



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:33 PM UTC

PDB ID : 4PIR / pdb_00004pir
Title : X-ray structure of the mouse serotonin 5-HT₃ receptor
Authors : Hassaine, G.; Deluz, C.; Grasso, L.; Wyss, R.; Tol, M.B.; Hovius, R.; Graff, A.; Stahlberg, H.; Tomizaki, T.; Desmyter, A.; Moreau, C.; Li, X.-D.; Poitevin, F.; Vogel, H.; Nury, H.
Deposited on : 2014-05-09
Resolution : 3.50 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

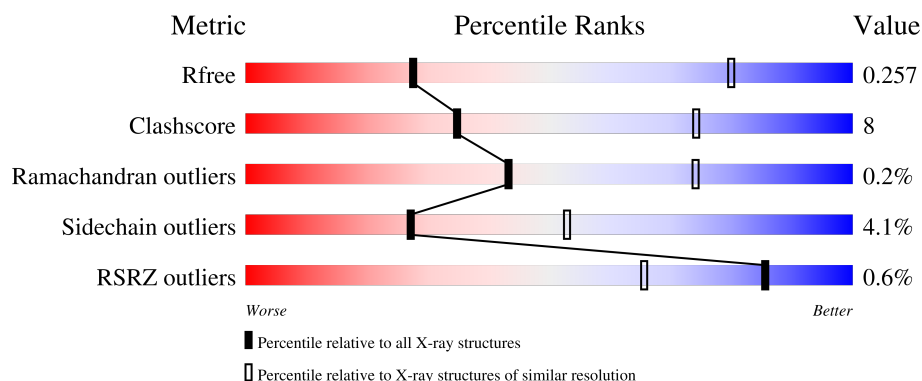
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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.



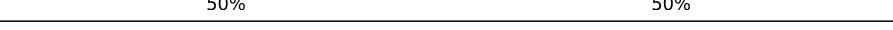
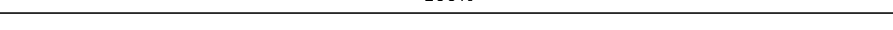
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	456	% 62% 20% • 16%
1	B	456	63% 20% • 16%
1	C	456	64% 20% • 16%
1	D	456	% 63% 20% • 16%
1	E	456	% 62% 21% • 16%

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Mol	Chain	Length	Quality of chain
2	F	124	 85% 12% ..
2	G	124	 88% 11% .
2	H	124	 81% 18% .
2	I	124	 88% 11% .
2	J	124	 84% 15% .
3	K	2	 50% 50%
3	M	2	 50% 50%
3	O	2	 50% 50%
3	P	2	 50% 50%
3	R	2	 50% 50%
3	S	2	 50% 50%
3	U	2	 50% 50%
3	V	2	 50% 50%
3	X	2	 50% 50%
4	L	3	 67% 33%
4	N	3	 67% 33%
4	Q	3	 67% 33%
4	T	3	 33% 67%
4	W	3	 100%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 20739 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 5-hydroxytryptamine receptor 3A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	383	Total	C	N	O	S	0	0	0
			3108	2045	507	547	9			
1	B	383	Total	C	N	O	S	0	0	0
			3111	2046	507	549	9			
1	C	384	Total	C	N	O	S	0	0	0
			3125	2057	509	550	9			
1	D	384	Total	C	N	O	S	0	0	0
			3125	2057	509	550	9			
1	E	383	Total	C	N	O	S	0	0	0
			3111	2046	507	549	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	ALA	GLU	conflict	UNP P23979
A	5	ARG	-	insertion	UNP P23979
A	277	ALA	-	insertion	UNP P23979
A	461	SER	TYR	conflict	UNP P23979
B	4	ALA	GLU	conflict	UNP P23979
B	5	ARG	-	insertion	UNP P23979
B	277	ALA	-	insertion	UNP P23979
B	461	SER	TYR	conflict	UNP P23979
C	4	ALA	GLU	conflict	UNP P23979
C	5	ARG	-	insertion	UNP P23979
C	277	ALA	-	insertion	UNP P23979
C	461	SER	TYR	conflict	UNP P23979
D	4	ALA	GLU	conflict	UNP P23979
D	5	ARG	-	insertion	UNP P23979
D	277	ALA	-	insertion	UNP P23979
D	461	SER	TYR	conflict	UNP P23979
E	4	ALA	GLU	conflict	UNP P23979
E	5	ARG	-	insertion	UNP P23979
E	277	ALA	-	insertion	UNP P23979

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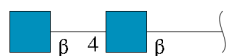
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Chain	Residue	Modelled	Actual	Comment	Reference
E	461	SER	TYR	conflict	UNP P23979

- Molecule 2 is a protein called VHH15.

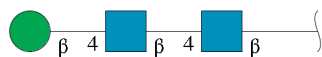
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	122	Total	C	N	O	S	0	0	0
			920	572	162	182	4			
2	G	124	Total	C	N	O	S	0	0	0
			936	579	164	189	4			
2	H	124	Total	C	N	O	S	0	0	0
			936	579	164	189	4			
2	I	124	Total	C	N	O	S	0	0	0
			936	579	164	189	4			
2	J	124	Total	C	N	O	S	0	0	0
			936	579	164	189	4			

- Molecule 3 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
3	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	M	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	O	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	P	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	R	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	S	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	U	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	V	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	X	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 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
4	L	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	N	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	Q	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	T	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	W	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

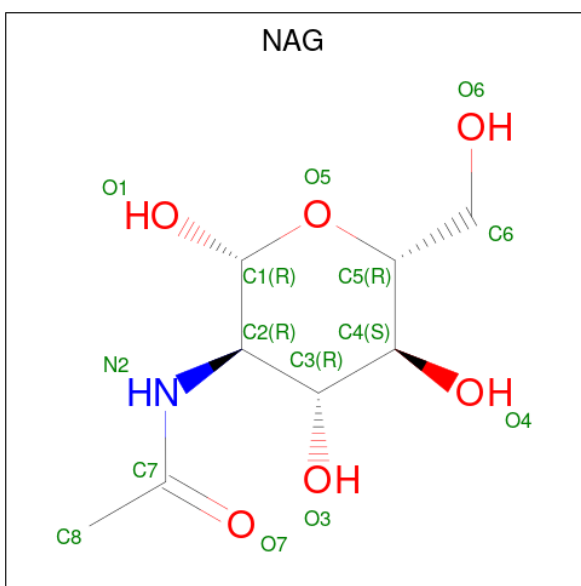
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		
5	B	1	Total	Cl	0	0
			1	1		
5	C	1	Total	Cl	0	0
			1	1		
5	D	1	Total	Cl	0	0
			1	1		
5	E	1	Total	Cl	0	0
			1	1		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	O	S	0	0
			5	4	1		
6	H	1	Total	O	S	0	0
			5	4	1		
6	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

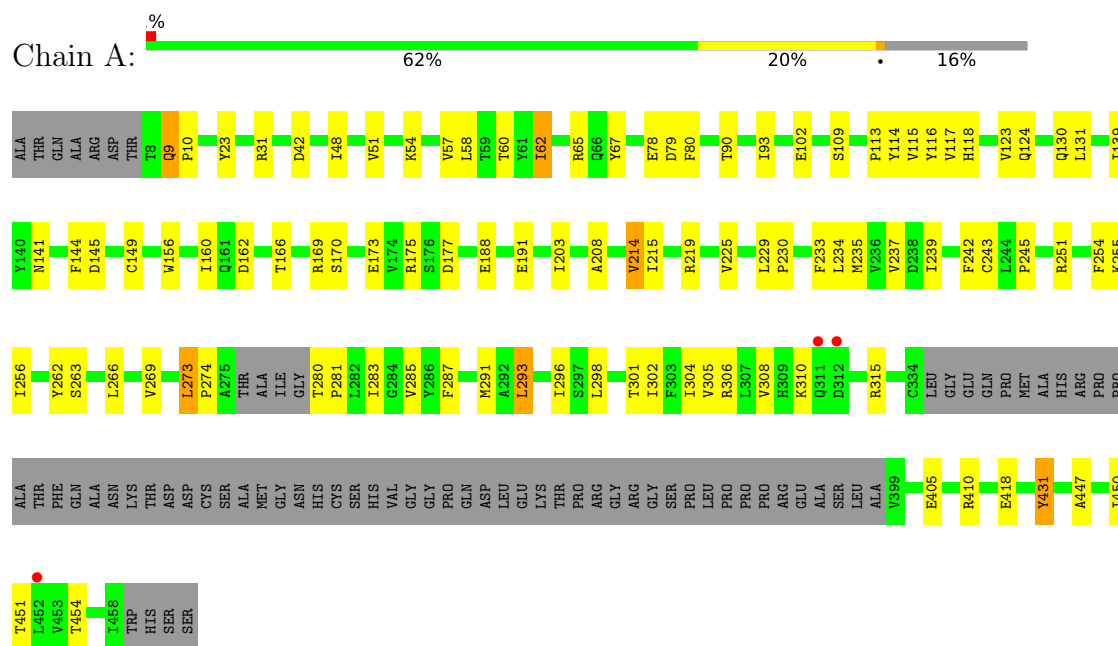


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	E	1	Total	C	N	O	0	0
			14	8	1	5		
7	E	1	Total	C	N	O	0	0
			14	8	1	5		

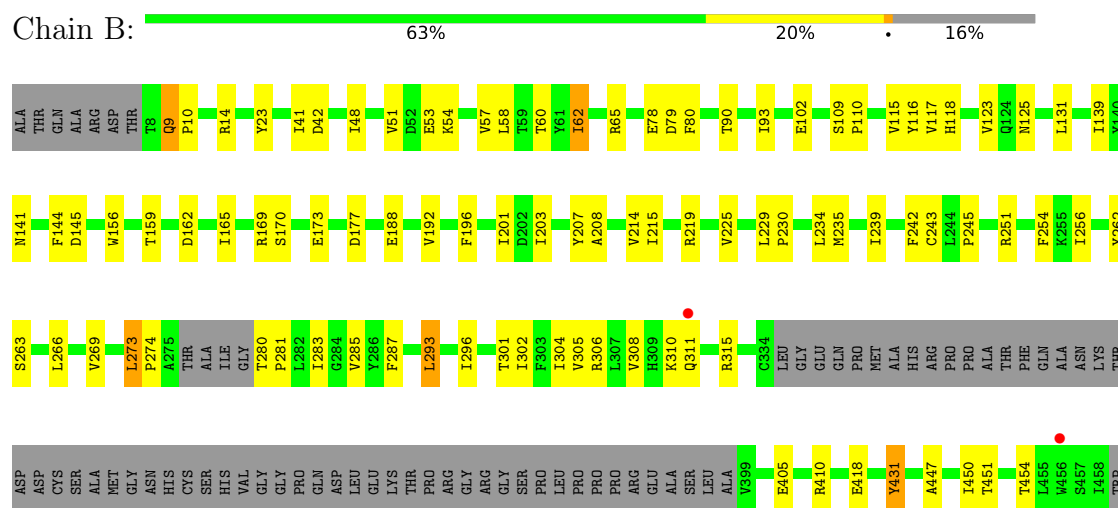
3 Residue-property plots

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: 5-hydroxytryptamine receptor 3A



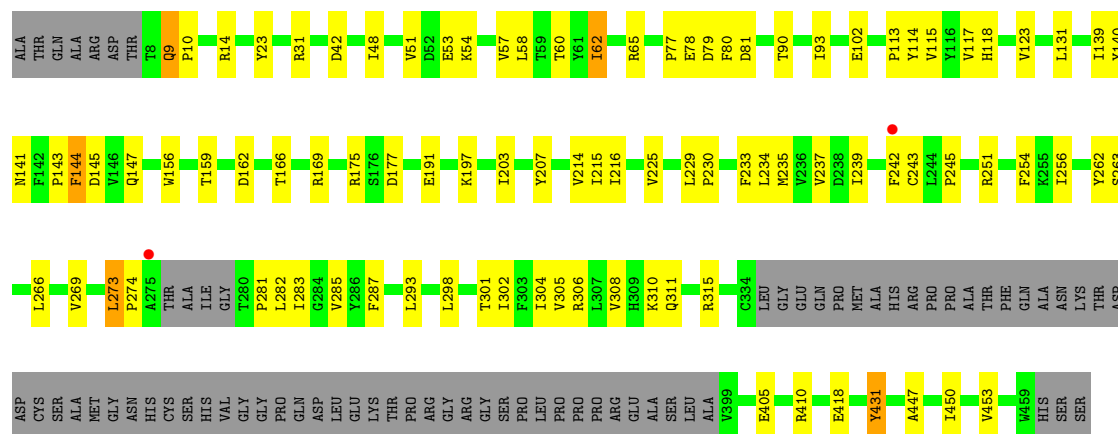
• Molecule 1: 5-hydroxytryptamine receptor 3A



HIS
SER
SER

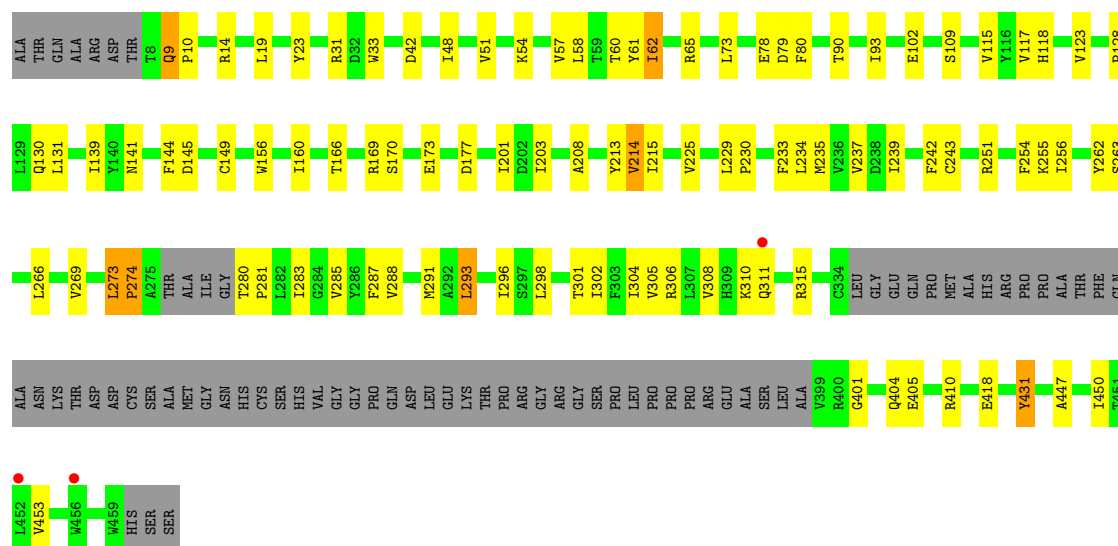
• Molecule 1: 5-hydroxytryptamine receptor 3A

Chain C: 



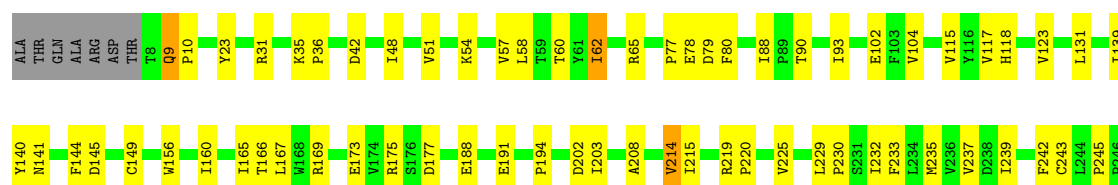
• Molecule 1: 5-hydroxytryptamine receptor 3A

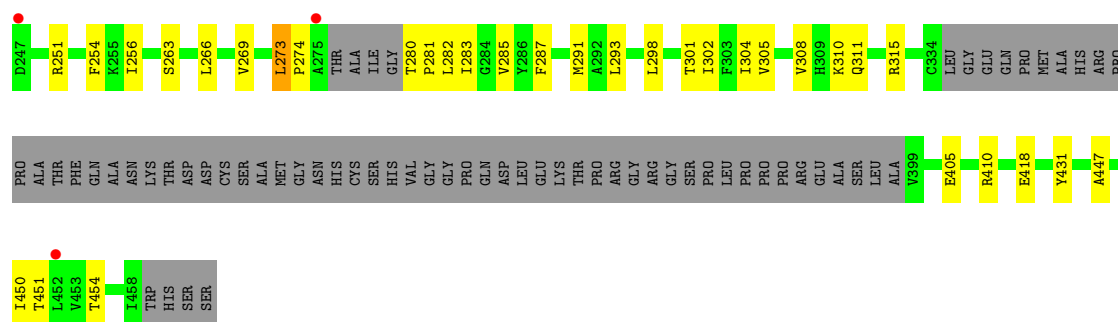
Chain D: 



• Molecule 1: 5-hydroxytryptamine receptor 3A

Chain E: 





● Molecule 2: VHH15

Chain F: 85% 12% ..



