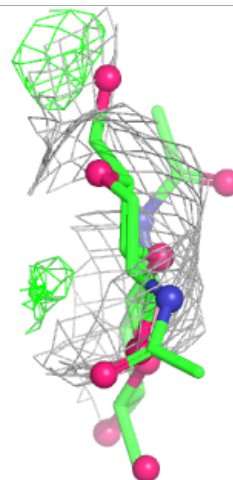
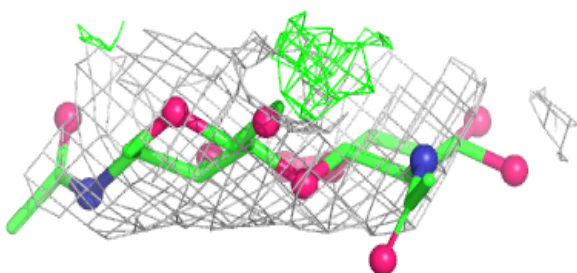
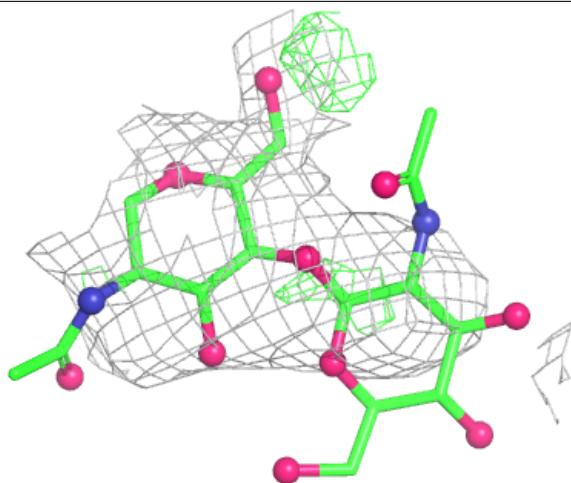
Electron density around Chain M:

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



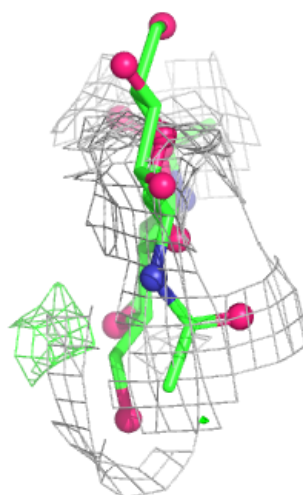
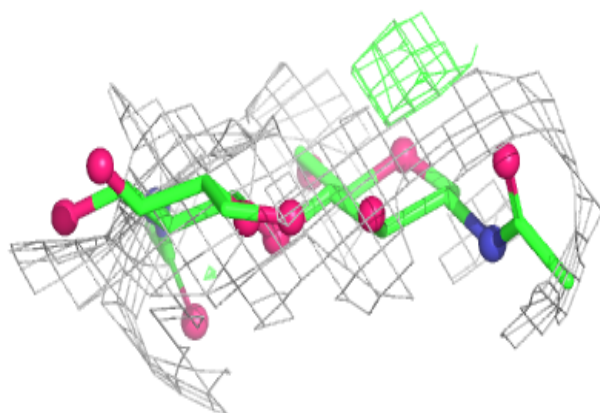
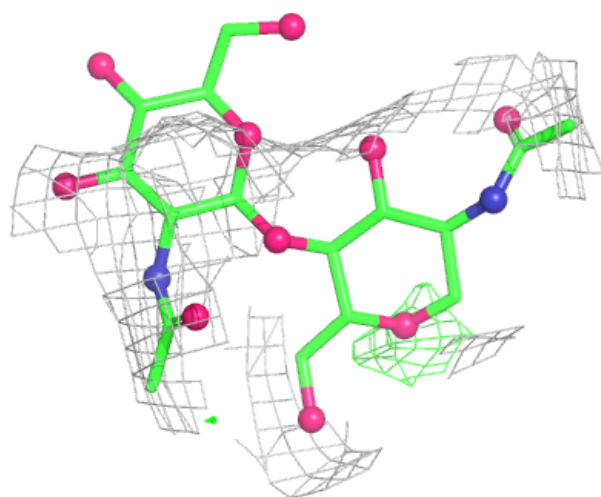
Electron density around Chain P:

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



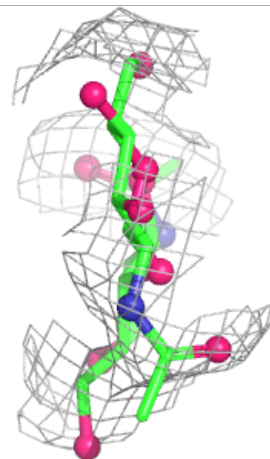
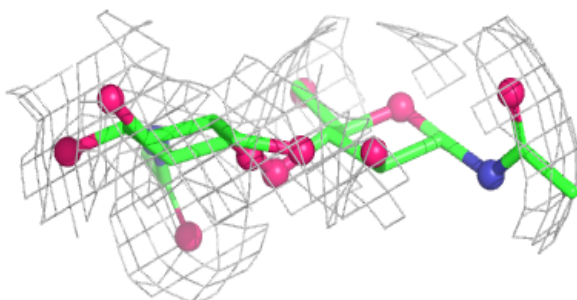
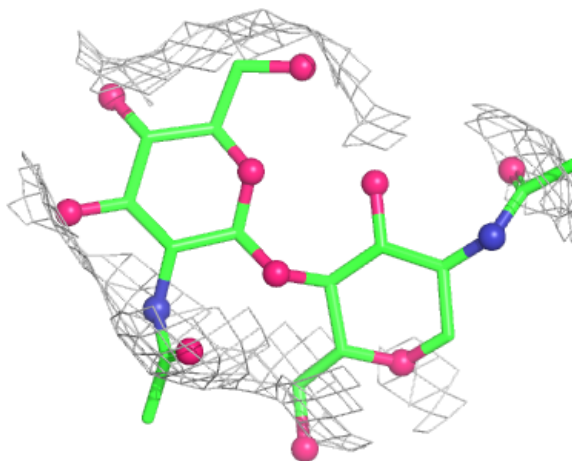
Electron density around Chain Q:

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



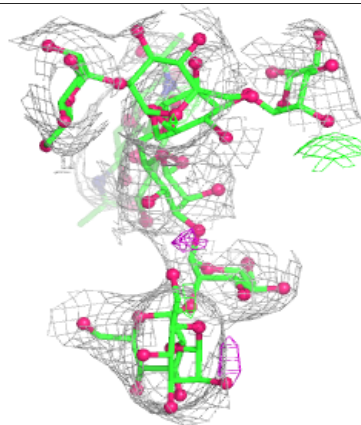
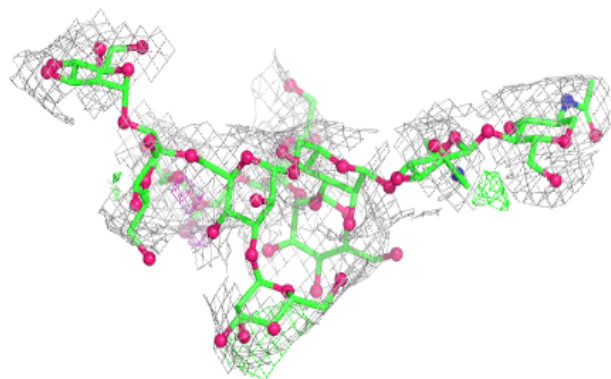
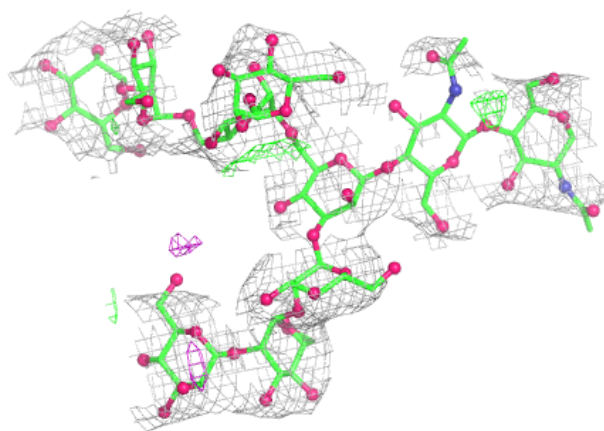
Electron density around Chain V:

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



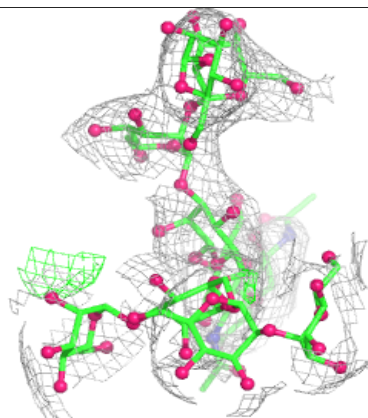
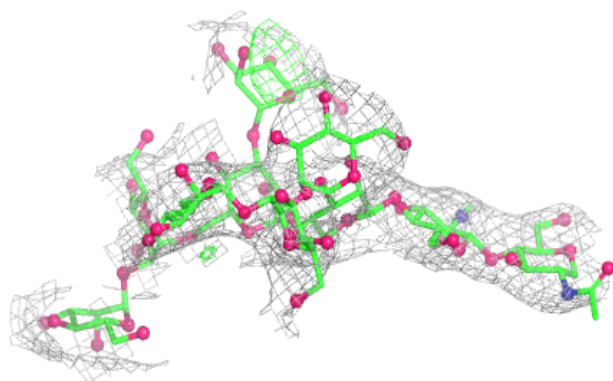
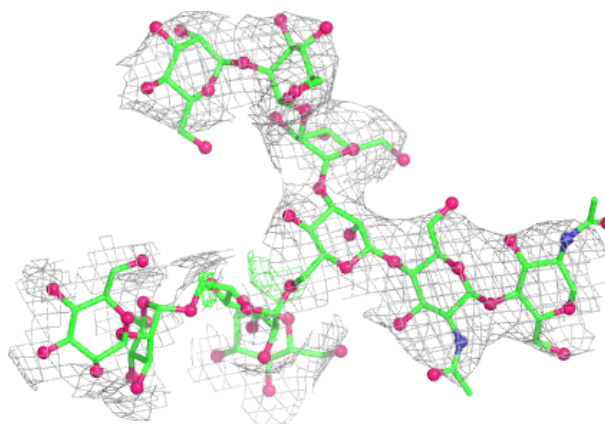
Electron density around Chain N:

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



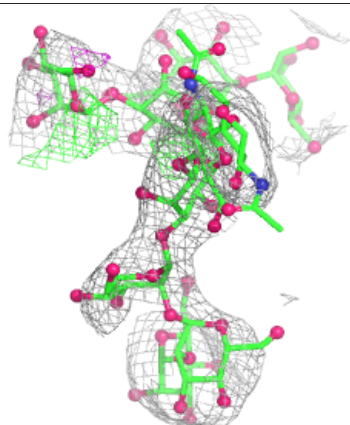
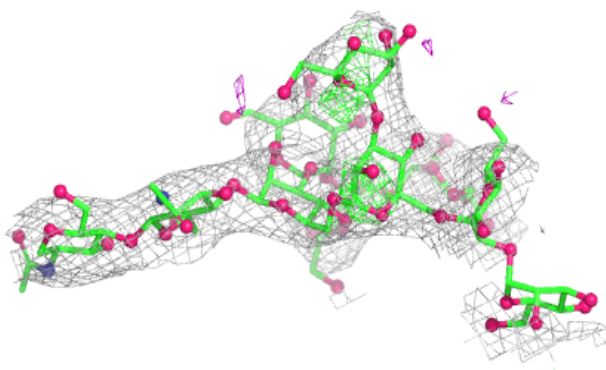
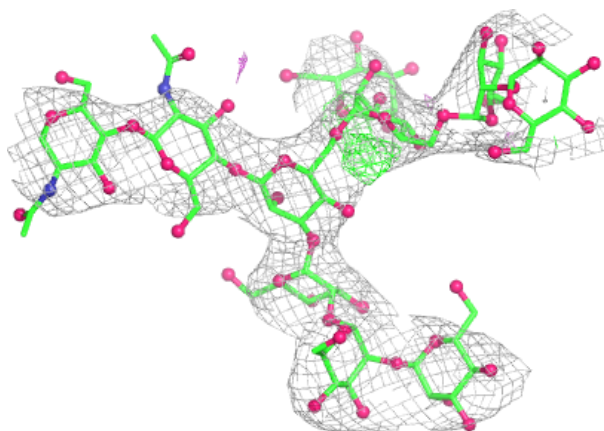
Electron density around Chain S:

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

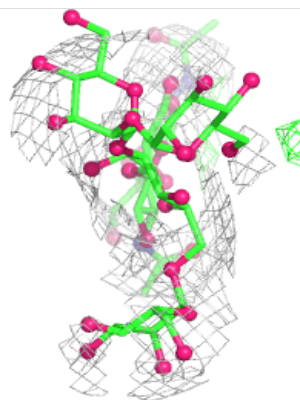
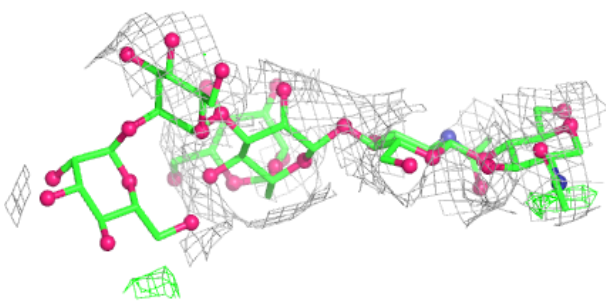
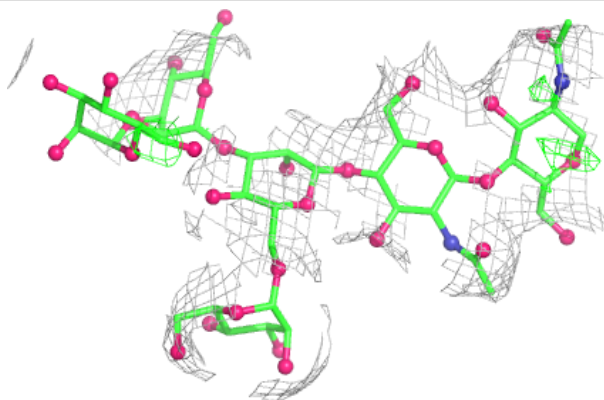


Electron density around Chain Z:

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

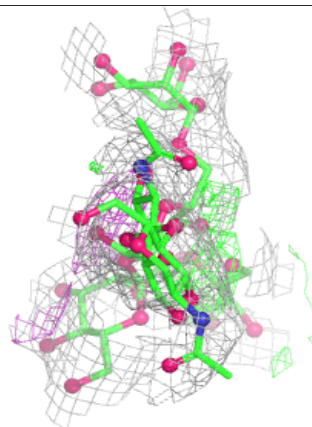
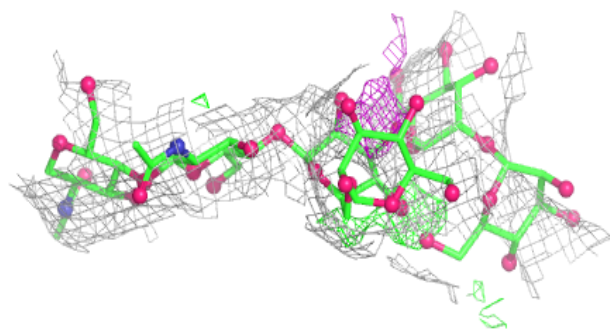
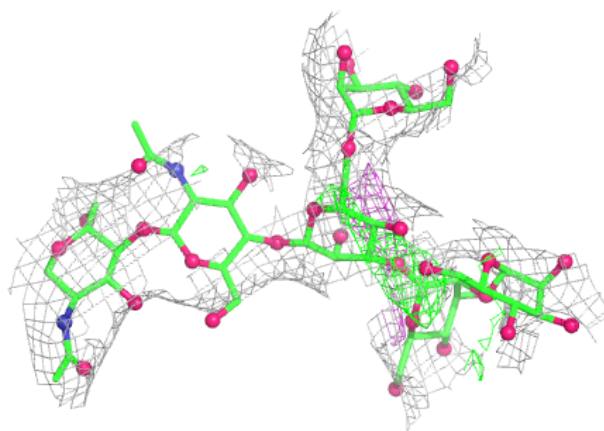
**Electron density around Chain O:**