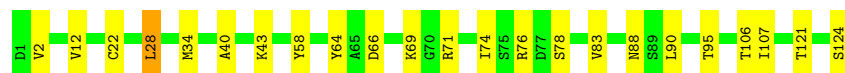
● Molecule 2: VHH15

Chain G: 88% 11% .



● Molecule 2: VHH15

Chain H: 81% 18% .



● Molecule 2: VHH15

Chain I: 88% 11% .



● Molecule 2: VHH15

Chain J: 84% 15% .



● Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K: 50% 50%



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

Chain M:  50%  50%

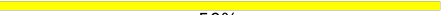


- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  50%  50%



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

Chain P:  50%  50%



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

Chain R:  50%  50%



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

Chain S:  50%  50%



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

Chain U:  50%  50%



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

Chain V:  50% 50%



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

Chain X:  50% 50%



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

Chain L:  67% 33%



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

Chain N:  67% 33%

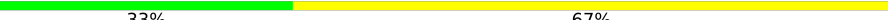


- Molecule 4: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  67% 33%

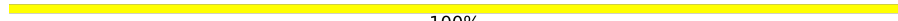


- Molecule 4: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:  33% 67%



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

Chain W:  100%

MAG1
MAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	121.54Å 187.09Å 240.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.50 20.00 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-3.50) 99.2 (20.00-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 3.52Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1685)	Depositor
R, R_{free}	0.217 , 0.257 0.219 , 0.257	Depositor DCC
R_{free} test set	3508 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	97.1	Xtriage
Anisotropy	0.534	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 84.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20739	wwPDB-VP
Average B, all atoms (Å ²)	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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: BMA, SO4, NAG, CL

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	0.28	0/3190	0.72	0/4358
1	B	0.28	0/3193	0.73	0/4362
1	C	0.28	0/3209	0.73	0/4385
1	D	0.28	0/3209	0.73	0/4385
1	E	0.28	0/3193	0.72	0/4362
2	F	0.25	0/936	0.65	0/1269
2	G	0.25	0/952	0.65	0/1289
2	H	0.25	0/952	0.65	0/1289
2	I	0.26	0/952	0.65	0/1289
2	J	0.25	0/952	0.66	0/1289
All	All	0.27	0/20738	0.71	0/28277

There are no bond length outliers.

There are no bond angle outliers.

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	3108	0	3087	58	0
1	B	3111	0	3089	57	0
1	C	3125	0	3099	59	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3125	0	3099	63	0
1	E	3111	0	3089	59	0
2	F	920	0	886	11	0
2	G	936	0	898	10	0
2	H	936	0	898	15	0
2	I	936	0	898	9	0
2	J	936	0	900	12	0
3	K	28	0	25	0	0
3	M	28	0	25	1	0
3	O	28	0	25	0	0
3	P	28	0	25	1	0
3	R	28	0	25	0	0
3	S	28	0	25	1	0
3	U	28	0	25	0	0
3	V	28	0	25	2	0
3	X	28	0	25	1	0
4	L	39	0	34	1	0
4	N	39	0	34	1	0
4	Q	39	0	34	1	0
4	T	39	0	34	1	0
4	W	39	0	34	2	0
5	A	1	0	0	1	0
5	B	1	0	0	0	0
5	C	1	0	0	1	0
5	D	1	0	0	1	0
5	E	1	0	0	1	0
6	D	5	0	0	0	0
6	H	10	0	0	1	0
7	E	28	0	26	0	0
All	All	20739	0	20364	310	0

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

All (310) 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:E:31:ARG:NH1	5:E:508:CL:CL	2.35	0.95
1:C:31:ARG:NH1	5:C:508:CL:CL	2.50	0.81
1:B:90:THR:HG22	1:B:115:VAL:HG13	1.64	0.80
1:D:90:THR:HG22	1:D:115:VAL:HG13	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:90:THR:HG22	1:E:115:VAL:HG13	1.69	0.73
1:C:215:ILE:HG12	3:S:1:NAG:H82	1.72	0.72
1:E:215:ILE:HG12	3:X:1:NAG:H82	1.72	0.72
1:A:215:ILE:HG12	3:M:1:NAG:H82	1.73	0.71
1:C:90:THR:HG22	1:C:115:VAL:HG13	1.72	0.70
1:D:215:ILE:HG12	3:V:1:NAG:H82	1.73	0.69
1:A:90:THR:HG22	1:A:115:VAL:HG13	1.73	0.69
1:E:117:VAL:HG22	1:E:123:VAL:HG22	1.76	0.68
1:B:215:ILE:HG12	3:P:1:NAG:H82	1.77	0.67
1:D:42:ASP:OD1	1:D:169:ARG:NH1	2.29	0.66
1:B:42:ASP:OD1	1:B:169:ARG:NH1	2.30	0.64
1:A:31:ARG:NH1	5:A:508:CL:CL	2.62	0.64
2:H:95:THR:HG23	2:H:121:THR:HA	1.80	0.64
1:C:42:ASP:OD1	1:C:169:ARG:NH1	2.31	0.63
1:D:117:VAL:HG22	1:D:123:VAL:HG22	1.81	0.62
2:J:34:MET:HG2	2:J:83:VAL:HG11	1.81	0.62
1:A:203:ILE:HG23	2:F:76:ARG:NH1	2.13	0.62
2:H:34:MET:HG2	2:H:83:VAL:HG11	1.82	0.61
2:I:34:MET:HG2	2:I:83:VAL:HG11	1.82	0.61
1:C:117:VAL:HG22	1:C:123:VAL:HG22	1.83	0.61
1:A:67:TYR:HE1	1:A:124:GLN:HG3	1.65	0.61
2:I:95:THR:HG23	2:I:121:THR:HA	1.82	0.60
1:A:42:ASP:OD1	1:A:169:ARG:NH1	2.34	0.60
1:C:42:ASP:HB3	1:C:65:ARG:HB2	1.82	0.60
2:J:95:THR:HG23	2:J:121:THR:HA	1.84	0.60
1:E:42:ASP:OD1	1:E:169:ARG:NH1	2.35	0.59
1:A:117:VAL:HG22	1:A:123:VAL:HG22	1.83	0.59
1:A:306:ARG:HG3	1:E:245:PRO:HD3	1.84	0.59
1:B:117:VAL:HG22	1:B:123:VAL:HG22	1.84	0.58
1:C:410:ARG:NH1	1:D:405:GLU:OE1	2.37	0.58
1:A:405:GLU:OE1	1:E:410:ARG:NH1	2.36	0.58
1:D:42:ASP:HB3	1:D:65:ARG:HB2	1.85	0.58
2:G:95:THR:HG23	2:G:121:THR:HA	1.85	0.58
1:C:203:ILE:HG23	2:H:76:ARG:NH1	2.18	0.58
1:D:203:ILE:HG23	2:I:76:ARG:NH1	2.19	0.57
1:E:42:ASP:HB3	1:E:65:ARG:HB2	1.84	0.57
1:C:48:ILE:HA	1:C:60:THR:HG22	1.87	0.57
1:D:42:ASP:CG	1:D:169:ARG:HH11	2.13	0.57
2:F:76:ARG:HH12	2:F:78:SER:HB3	1.69	0.57
2:F:95:THR:HG23	2:F:121:THR:HA	1.86	0.57
1:C:251:ARG:NH1	1:C:304:ILE:HD11	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:48:ILE:HA	1:D:60:THR:HG22	1.86	0.56
1:E:315:ARG:NH2	1:E:418:GLU:OE2	2.36	0.56
1:A:23:TYR:HE1	1:A:93:ILE:HA	1.71	0.56
1:D:255:LYS:NZ	1:D:301:THR:OG1	2.35	0.56
1:A:315:ARG:NH2	1:A:418:GLU:OE2	2.35	0.56
1:B:243:CYS:HB2	1:C:302:ILE:HD13	1.88	0.55
1:B:410:ARG:NH1	1:C:405:GLU:OE1	2.38	0.55
1:D:315:ARG:NH2	1:D:418:GLU:OE2	2.37	0.55
1:A:255:LYS:NZ	1:A:301:THR:OG1	2.35	0.55
1:B:245:PRO:HD3	1:C:306:ARG:HG3	1.89	0.55
1:C:169:ARG:NH2	1:C:177:ASP:OD2	2.40	0.55
1:A:175:ARG:NE	1:A:191:GLU:OE2	2.39	0.55
1:D:51:VAL:HG12	1:D:58:LEU:HD13	1.89	0.55
1:D:234:LEU:HD13	1:D:262:TYR:HD1	1.72	0.55
1:C:245:PRO:HD3	1:D:306:ARG:HG3	1.88	0.55
1:E:203:ILE:HG23	2:J:76:ARG:HD2	1.89	0.55
2:I:22:CYS:HB3	2:I:83:VAL:HG13	1.89	0.54
1:A:302:ILE:HD13	1:E:243:CYS:HB2	1.88	0.54
1:B:42:ASP:HB3	1:B:65:ARG:HB2	1.88	0.54
2:F:34:MET:HG2	2:F:83:VAL:HG11	1.89	0.54
2:H:22:CYS:HB3	2:H:83:VAL:HG13	1.88	0.54
1:D:31:ARG:NH1	5:D:509:CL:CL	2.75	0.54
1:C:42:ASP:CG	1:C:169:ARG:HH11	2.16	0.54
1:B:254:PHE:CE2	1:C:256:ILE:HD11	2.42	0.54
1:B:234:LEU:HD13	1:B:262:TYR:HD1	1.73	0.53
2:J:22:CYS:HB3	2:J:83:VAL:HG13	1.90	0.53
1:B:42:ASP:CG	1:B:169:ARG:HH11	2.16	0.53
1:E:188:GLU:HB2	1:E:219:ARG:HD3	1.91	0.53
1:A:256:ILE:HD11	1:E:254:PHE:CE2	2.43	0.53
1:C:23:TYR:HE1	1:C:93:ILE:HA	1.75	0.52
1:B:203:ILE:HG23	2:G:76:ARG:HD2	1.91	0.52
1:D:243:CYS:HB2	1:E:302:ILE:HD13	1.92	0.52
1:D:251:ARG:HH11	1:D:301:THR:HG23	1.75	0.51
1:A:234:LEU:HD13	1:A:262:TYR:HD1	1.75	0.51
1:B:48:ILE:HA	1:B:60:THR:HG22	1.91	0.51
1:B:266:LEU:HA	1:B:269:VAL:HG12	1.93	0.51
1:C:243:CYS:HB2	1:D:302:ILE:HD13	1.91	0.51
1:C:254:PHE:CE2	1:D:256:ILE:HD11	2.45	0.51
1:E:251:ARG:NH1	1:E:304:ILE:HD11	2.25	0.51
1:B:23:TYR:HE2	1:B:93:ILE:HA	1.76	0.51
1:B:192:VAL:HG22	1:B:214:VAL:HG12	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:254:PHE:CE2	1:E:256:ILE:HD11	2.45	0.51
1:A:251:ARG:HH11	1:A:301:THR:HG23	1.76	0.51
1:D:23:TYR:HE2	1:D:93:ILE:HA	1.75	0.50
1:D:273:LEU:HD13	1:D:274:PRO:HD2	1.92	0.50
1:A:243:CYS:HB2	1:B:302:ILE:HD13	1.93	0.50
1:B:10:PRO:HD2	1:B:14:ARG:HD3	1.94	0.50
1:A:410:ARG:NH1	1:B:405:GLU:OE1	2.44	0.50
1:A:431:TYR:OH	1:B:311:GLN:OE1	2.27	0.50
1:B:51:VAL:HG12	1:B:58:LEU:HD13	1.92	0.50
2:F:7:SER:OG	2:F:21:SER:OG	2.27	0.50
1:A:254:PHE:CE2	1:B:256:ILE:HD11	2.46	0.50
1:B:41:ILE:HD13	1:B:165:ILE:HD11	1.92	0.50
1:B:169:ARG:NH2	1:B:177:ASP:OD2	2.45	0.50
1:E:202:ASP:OD2	2:J:30:SER:OG	2.29	0.50
1:E:42:ASP:CG	1:E:169:ARG:HH11	2.20	0.49
1:E:239:ILE:HA	1:E:242:PHE:HD2	1.77	0.49
1:C:175:ARG:NE	1:C:191:GLU:OE1	2.45	0.49
1:A:166:THR:HG22	4:L:1:NAG:H62	1.94	0.49
1:C:266:LEU:HA	1:C:269:VAL:HG12	1.95	0.49
2:F:22:CYS:HB3	2:F:83:VAL:HG13	1.94	0.49
2:J:106:THR:OG1	2:J:107:ILE:N	2.46	0.49
1:E:23:TYR:HE2	1:E:93:ILE:HA	1.77	0.49
1:E:48:ILE:HA	1:E:60:THR:HG22	1.94	0.49
1:B:118:HIS:CE1	2:H:28:LEU:HD12	2.48	0.49
1:B:239:ILE:HA	1:B:242:PHE:HD2	1.78	0.49
1:A:48:ILE:HA	1:A:60:THR:HG22	1.94	0.48
1:C:51:VAL:HG12	1:C:58:LEU:HD13	1.95	0.48
1:E:149:CYS:HB2	1:E:214:VAL:HG13	1.95	0.48
1:A:42:ASP:CG	1:A:169:ARG:HH11	2.21	0.48
1:A:251:ARG:NH1	1:A:304:ILE:HD11	2.29	0.48
2:G:106:THR:OG1	2:G:107:ILE:N	2.46	0.48
1:E:169:ARG:NH2	1:E:177:ASP:OD2	2.47	0.48
1:E:118:HIS:CE1	2:F:28:LEU:HD12	2.49	0.48
2:F:106:THR:OG1	2:F:107:ILE:N	2.47	0.48
2:I:106:THR:OG1	2:I:107:ILE:N	2.45	0.47
1:B:141:ASN:O	1:B:145:ASP:HB3	2.13	0.47
1:A:141:ASN:O	1:A:145:ASP:HB3	2.13	0.47
1:A:239:ILE:HA	1:A:242:PHE:HD2	1.80	0.47
1:D:141:ASN:O	1:D:145:ASP:HB3	2.14	0.47
1:D:160:ILE:HG13	1:D:208:ALA:HB2	1.97	0.47
1:B:102:GLU:OE2	1:B:102:GLU:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:90:LEU:HB3	2:G:122:VAL:HG21	1.95	0.47
1:A:266:LEU:HA	1:A:269:VAL:HG12	1.97	0.47
1:D:10:PRO:HD2	1:D:14:ARG:HD3	1.96	0.47
2:H:106:THR:OG1	2:H:107:ILE:N	2.48	0.47
1:C:141:ASN:O	1:C:145:ASP:HB3	2.15	0.47
1:C:234:LEU:HD13	1:C:262:TYR:HD1	1.79	0.47
1:D:239:ILE:HA	1:D:242:PHE:HD2	1.78	0.47
2:H:76:ARG:HH12	2:H:78:SER:HB3	1.80	0.47
1:A:109:SER:HG	1:A:130:GLN:H	1.62	0.46
1:C:431:TYR:OH	1:D:311:GLN:OE1	2.27	0.46
1:D:266:LEU:HA	1:D:269:VAL:HG12	1.95	0.46
1:E:266:LEU:HA	1:E:269:VAL:HG12	1.97	0.46
1:D:251:ARG:NH1	1:D:304:ILE:HD11	2.31	0.46
2:H:66:ASP:HA	2:H:69:LYS:HD3	1.97	0.46
1:A:51:VAL:HG12	1:A:58:LEU:HD13	1.96	0.46
1:E:273:LEU:HD13	1:E:274:PRO:HD2	1.98	0.46
1:B:251:ARG:NH1	1:B:304:ILE:HD11	2.31	0.46
1:E:285:VAL:C	1:E:287:PHE:H	2.24	0.46
1:A:42:ASP:HB3	1:A:65:ARG:HB2	1.97	0.46
1:B:285:VAL:C	1:B:287:PHE:H	2.24	0.46
1:C:273:LEU:HD13	1:C:274:PRO:HD2	1.98	0.46
1:C:285:VAL:C	1:C:287:PHE:H	2.24	0.46
2:H:12:VAL:HG21	2:H:90:LEU:HD13	1.97	0.45
1:A:188:GLU:HB2	1:A:219:ARG:HD3	1.98	0.45
1:A:266:LEU:HD21	1:A:291:MET:HE3	1.98	0.45
1:A:273:LEU:HD13	1:A:274:PRO:HD2	1.98	0.45
1:E:141:ASN:O	1:E:145:ASP:HB3	2.16	0.45
1:E:280:THR:HA	1:E:281:PRO:HD3	1.84	0.45
2:G:20:LEU:HD13	2:G:87:MET:HE2	1.98	0.45
1:C:239:ILE:HA	1:C:242:PHE:HD2	1.81	0.45
1:D:266:LEU:HD21	1:D:291:MET:HE3	1.99	0.45
2:J:52:THR:O	2:J:61:SER:OG	2.27	0.45
1:C:78:GLU:C	1:C:80:PHE:H	2.24	0.45
1:C:118:HIS:CE1	2:I:28:LEU:HD12	2.50	0.45
1:C:102:GLU:N	1:C:102:GLU:OE2	2.49	0.45
1:D:169:ARG:NH2	1:D:177:ASP:OD2	2.49	0.45
1:A:160:ILE:HG13	1:A:208:ALA:HB2	1.99	0.45
1:A:280:THR:HA	1:A:281:PRO:HD3	1.83	0.45
1:A:285:VAL:C	1:A:287:PHE:H	2.23	0.45
1:D:61:TYR:HE1	1:D:128:PRO:HB2	1.82	0.45
1:D:62:ILE:HD13	1:D:131:LEU:HD12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:ILE:HG23	2:G:76:ARG:NH1	2.31	0.45
1:B:281:PRO:C	1:B:283:ILE:H	2.25	0.45
1:B:451:THR:O	1:B:454:THR:HG22	2.17	0.45
1:E:51:VAL:HG12	1:E:58:LEU:HD13	1.98	0.44
1:E:160:ILE:HG13	1:E:208:ALA:HB2	1.98	0.44
1:B:53:GLU:HG2	1:B:273:LEU:HD11	1.99	0.44
1:B:196:PHE:CE2	4:N:1:NAG:H5	2.52	0.44
1:A:451:THR:O	1:A:454:THR:HG22	2.16	0.44
1:D:285:VAL:C	1:D:287:PHE:H	2.25	0.44
2:I:52:THR:O	2:I:61:SER:OG	2.28	0.44
1:E:251:ARG:HH11	1:E:301:THR:HG23	1.81	0.44
1:B:159:THR:HG22	1:B:207:TYR:HE1	1.83	0.44
1:D:61:TYR:CE1	1:D:128:PRO:HB2	2.53	0.44
1:E:88:ILE:HD11	1:E:117:VAL:HG21	2.00	0.44
1:E:281:PRO:C	1:E:283:ILE:H	2.25	0.44
1:A:102:GLU:N	1:A:102:GLU:OE2	2.50	0.44
1:D:431:TYR:OH	1:E:311:GLN:OE1	2.29	0.44
1:E:77:PRO:HG2	1:E:78:GLU:OE2	2.17	0.44
1:E:266:LEU:HD21	1:E:291:MET:HE3	2.00	0.44
1:B:263:SER:O	1:B:266:LEU:HG	2.17	0.44
1:C:203:ILE:HG23	2:H:76:ARG:HD2	2.00	0.44
1:D:102:GLU:OE2	1:D:102:GLU:N	2.48	0.43
1:D:118:HIS:CE1	2:J:28:LEU:HD12	2.53	0.43
1:E:447:ALA:O	1:E:450:ILE:HG22	2.18	0.43
1:B:235:MET:O	1:B:239:ILE:HG12	2.19	0.43
1:B:251:ARG:HH11	1:B:301:THR:HG23	1.81	0.43
1:D:170:SER:HB3	1:D:173:GLU:HG3	2.00	0.43
1:D:251:ARG:HD3	1:D:301:THR:HG23	2.00	0.43
1:E:173:GLU:OE1	2:F:105:ARG:NH2	2.52	0.43
1:B:188:GLU:HB2	1:B:219:ARG:HD3	2.00	0.43
1:C:235:MET:O	1:C:239:ILE:HG12	2.19	0.43
1:D:61:TYR:CD2	1:E:104:VAL:HA	2.53	0.43
1:E:451:THR:O	1:E:454:THR:HG22	2.19	0.43
4:W:1:NAG:H61	4:W:2:NAG:HN2	1.82	0.43
1:B:280:THR:HA	1:B:281:PRO:HD3	1.85	0.43
1:E:35:LYS:HA	1:E:36:PRO:HD3	1.87	0.43
1:C:81:ASP:OD2	1:D:33:TRP:NE1	2.48	0.43
1:D:166:THR:HG22	4:T:1:NAG:H62	1.98	0.43
1:E:102:GLU:OE2	1:E:102:GLU:N	2.49	0.43
1:E:235:MET:O	1:E:239:ILE:HG12	2.18	0.43
2:F:52:THR:O	2:F:61:SER:OG	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:ILE:HG23	2:F:76:ARG:HH11	1.84	0.43
1:D:410:ARG:NH1	1:E:405:GLU:OE1	2.49	0.43
2:G:64:TYR:CZ	2:G:74:ILE:HG22	2.54	0.43
1:A:118:HIS:CE1	2:G:28:LEU:HD12	2.53	0.43
1:A:169:ARG:NH2	1:A:177:ASP:OD2	2.52	0.43
1:C:263:SER:O	1:C:266:LEU:HG	2.19	0.43
1:E:62:ILE:HD13	1:E:131:LEU:HD12	1.99	0.43
1:E:263:SER:O	1:E:266:LEU:HG	2.18	0.43
1:C:77:PRO:HG2	1:C:78:GLU:OE2	2.19	0.43
1:C:281:PRO:C	1:C:283:ILE:H	2.26	0.43
1:D:149:CYS:HB2	1:D:214:VAL:HG13	2.00	0.43
1:E:78:GLU:C	1:E:80:PHE:H	2.27	0.43
1:A:447:ALA:O	1:A:450:ILE:HG22	2.19	0.43
1:B:62:ILE:HD13	1:B:131:LEU:HD12	2.01	0.43
1:C:62:ILE:HD13	1:C:131:LEU:HD12	2.00	0.43
1:D:229:LEU:HB3	1:D:230:PRO:HD3	2.01	0.43
1:A:229:LEU:HB3	1:A:230:PRO:HD3	1.99	0.43
1:D:280:THR:HA	1:D:281:PRO:HD3	1.85	0.43
2:G:87:MET:HB3	2:G:90:LEU:HD21	2.01	0.43
1:A:298:LEU:HD21	1:E:254:PHE:HE2	1.84	0.42
1:D:281:PRO:C	1:D:283:ILE:H	2.27	0.42
1:B:78:GLU:C	1:B:80:PHE:H	2.26	0.42
2:G:71:ARG:NH1	2:G:91:LYS:HE3	2.34	0.42
1:B:273:LEU:HD13	1:B:274:PRO:HD2	2.00	0.42
1:C:197:LYS:HD3	2:H:58:TYR:CZ	2.54	0.42
1:C:447:ALA:O	1:C:450:ILE:HG22	2.19	0.42
1:B:447:ALA:O	1:B:450:ILE:HG22	2.20	0.42
1:E:229:LEU:HB3	1:E:230:PRO:HD3	2.01	0.42
2:I:76:ARG:HH12	2:I:78:SER:HB3	1.85	0.42
1:C:254:PHE:HE2	1:D:298:LEU:HD21	1.84	0.42
1:E:167:LEU:HD11	1:E:194:PRO:HB2	2.02	0.42
1:B:115:VAL:HG12	1:B:125:ASN:HA	2.01	0.42
1:C:147:GLN:HB2	1:C:216:ILE:HG13	2.02	0.42
1:C:159:THR:HG22	1:C:207:TYR:HE1	1.85	0.42
1:D:447:ALA:O	1:D:450:ILE:HG22	2.19	0.42
1:E:166:THR:HG22	4:W:1:NAG:H62	2.00	0.42
1:A:235:MET:O	1:A:239:ILE:HG12	2.19	0.42
1:A:245:PRO:HD3	1:B:306:ARG:HG3	2.02	0.42
1:A:281:PRO:C	1:A:283:ILE:H	2.27	0.42
2:J:71:ARG:HB3	2:J:88:ASN:O	2.19	0.42
1:A:113:PRO:HG2	1:A:114:TYR:CD2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:166:THR:HG22	4:Q:1:NAG:H62	2.02	0.42
1:A:293:LEU:O	1:A:296:ILE:HG13	2.19	0.42
1:B:165:ILE:HG22	1:B:208:ALA:HB1	2.02	0.42
1:B:201:ILE:HD12	1:B:201:ILE:HA	1.94	0.42
1:C:315:ARG:NH2	1:C:418:GLU:OE2	2.51	0.42
1:C:450:ILE:HA	1:C:453:VAL:HG22	2.02	0.42
1:E:219:ARG:HA	1:E:220:PRO:HD3	1.86	0.42
1:C:251:ARG:HH11	1:C:301:THR:HG23	1.84	0.42
1:D:9:GLN:N	1:D:10:PRO:HD3	2.35	0.42
1:D:235:MET:O	1:D:239:ILE:HG12	2.18	0.42
1:D:293:LEU:O	1:D:296:ILE:HG13	2.20	0.42
1:B:229:LEU:HB3	1:B:230:PRO:HD3	2.01	0.41
1:B:315:ARG:NH2	1:B:418:GLU:OE2	2.51	0.41
1:E:140:TYR:CZ	1:E:282:LEU:HD22	2.55	0.41
1:E:175:ARG:NE	1:E:191:GLU:OE1	2.52	0.41
2:H:71:ARG:HB3	2:H:88:ASN:O	2.19	0.41
2:H:124:SER:OG	6:H:302:SO4:O2	2.37	0.41
1:C:53:GLU:HG2	1:C:273:LEU:HD11	2.02	0.41
1:D:78:GLU:C	1:D:80:PHE:H	2.27	0.41
1:E:229:LEU:O	1:E:232:ILE:HG13	2.20	0.41
2:I:66:ASP:HA	2:I:69:LYS:HD3	2.02	0.41
2:J:66:ASP:HA	2:J:69:LYS:HD3	2.01	0.41
1:D:109:SER:OG	1:D:130:GLN:N	2.49	0.41
1:B:293:LEU:O	1:B:296:ILE:HG13	2.20	0.41
1:C:233:PHE:O	1:C:237:VAL:HG23	2.20	0.41
1:D:263:SER:O	1:D:266:LEU:HG	2.19	0.41
1:A:9:GLN:N	1:A:10:PRO:HD3	2.34	0.41
1:D:19:LEU:HD23	1:D:73:LEU:HD22	2.02	0.41
1:D:201:ILE:HD12	1:D:201:ILE:HA	1.94	0.41
2:J:87:MET:HB3	2:J:90:LEU:HD21	2.02	0.41
2:H:64:TYR:CZ	2:H:74:ILE:HG22	2.56	0.41
1:A:78:GLU:C	1:A:80:PHE:H	2.28	0.41
1:A:263:SER:O	1:A:266:LEU:HG	2.20	0.41
1:B:109:SER:HA	1:B:110:PRO:HD3	1.93	0.41
1:C:9:GLN:N	1:C:10:PRO:HD3	2.35	0.41
1:A:233:PHE:O	1:A:237:VAL:HG23	2.20	0.41
1:C:256:ILE:HD13	1:C:298:LEU:HD11	2.03	0.41
1:A:170:SER:HB3	1:A:173:GLU:HG3	2.03	0.41
1:C:113:PRO:HG2	1:C:114:TYR:CD2	2.56	0.41
1:C:143:PRO:HG2	1:C:144:PHE:CE1	2.56	0.41
1:D:213:TYR:CE1	3:V:1:NAG:H5	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:233:PHE:O	1:D:237:VAL:HG23	2.20	0.41
1:E:9:GLN:N	1:E:10:PRO:HD3	2.35	0.41
1:E:256:ILE:HD13	1:E:298:LEU:HD11	2.03	0.41
1:A:149:CYS:HB2	1:A:214:VAL:HG13	2.02	0.41
1:B:9:GLN:N	1:B:10:PRO:HD3	2.36	0.41
1:E:233:PHE:O	1:E:237:VAL:HG23	2.21	0.41
1:B:170:SER:HB3	1:B:173:GLU:HG3	2.03	0.40
1:C:140:TYR:CZ	1:C:282:LEU:HD22	2.57	0.40
1:D:450:ILE:HA	1:D:453:VAL:HG22	2.02	0.40
1:A:62:ILE:HD13	1:A:131:LEU:HD12	2.02	0.40
1:B:254:PHE:HE2	1:C:298:LEU:HD21	1.86	0.40
1:B:431:TYR:OH	1:C:311:GLN:OE1	2.31	0.40
1:D:401:GLY:O	1:D:404:GLN:HG2	2.22	0.40
2:J:12:VAL:HG21	2:J:90:LEU:HD13	2.03	0.40
1:C:10:PRO:HD2	1:C:14:ARG:HD3	2.03	0.40
1:A:109:SER:OG	1:A:130:GLN:N	2.46	0.40
1:C:229:LEU:HB3	1:C:230:PRO:HD3	2.03	0.40
1:D:285:VAL:HA	1:D:288:VAL:HG12	2.04	0.40
2:H:40:ALA:HB3	2:H:43:LYS:HB2	2.04	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	377/456 (83%)	352 (93%)	24 (6%)	1 (0%)	36 67
1	B	377/456 (83%)	352 (93%)	24 (6%)	1 (0%)	36 67
1	C	378/456 (83%)	353 (93%)	24 (6%)	1 (0%)	36 67
1	D	378/456 (83%)	352 (93%)	24 (6%)	2 (0%)	24 57
1	E	377/456 (83%)	352 (93%)	24 (6%)	1 (0%)	36 67

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	120/124 (97%)	116 (97%)	4 (3%)	0	100	100
2	G	122/124 (98%)	117 (96%)	5 (4%)	0	100	100
2	H	122/124 (98%)	118 (97%)	4 (3%)	0	100	100
2	I	122/124 (98%)	117 (96%)	5 (4%)	0	100	100
2	J	122/124 (98%)	118 (97%)	4 (3%)	0	100	100
All	All	2495/2900 (86%)	2347 (94%)	142 (6%)	6 (0%)	43	74

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	79	ASP
1	A	79	ASP
1	B	79	ASP
1	D	79	ASP
1	E	79	ASP
1	D	274	PRO

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	341/417 (82%)	324 (95%)	17 (5%)	22	48
1	B	342/417 (82%)	326 (95%)	16 (5%)	23	49
1	C	343/417 (82%)	327 (95%)	16 (5%)	23	49
1	D	343/417 (82%)	328 (96%)	15 (4%)	25	51
1	E	342/417 (82%)	326 (95%)	16 (5%)	23	49
2	F	94/99 (95%)	92 (98%)	2 (2%)	47	66
2	G	97/99 (98%)	95 (98%)	2 (2%)	47	66
2	H	97/99 (98%)	95 (98%)	2 (2%)	47	66
2	I	97/99 (98%)	95 (98%)	2 (2%)	47	66
2	J	97/99 (98%)	95 (98%)	2 (2%)	47	66