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

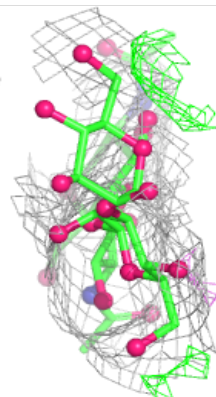
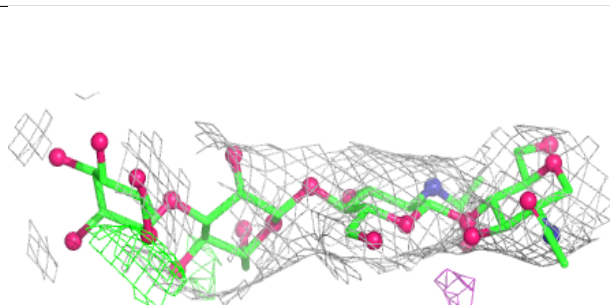
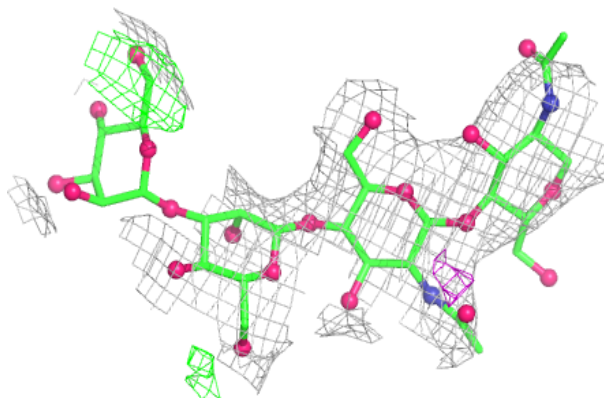


Electron density around Chain U:

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

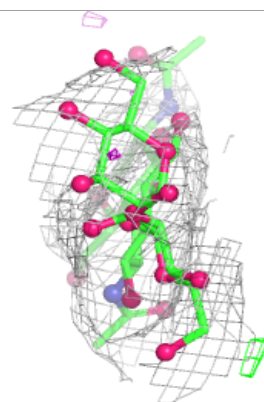
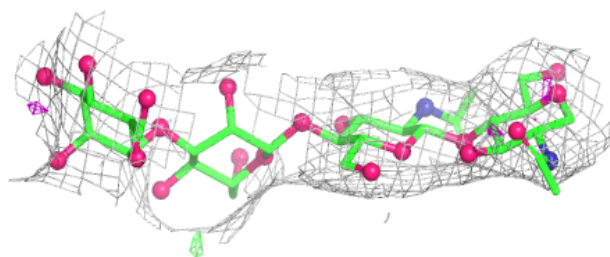
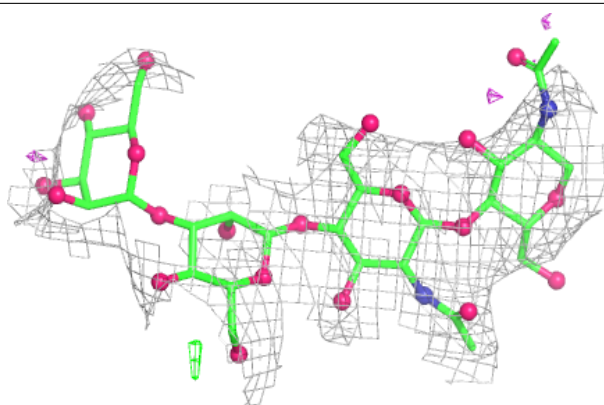
**Electron density around Chain R:**

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

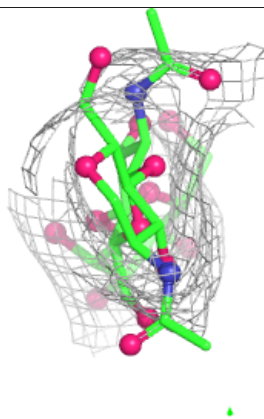
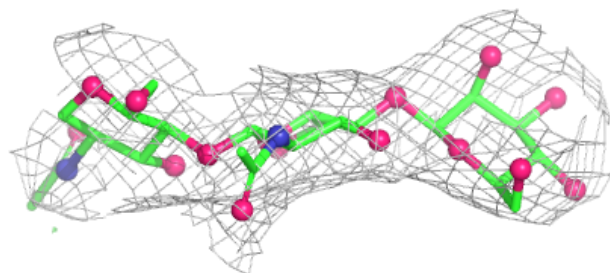
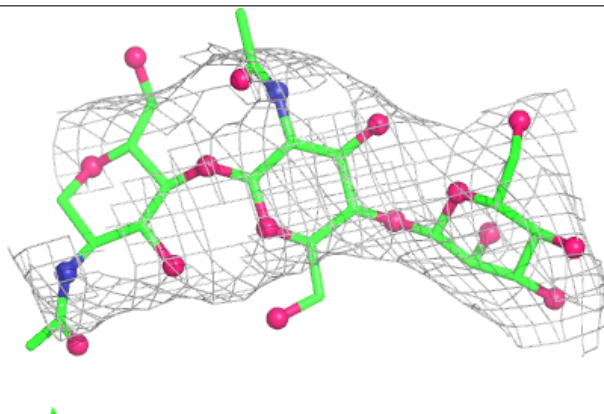