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2193/2580 (85%)	2103 (96%)	90 (4%)	27 53

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	54	LYS
1	A	57	VAL
1	A	62	ILE
1	A	116	TYR
1	A	139	ILE
1	A	144	PHE
1	A	156	TRP
1	A	162	ASP
1	A	214	VAL
1	A	225	VAL
1	A	273	LEU
1	A	293	LEU
1	A	305	VAL
1	A	308	VAL
1	A	310	LYS
1	A	431	TYR
1	B	9	GLN
1	B	54	LYS
1	B	57	VAL
1	B	62	ILE
1	B	116	TYR
1	B	139	ILE
1	B	144	PHE
1	B	156	TRP
1	B	162	ASP
1	B	225	VAL
1	B	273	LEU
1	B	293	LEU
1	B	305	VAL
1	B	308	VAL
1	B	310	LYS
1	B	431	TYR
1	C	9	GLN
1	C	54	LYS
1	C	57	VAL
1	C	62	ILE

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Mol	Chain	Res	Type
1	C	139	ILE
1	C	144	PHE
1	C	156	TRP
1	C	162	ASP
1	C	214	VAL
1	C	225	VAL
1	C	273	LEU
1	C	293	LEU
1	C	305	VAL
1	C	308	VAL
1	C	310	LYS
1	C	431	TYR
1	D	9	GLN
1	D	54	LYS
1	D	57	VAL
1	D	62	ILE
1	D	139	ILE
1	D	144	PHE
1	D	156	TRP
1	D	214	VAL
1	D	225	VAL
1	D	273	LEU
1	D	293	LEU
1	D	305	VAL
1	D	308	VAL
1	D	310	LYS
1	D	431	TYR
1	E	9	GLN
1	E	54	LYS
1	E	57	VAL
1	E	62	ILE
1	E	139	ILE
1	E	144	PHE
1	E	156	TRP
1	E	165	ILE
1	E	214	VAL
1	E	225	VAL
1	E	273	LEU
1	E	293	LEU
1	E	305	VAL
1	E	308	VAL
1	E	310	LYS

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Mol	Chain	Res	Type
1	E	431	TYR
2	F	2	VAL
2	F	28	LEU
2	G	2	VAL
2	G	28	LEU
2	H	2	VAL
2	H	28	LEU
2	I	2	VAL
2	I	28	LEU
2	J	2	VAL
2	J	28	LEU

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

Mol	Chain	Res	Type
1	B	205	ASN
1	C	66	GLN
1	C	205	ASN
1	D	205	ASN
1	E	124	GLN
1	E	205	ASN
2	F	81	ASN
2	G	81	ASN
2	G	88	ASN
2	G	101	ASN
2	H	81	ASN
2	I	81	ASN
2	I	101	ASN
2	J	81	ASN
2	J	86	GLN
2	J	101	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 ⓘ

33 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
3	NAG	K	1	3,1	14,14,15	0.26	0	17,19,21	0.49	0
3	NAG	K	2	3	14,14,15	0.72	1 (7%)	17,19,21	1.02	1 (5%)
4	NAG	L	1	4,1	14,14,15	0.26	0	17,19,21	0.40	0
4	NAG	L	2	4	14,14,15	0.22	0	17,19,21	0.36	0
4	BMA	L	3	4	11,11,12	0.61	0	15,15,17	0.72	0
3	NAG	M	1	3,1	14,14,15	0.27	0	17,19,21	0.46	0
3	NAG	M	2	3	14,14,15	0.24	0	17,19,21	0.46	0
4	NAG	N	1	4,1	14,14,15	0.22	0	17,19,21	0.52	0
4	NAG	N	2	4	14,14,15	0.23	0	17,19,21	0.42	0
4	BMA	N	3	4	11,11,12	0.58	0	15,15,17	0.73	0
3	NAG	O	1	3,1	14,14,15	0.25	0	17,19,21	0.47	0
3	NAG	O	2	3	14,14,15	0.71	1 (7%)	17,19,21	1.00	1 (5%)
3	NAG	P	1	3,1	14,14,15	0.31	0	17,19,21	0.48	0
3	NAG	P	2	3	14,14,15	0.27	0	17,19,21	0.46	0
4	NAG	Q	1	4,1	14,14,15	0.27	0	17,19,21	0.41	0
4	NAG	Q	2	4	14,14,15	0.20	0	17,19,21	0.39	0
4	BMA	Q	3	4	11,11,12	0.58	0	15,15,17	0.71	0
3	NAG	R	1	3,1	14,14,15	0.25	0	17,19,21	0.45	0
3	NAG	R	2	3	14,14,15	0.71	1 (7%)	17,19,21	1.01	1 (5%)
3	NAG	S	1	3,1	14,14,15	0.30	0	17,19,21	0.45	0
3	NAG	S	2	3	14,14,15	0.24	0	17,19,21	0.46	0
4	NAG	T	1	4,1	14,14,15	0.26	0	17,19,21	0.36	0
4	NAG	T	2	4	14,14,15	0.20	0	17,19,21	0.43	0
4	BMA	T	3	4	11,11,12	0.83	0	15,15,17	1.28	2 (13%)
3	NAG	U	1	3,1	14,14,15	0.29	0	17,19,21	0.48	0
3	NAG	U	2	3	14,14,15	0.72	1 (7%)	17,19,21	1.01	1 (5%)
3	NAG	V	1	3,1	14,14,15	0.30	0	17,19,21	0.49	0
3	NAG	V	2	3	14,14,15	0.25	0	17,19,21	0.45	0
4	NAG	W	1	4,1	14,14,15	0.26	0	17,19,21	0.41	0
4	NAG	W	2	4	14,14,15	0.20	0	17,19,21	0.41	0
4	BMA	W	3	4	11,11,12	1.37	2 (18%)	15,15,17	1.33	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	X	1	3,1	14,14,15	0.28	0	17,19,21	0.50	0
3	NAG	X	2	3	14,14,15	0.25	0	17,19,21	0.46	0

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
3	NAG	K	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	K	2	3	-	0/6/23/26	0/1/1/1
4	NAG	L	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	L	2	4	-	0/6/23/26	0/1/1/1
4	BMA	L	3	4	-	0/2/19/22	0/1/1/1
3	NAG	M	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	M	2	3	-	1/6/23/26	0/1/1/1
4	NAG	N	1	4,1	-	1/6/23/26	0/1/1/1
4	NAG	N	2	4	-	2/6/23/26	0/1/1/1
4	BMA	N	3	4	-	0/2/19/22	0/1/1/1
3	NAG	O	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	O	2	3	-	2/6/23/26	0/1/1/1
3	NAG	P	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	P	2	3	-	1/6/23/26	0/1/1/1
4	NAG	Q	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	Q	2	4	-	2/6/23/26	0/1/1/1
4	BMA	Q	3	4	-	0/2/19/22	0/1/1/1
3	NAG	R	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	R	2	3	-	1/6/23/26	0/1/1/1
3	NAG	S	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	S	2	3	-	0/6/23/26	0/1/1/1
4	NAG	T	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	T	2	4	-	2/6/23/26	0/1/1/1
4	BMA	T	3	4	-	0/2/19/22	0/1/1/1
3	NAG	U	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	U	2	3	-	2/6/23/26	0/1/1/1
3	NAG	V	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	V	2	3	-	0/6/23/26	0/1/1/1
4	NAG	W	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	W	2	4	-	2/6/23/26	0/1/1/1
4	BMA	W	3	4	-	0/2/19/22	0/1/1/1
3	NAG	X	1	3,1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	X	2	3	-	1/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	W	3	BMA	O5-C1	-2.56	1.39	1.43
3	U	2	NAG	O5-C1	2.53	1.47	1.43
3	K	2	NAG	O5-C1	2.51	1.47	1.43
4	W	3	BMA	C4-C3	2.49	1.58	1.52
3	O	2	NAG	O5-C1	2.45	1.47	1.43
3	R	2	NAG	O5-C1	2.44	1.47	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	2	NAG	C1-O5-C5	3.96	117.49	112.19
3	U	2	NAG	C1-O5-C5	3.96	117.49	112.19
3	R	2	NAG	C1-O5-C5	3.94	117.47	112.19
3	O	2	NAG	C1-O5-C5	3.90	117.41	112.19
4	W	3	BMA	C2-C3-C4	3.33	116.72	110.86
4	W	3	BMA	C3-C4-C5	3.10	115.85	110.23
4	T	3	BMA	C1-O5-C5	2.60	115.67	112.19
4	T	3	BMA	C1-C2-C3	2.11	112.72	109.64

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	U	1	NAG	O5-C5-C6-O6
4	N	2	NAG	O5-C5-C6-O6
3	R	1	NAG	O5-C5-C6-O6
4	W	2	NAG	O5-C5-C6-O6
3	O	2	NAG	C4-C5-C6-O6
3	U	1	NAG	C4-C5-C6-O6
3	R	1	NAG	C4-C5-C6-O6
3	O	1	NAG	O5-C5-C6-O6
4	T	2	NAG	O5-C5-C6-O6
3	U	2	NAG	C4-C5-C6-O6
4	N	2	NAG	C4-C5-C6-O6
3	O	1	NAG	C4-C5-C6-O6
4	W	2	NAG	C4-C5-C6-O6

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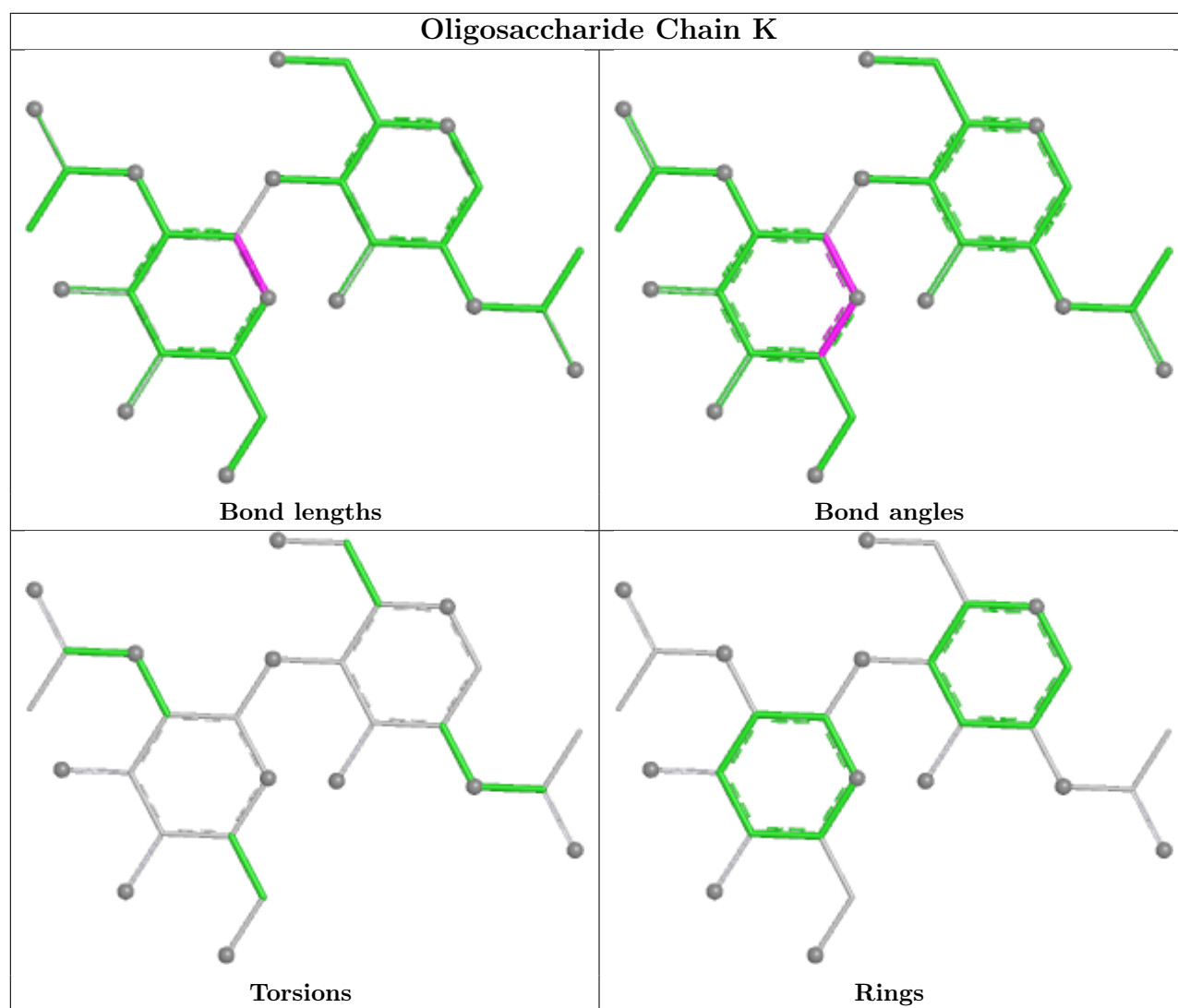
Mol	Chain	Res	Type	Atoms
3	O	2	NAG	O5-C5-C6-O6
4	T	2	NAG	C4-C5-C6-O6
3	U	2	NAG	O5-C5-C6-O6
4	Q	2	NAG	C4-C5-C6-O6
3	M	2	NAG	O5-C5-C6-O6
4	Q	2	NAG	O5-C5-C6-O6
3	V	1	NAG	C4-C5-C6-O6
3	P	2	NAG	O5-C5-C6-O6
3	P	1	NAG	C4-C5-C6-O6
3	R	2	NAG	C4-C5-C6-O6
3	X	2	NAG	O5-C5-C6-O6
4	N	1	NAG	C4-C5-C6-O6

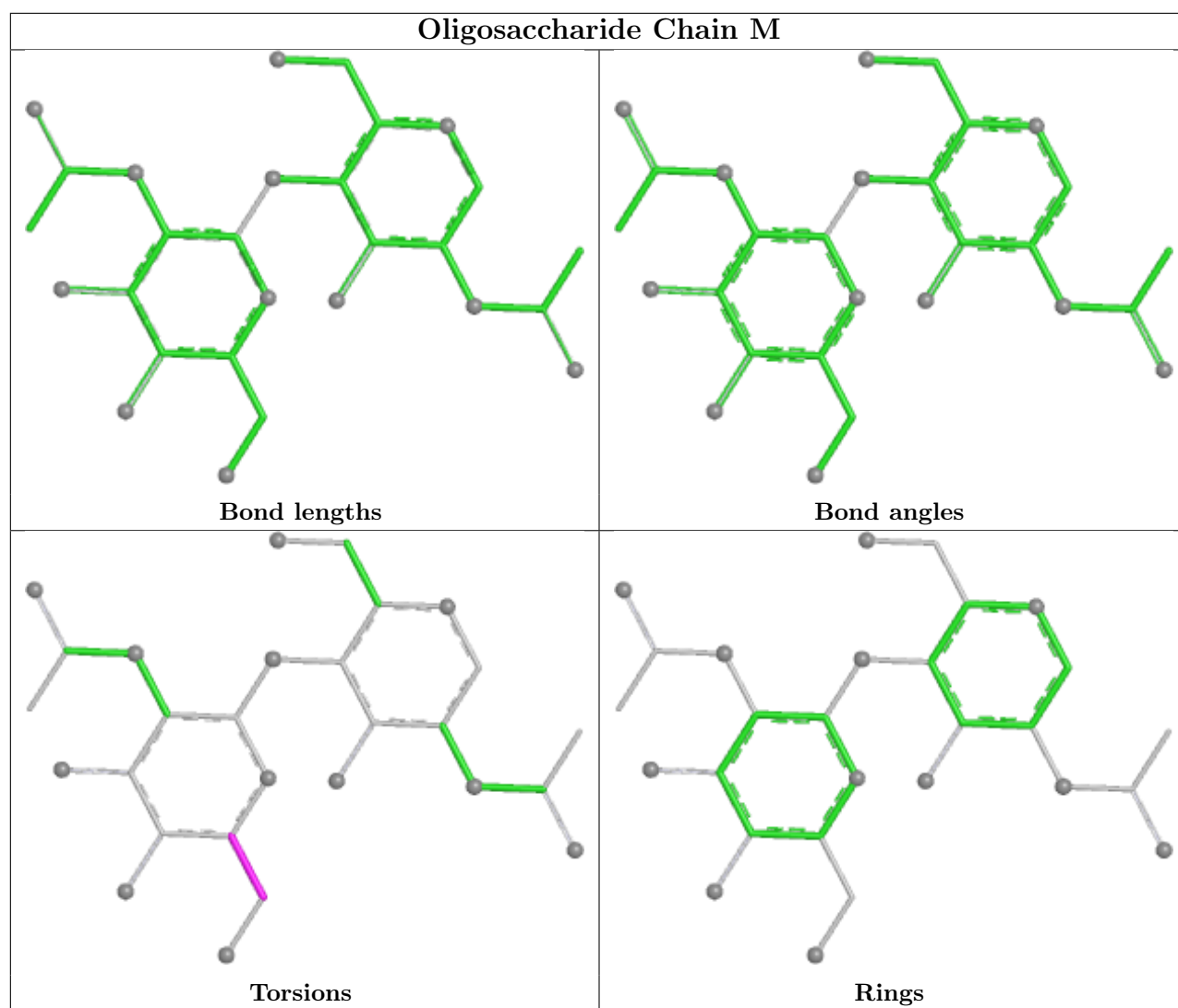
There are no ring outliers.