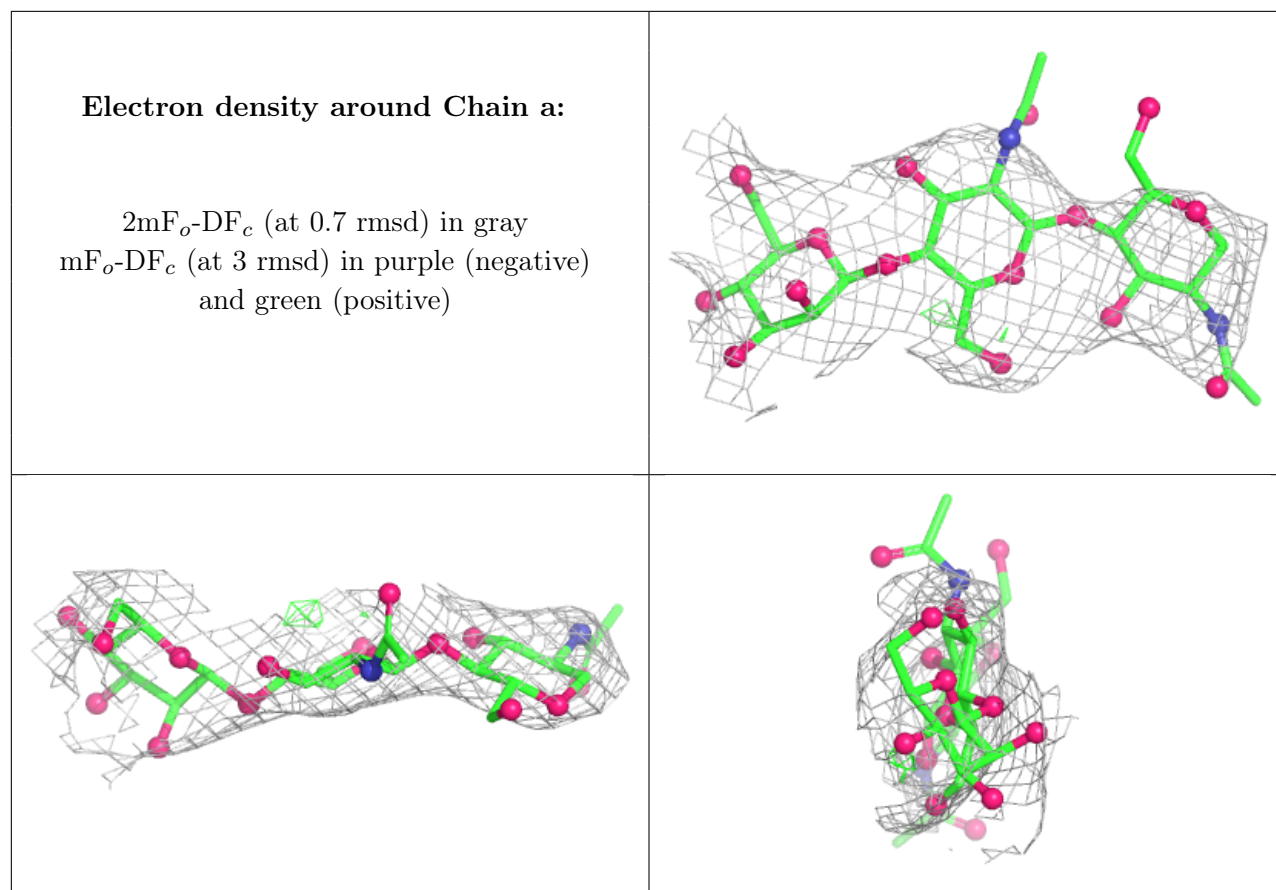
Electron density around Chain W:

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

**Electron density around Chain T:**

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





6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
11	NAG	A	1000	14/15	0.36	0.18	157,166,174,175	14
11	NAG	Y	1040	14/15	0.59	0.12	167,180,199,208	0
11	NAG	J	1042	14/15	0.70	0.15	115,125,139,145	14
11	NAG	K	1040	14/15	0.77	0.14	135,156,179,185	14
11	NAG	Y	1041	14/15	0.81	0.15	106,127,145,151	14
11	NAG	B	1040	14/15	0.85	0.10	112,132,141,143	0

6.5 Other polymers [i](#)

There are no such residues in this entry.