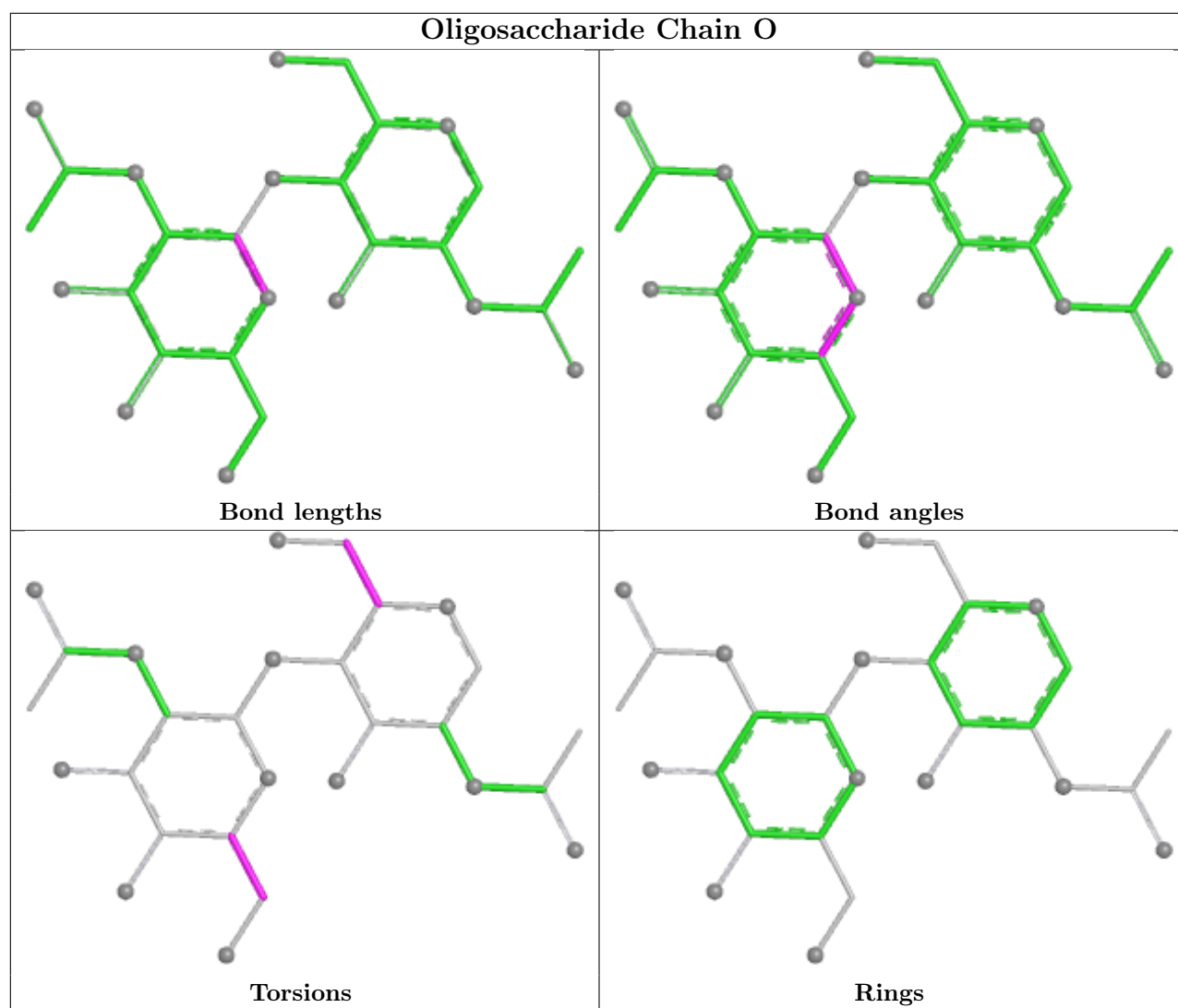
11 monomers are involved in 12 short contacts:

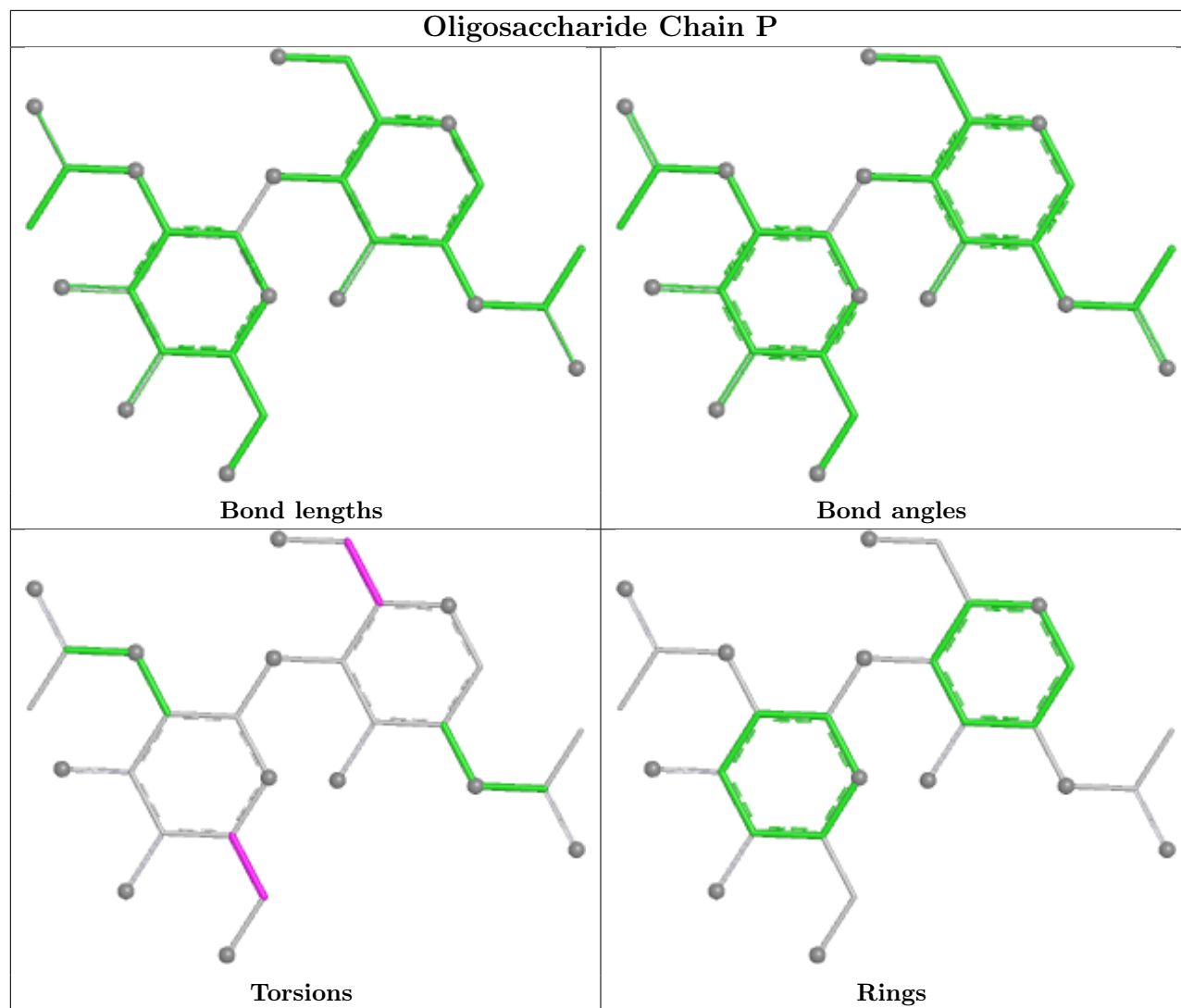
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	1	NAG	1	0
4	Q	1	NAG	1	0
3	M	1	NAG	1	0
4	N	1	NAG	1	0
3	V	1	NAG	2	0
4	T	1	NAG	1	0
3	S	1	NAG	1	0
4	W	1	NAG	2	0
3	X	1	NAG	1	0
4	W	2	NAG	1	0
3	P	1	NAG	1	0

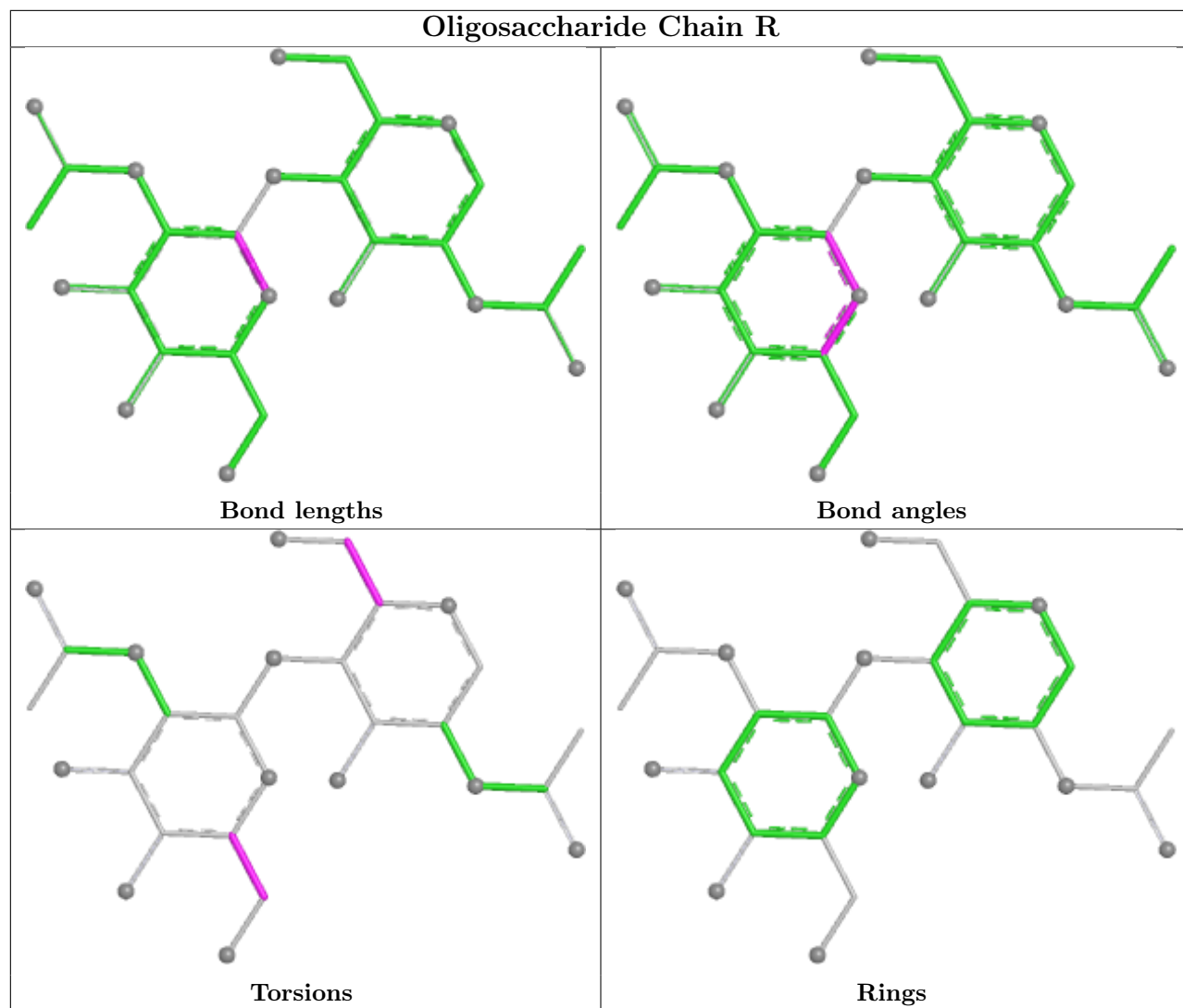
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

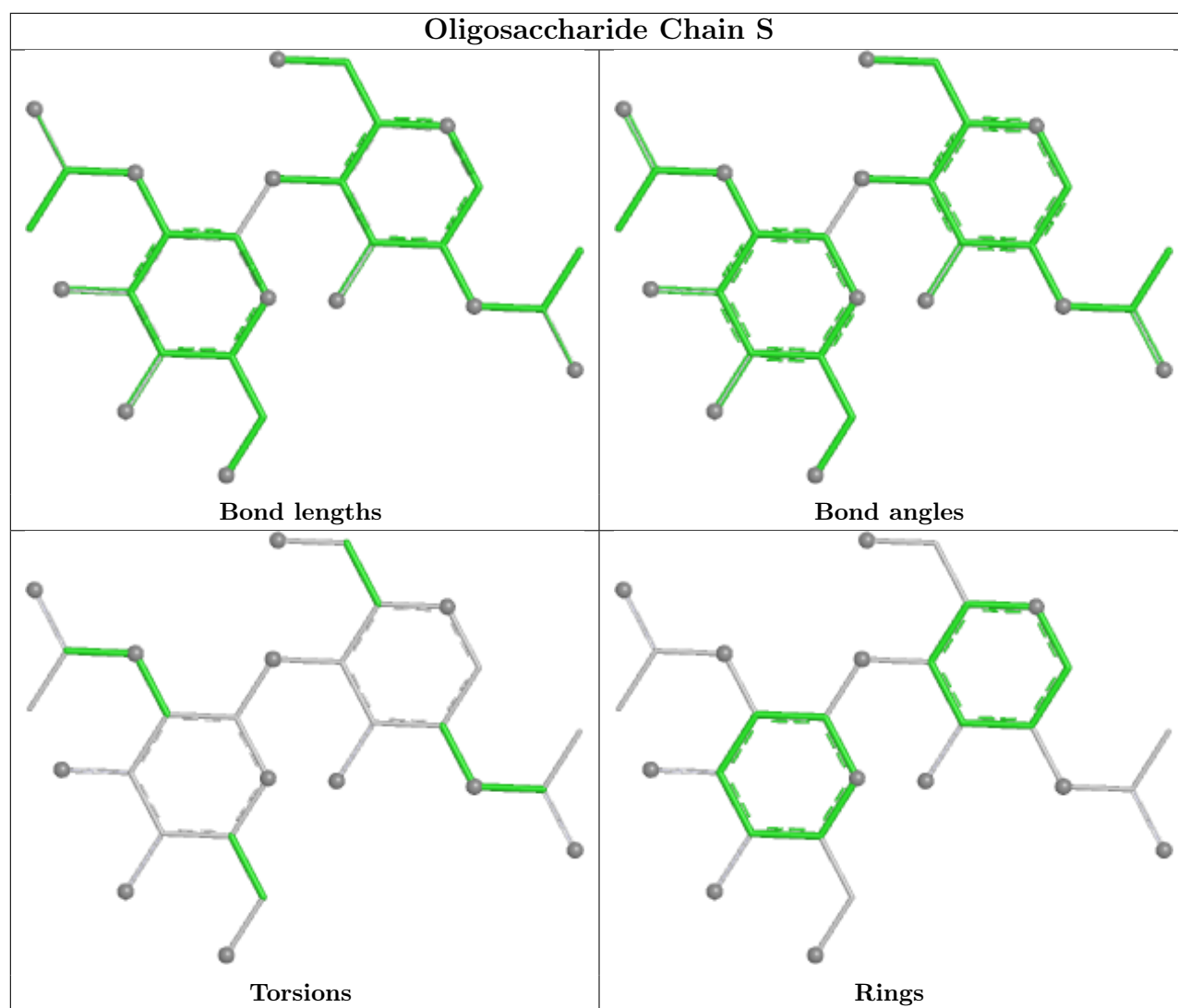


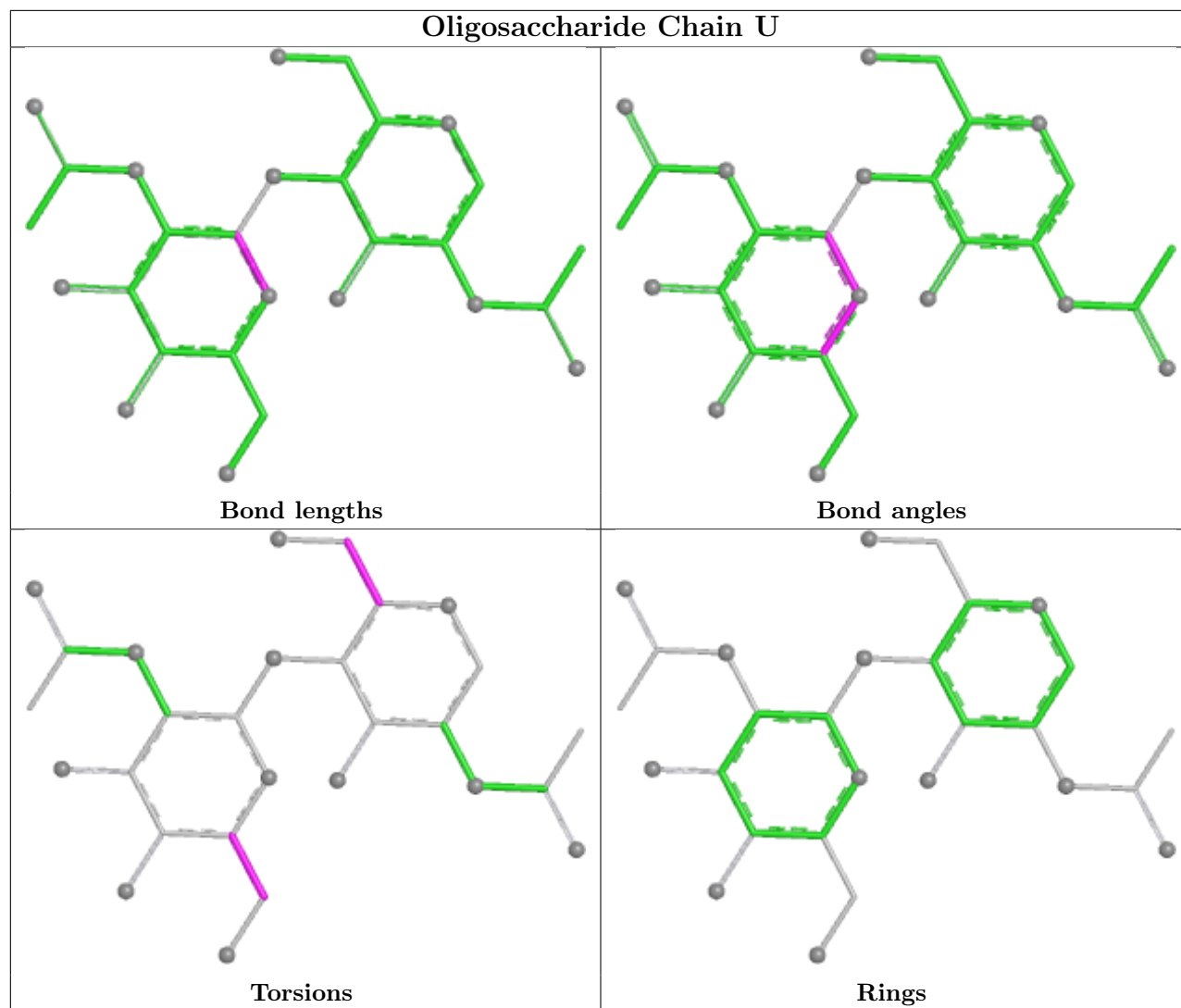


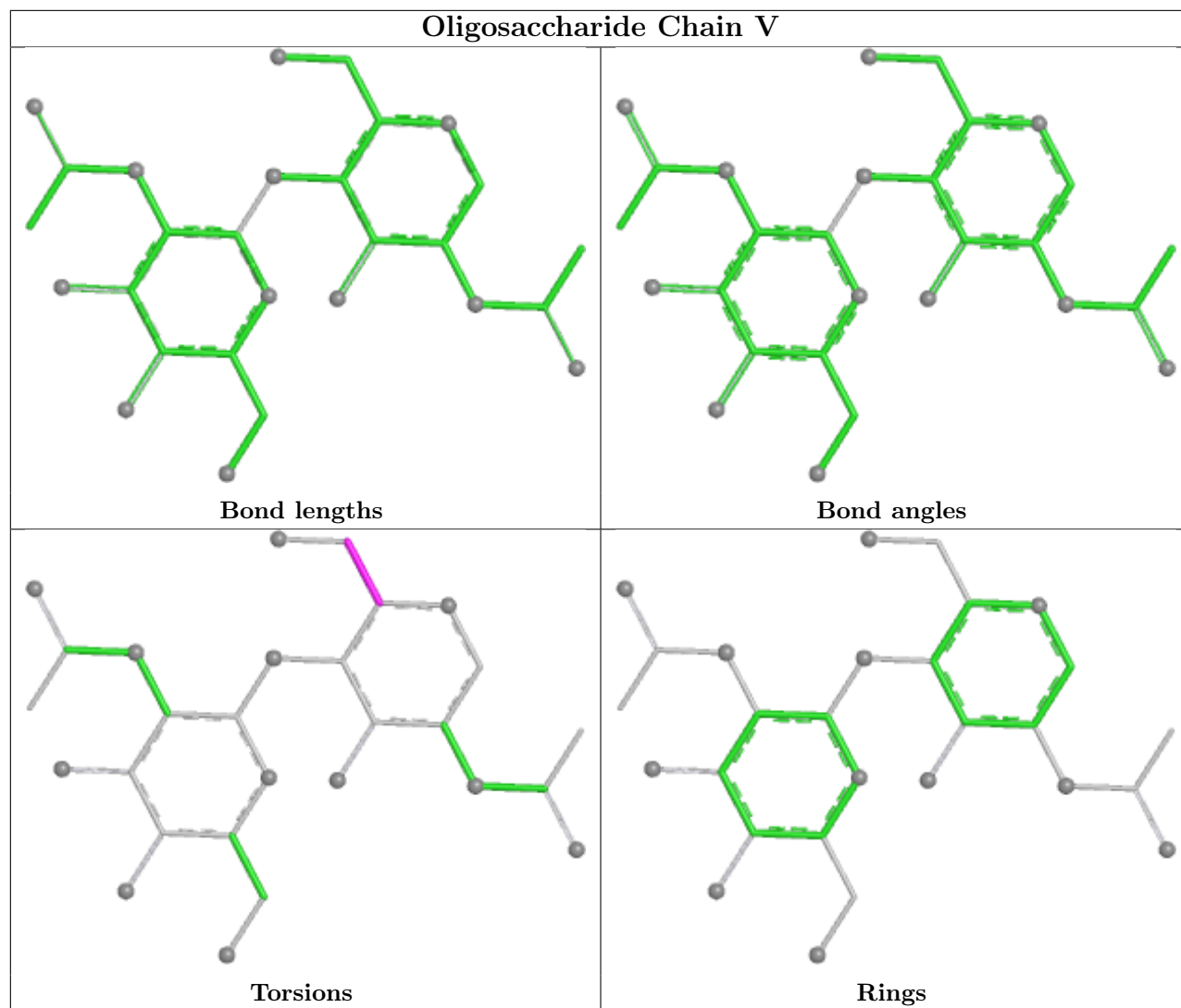


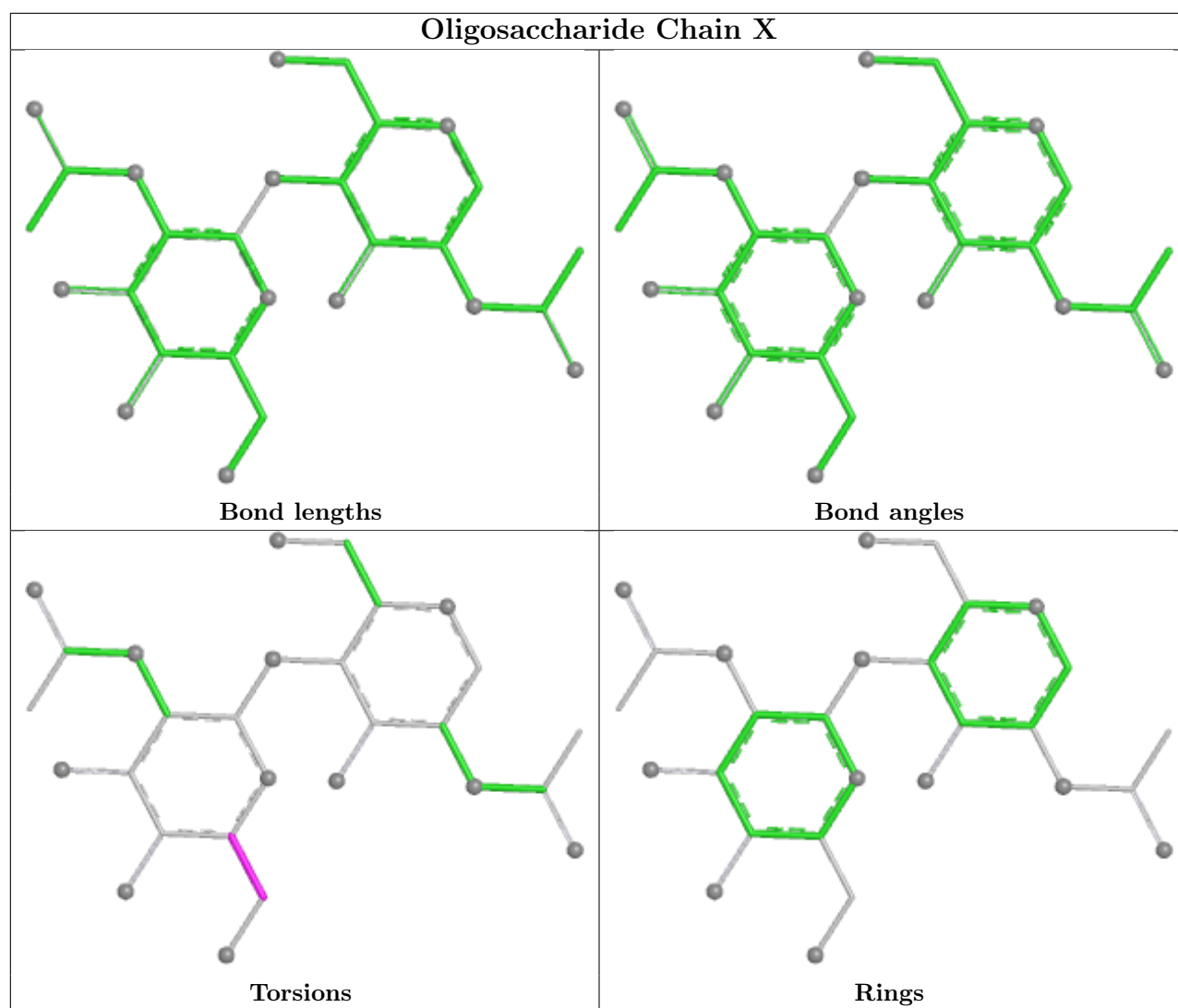


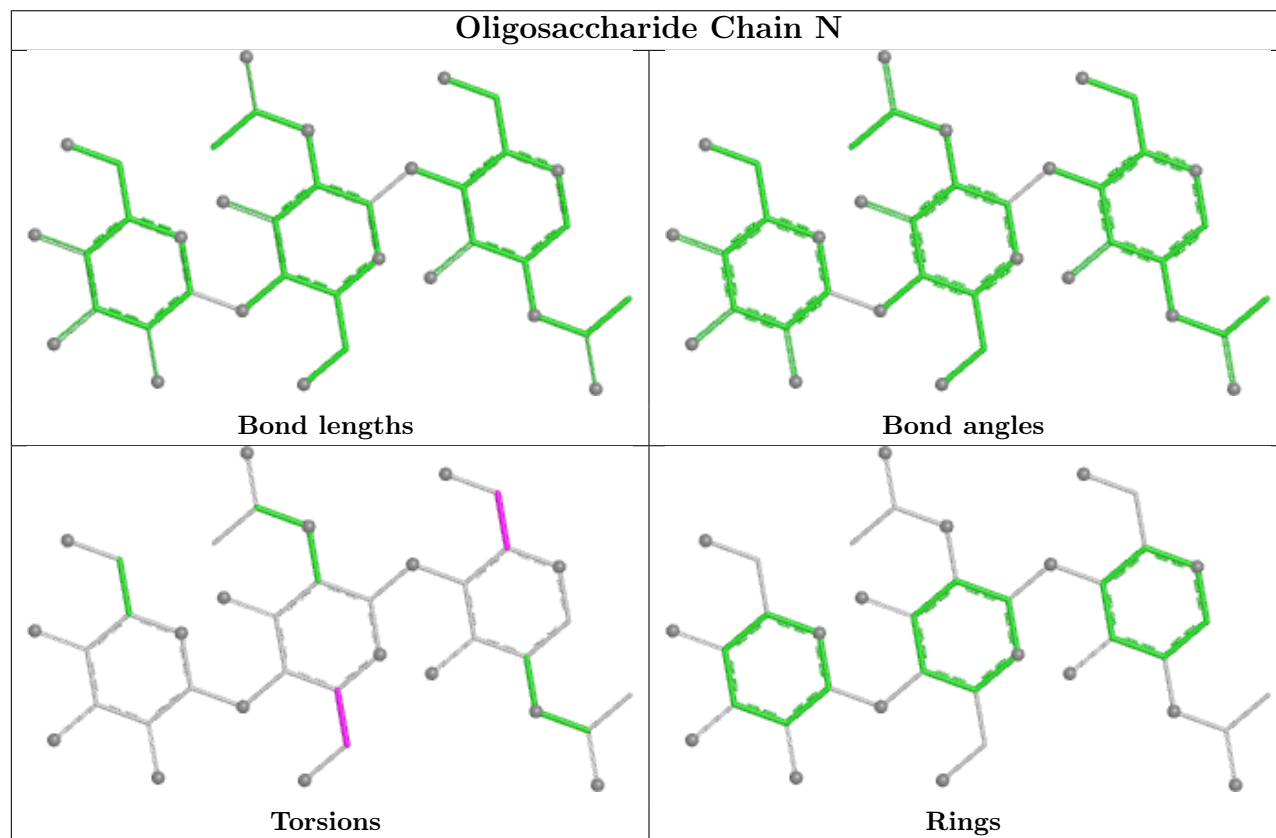
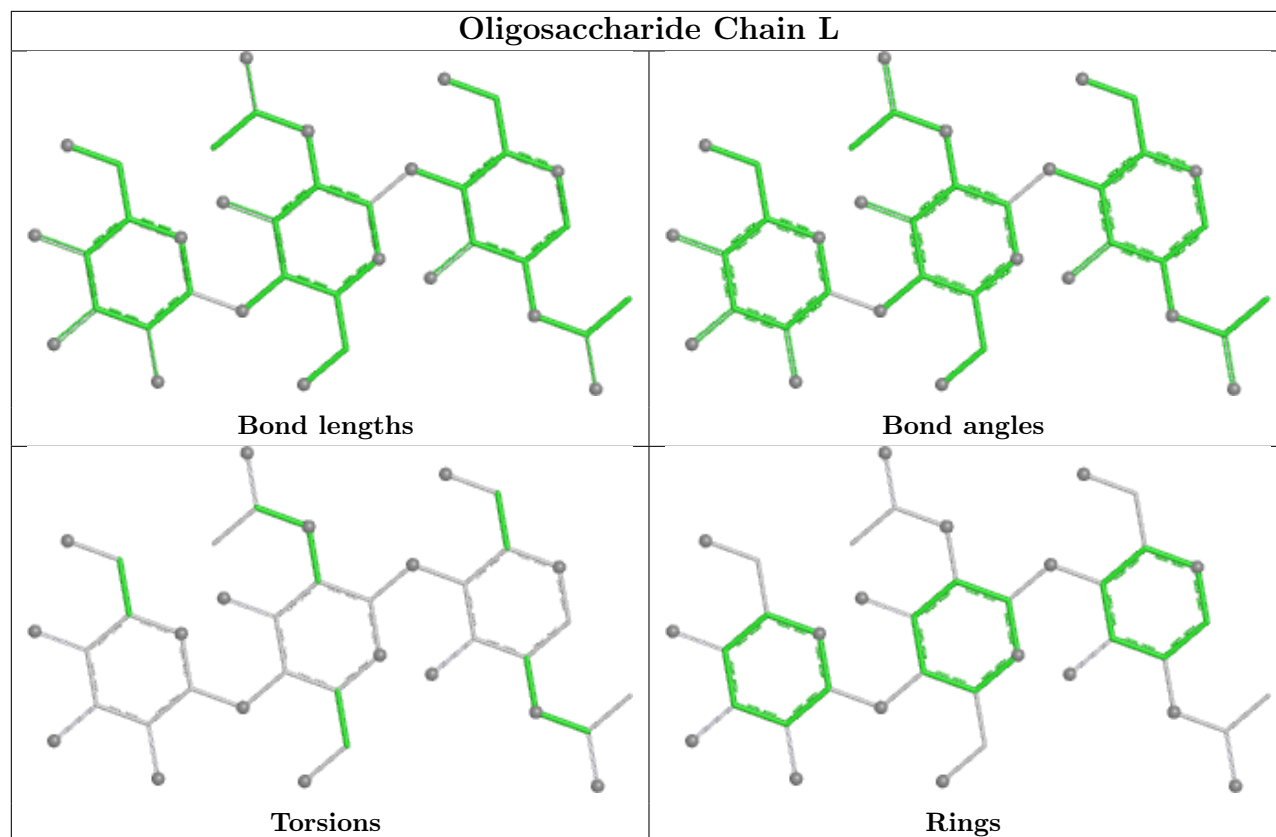


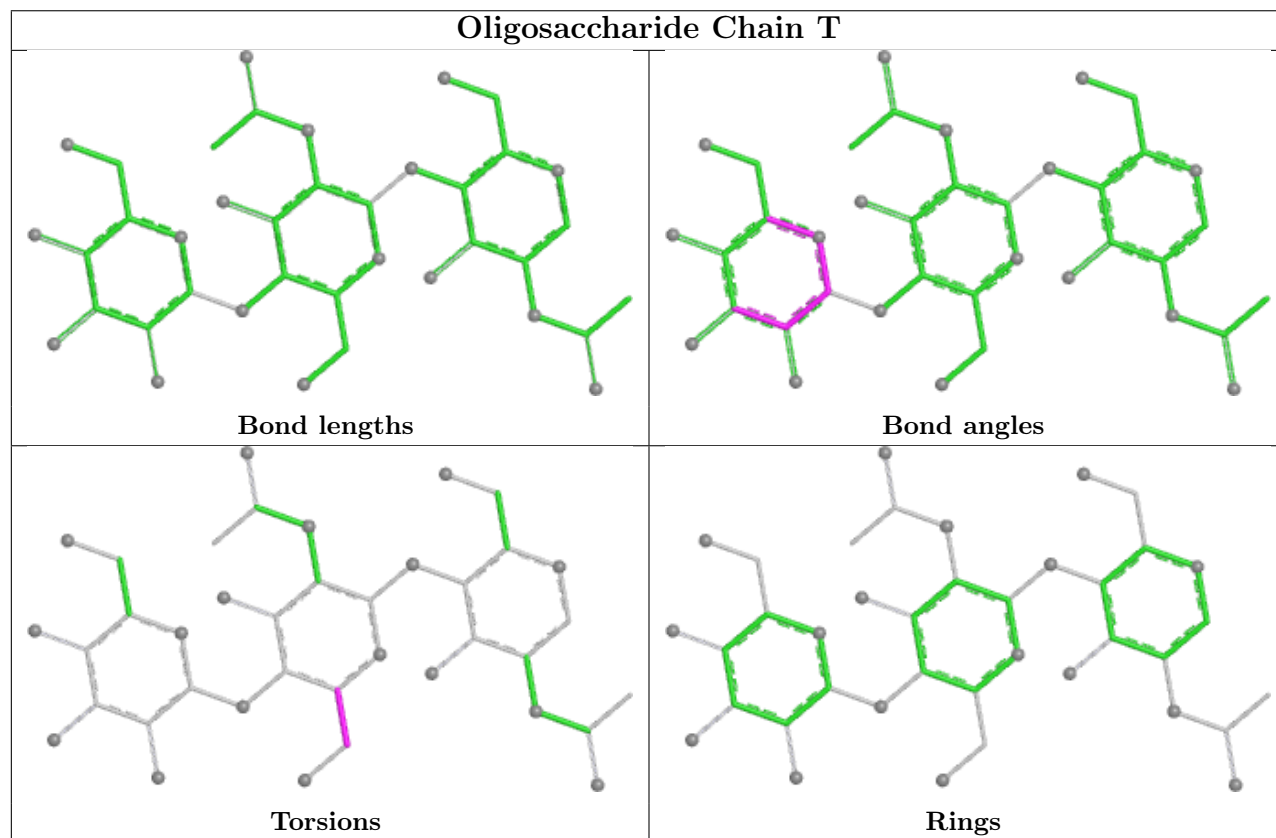
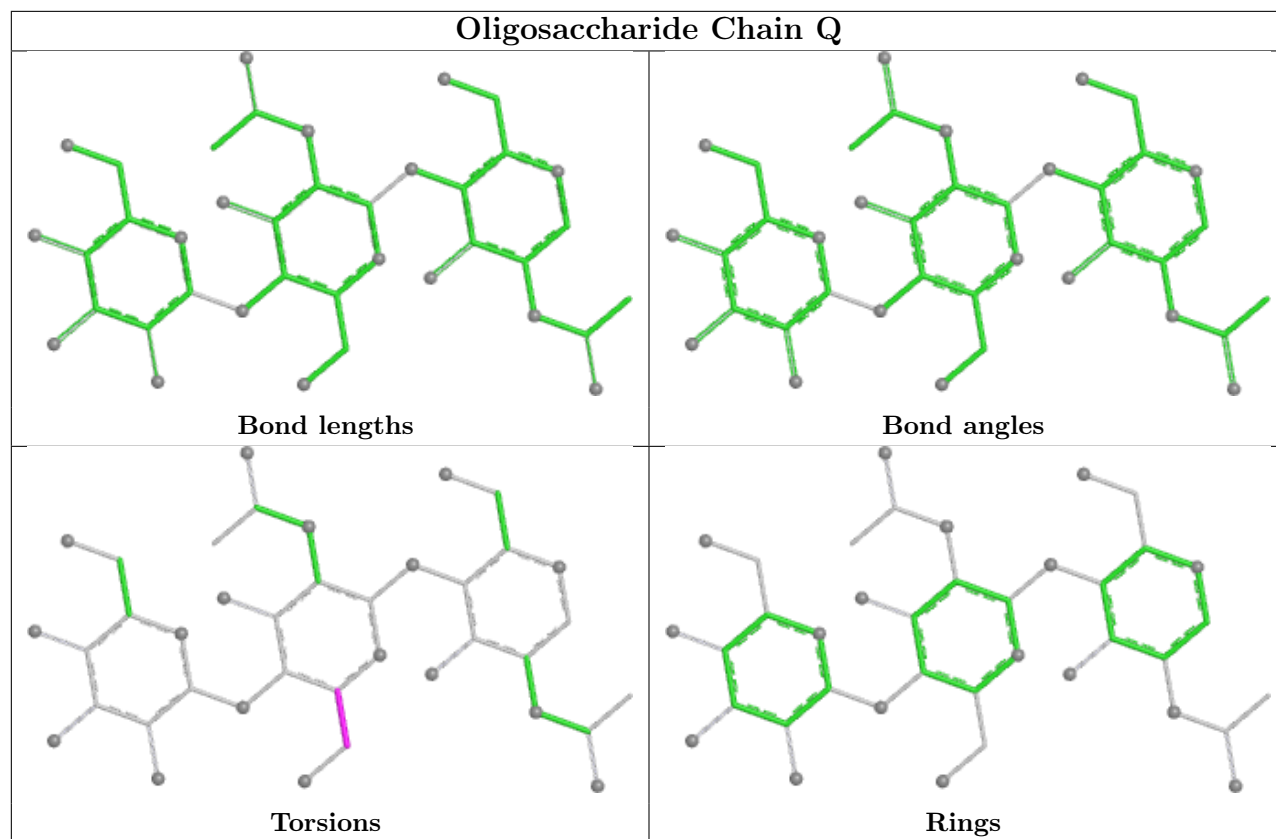


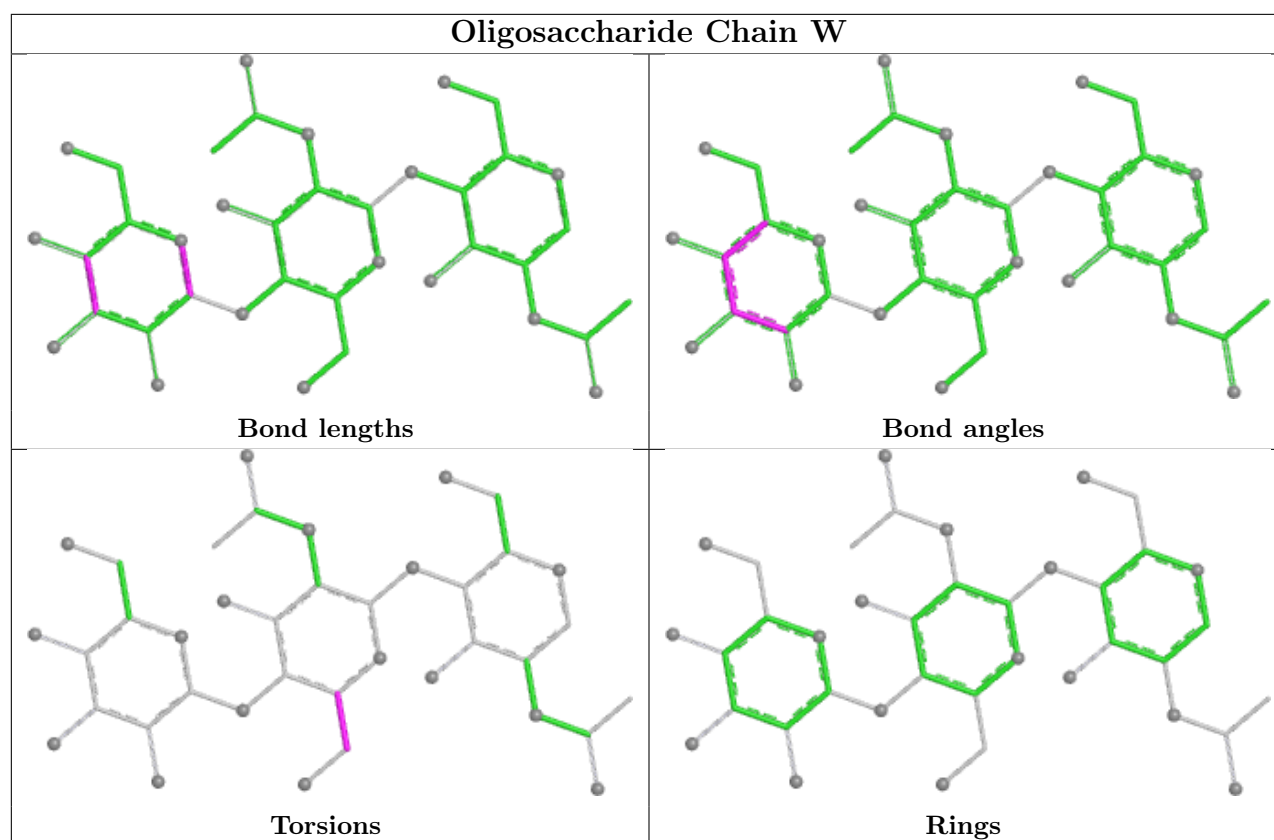












5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 for Mogul analysis.

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	$\# Z > 2$
6	SO4	H	301	-	4,4,4	0.24	0	6,6,6	0.07	0
7	NAG	E	504	-	14,14,15	0.21	0	17,19,21	0.42	0
6	SO4	H	302	-	4,4,4	0.23	0	6,6,6	0.07	0
6	SO4	D	508	-	4,4,4	0.23	0	6,6,6	0.08	0
7	NAG	E	505	1	14,14,15	0.21	0	17,19,21	0.40	0

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
7	NAG	E	504	-	-	2/6/23/26	0/1/1/1
7	NAG	E	505	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	E	504	NAG	O5-C5-C6-O6
7	E	504	NAG	C4-C5-C6-O6
7	E	505	NAG	O5-C5-C6-O6
7	E	505	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	H	302	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	383/456 (83%)	-0.19	3 (0%) 82 60	66, 108, 158, 186	0
1	B	383/456 (83%)	-0.22	2 (0%) 87 68	63, 102, 152, 189	0
1	C	384/456 (84%)	-0.18	2 (0%) 87 68	62, 103, 155, 184	0
1	D	384/456 (84%)	-0.23	3 (0%) 82 60	62, 103, 161, 189	0
1	E	383/456 (83%)	-0.24	3 (0%) 82 60	60, 103, 157, 195	0
2	F	122/124 (98%)	0.05	0 100 100	75, 120, 153, 178	1 (0%)
2	G	124/124 (100%)	-0.19	0 100 100	86, 108, 140, 156	1 (0%)
2	H	124/124 (100%)	-0.31	0 100 100	71, 95, 125, 147	1 (0%)
2	I	124/124 (100%)	-0.17	0 100 100	74, 119, 160, 178	1 (0%)
2	J	124/124 (100%)	-0.31	1 (0%) 82 60	73, 98, 131, 156	1 (0%)
All	All	2535/2900 (87%)	-0.21	14 (0%) 85 65	60, 104, 156, 195	5 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	312	ASP	5.7
1	A	311	GLN	3.1
1	B	311	GLN	2.6
1	E	452	LEU	2.6
1	E	275	ALA	2.4
1	A	452	LEU	2.4
1	D	456	TRP	2.3
1	D	452	LEU	2.2
1	C	242	PHE	2.2
2	J	100	CYS	2.1
1	C	275	ALA	2.1
1	D	311	GLN	2.1
1	E	247	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	456	TRP	2.1

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.

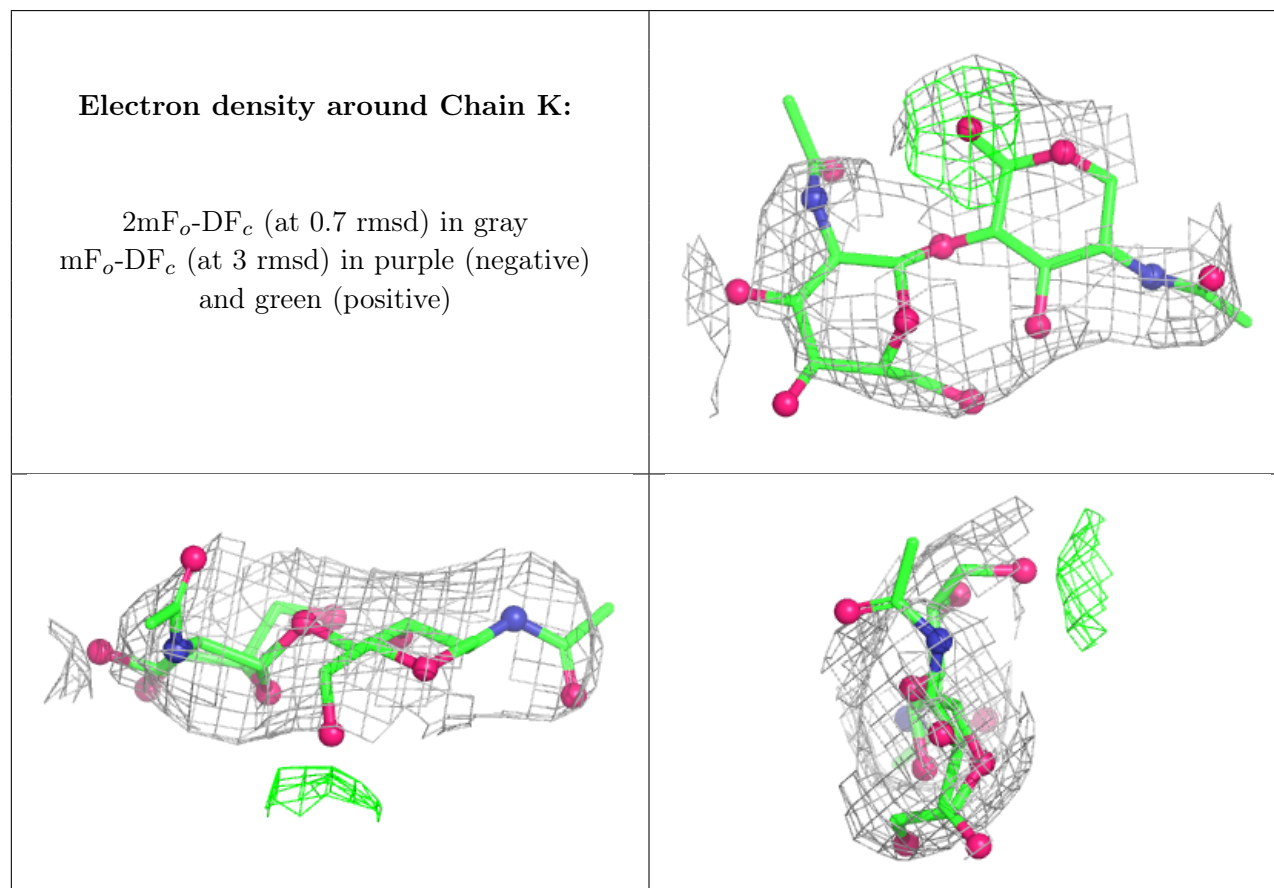
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	BMA	W	3	11/12	0.40	0.12	128,165,179,194	0
4	BMA	T	3	11/12	0.58	0.10	109,159,179,197	0
4	BMA	L	3	11/12	0.59	0.09	130,154,169,170	0
4	BMA	Q	3	11/12	0.59	0.10	122,145,162,174	0
4	BMA	N	3	11/12	0.62	0.09	107,129,155,168	0
3	NAG	U	2	14/15	0.65	0.10	126,137,167,187	0
3	NAG	K	1	14/15	0.65	0.11	109,140,165,167	0
3	NAG	R	2	14/15	0.70	0.11	140,171,185,192	0
3	NAG	K	2	14/15	0.71	0.11	131,161,174,178	0
3	NAG	M	2	14/15	0.72	0.09	97,143,159,161	0
3	NAG	S	2	14/15	0.76	0.10	111,139,151,157	0
3	NAG	O	2	14/15	0.79	0.10	116,132,167,172	0
3	NAG	O	1	14/15	0.80	0.11	112,123,142,144	0
3	NAG	V	2	14/15	0.81	0.09	107,130,143,150	0
3	NAG	P	2	14/15	0.82	0.11	124,151,174,180	0
3	NAG	U	1	14/15	0.84	0.10	105,127,139,152	0
4	NAG	Q	1	14/15	0.85	0.14	77,105,133,141	0
3	NAG	X	2	14/15	0.85	0.09	103,125,145,149	0
4	NAG	T	2	14/15	0.86	0.10	88,120,157,163	0
3	NAG	R	1	14/15	0.87	0.09	132,144,160,165	0
4	NAG	T	1	14/15	0.88	0.12	89,110,145,147	0
3	NAG	M	1	14/15	0.90	0.09	86,107,126,148	0
4	NAG	N	2	14/15	0.90	0.08	86,107,138,140	0
4	NAG	W	1	14/15	0.90	0.12	73,113,121,132	0
4	NAG	L	1	14/15	0.90	0.10	81,107,150,152	0
4	NAG	L	2	14/15	0.91	0.08	93,136,146,151	0
4	NAG	Q	2	14/15	0.91	0.08	86,106,135,149	0

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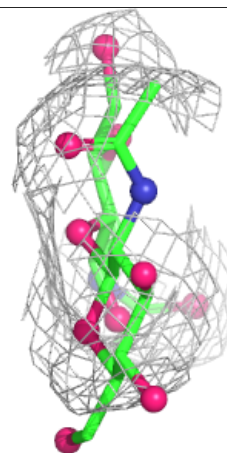
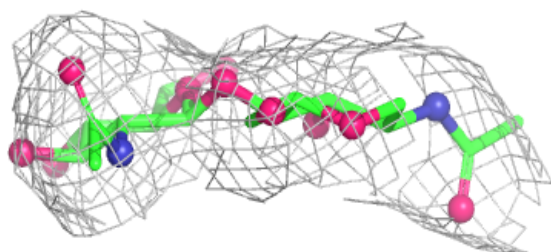
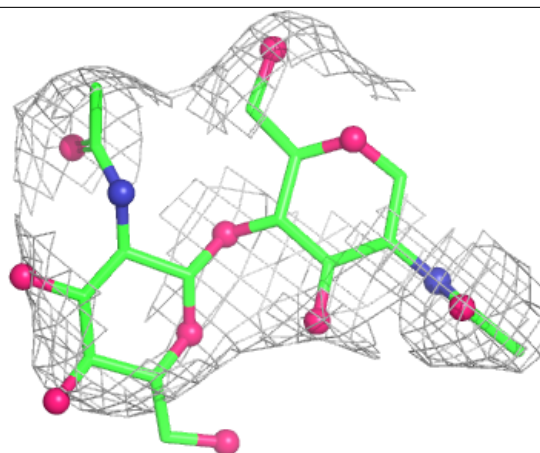
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	W	2	14/15	0.92	0.10	76,108,139,156	0
4	NAG	N	1	14/15	0.92	0.10	78,101,135,149	0
3	NAG	V	1	14/15	0.94	0.08	76,105,125,143	0
3	NAG	S	1	14/15	0.94	0.06	63,97,134,147	0
3	NAG	P	1	14/15	0.95	0.06	82,99,136,163	0
3	NAG	X	1	14/15	0.95	0.08	77,94,120,134	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.



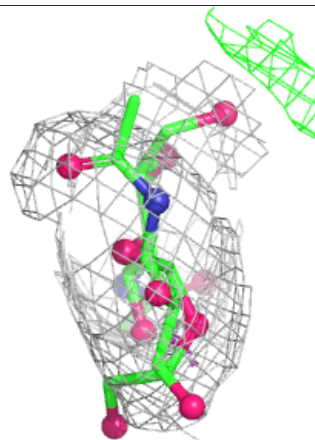
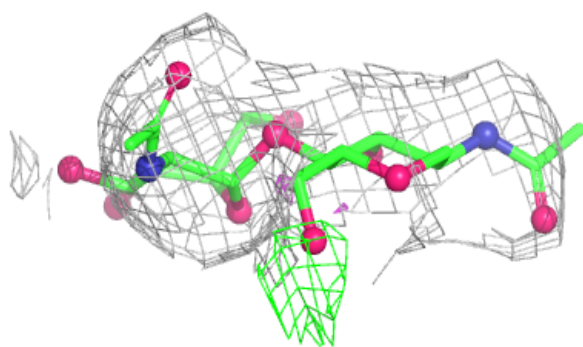
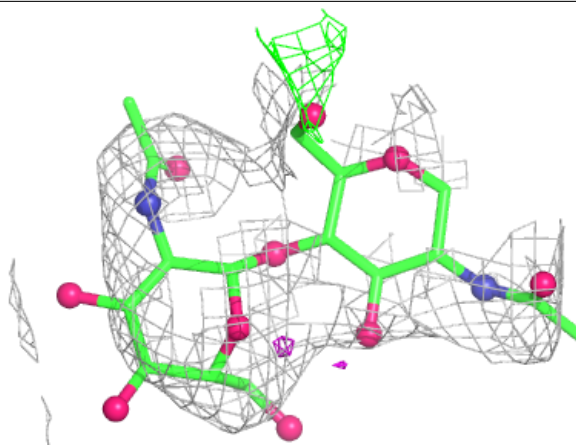
Electron density around Chain M:

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



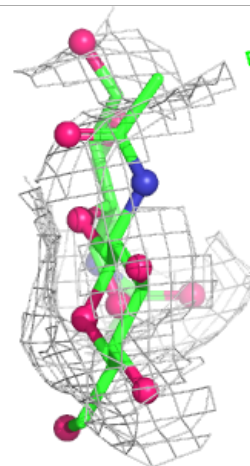
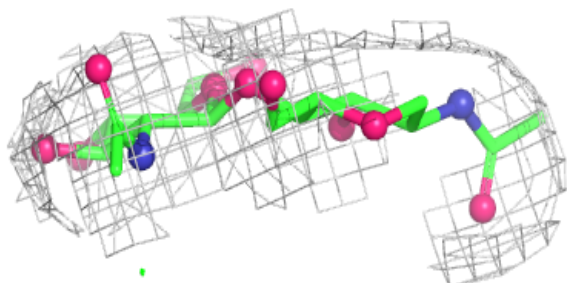
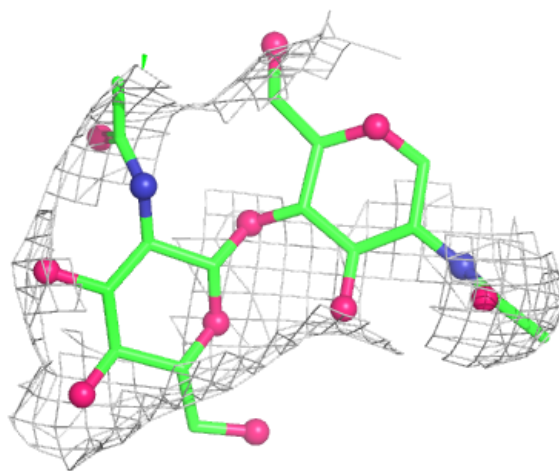
Electron density around Chain O:

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



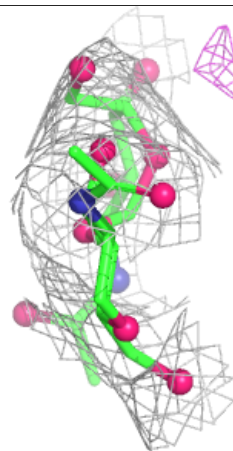
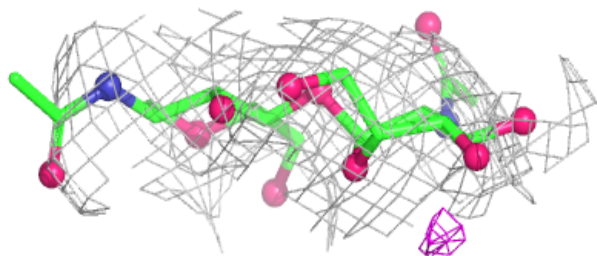
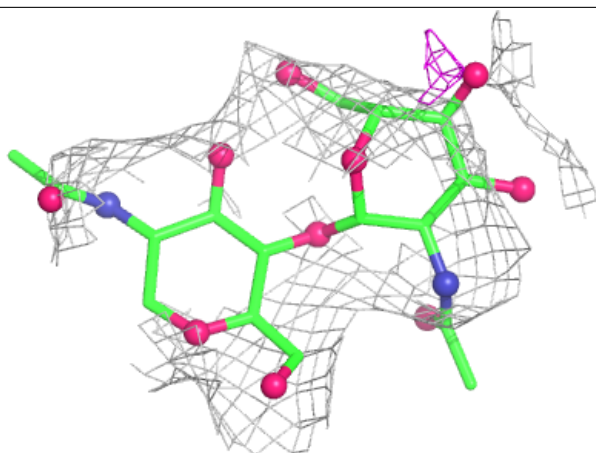
Electron density around Chain P:

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



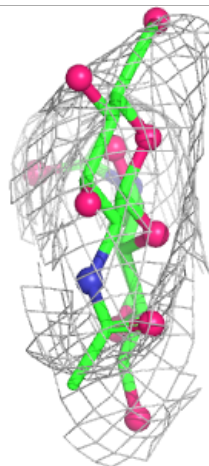
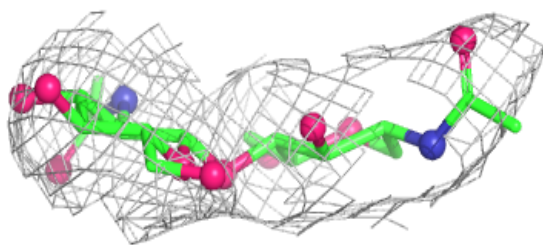
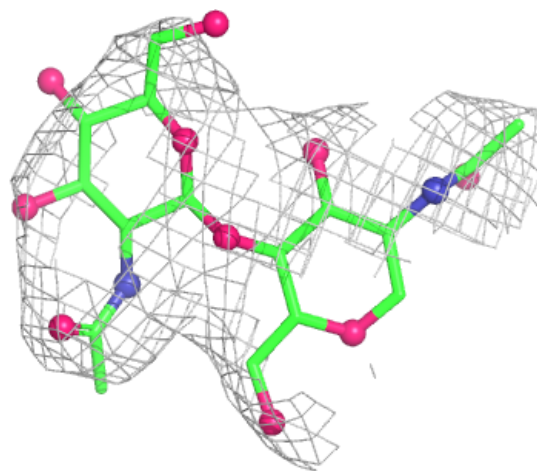
Electron density around Chain R:

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



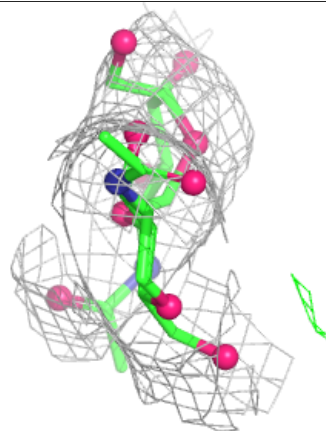
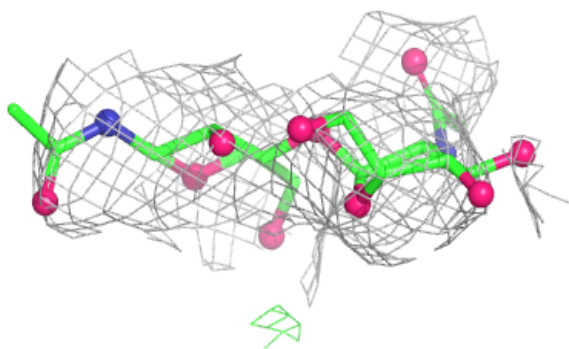
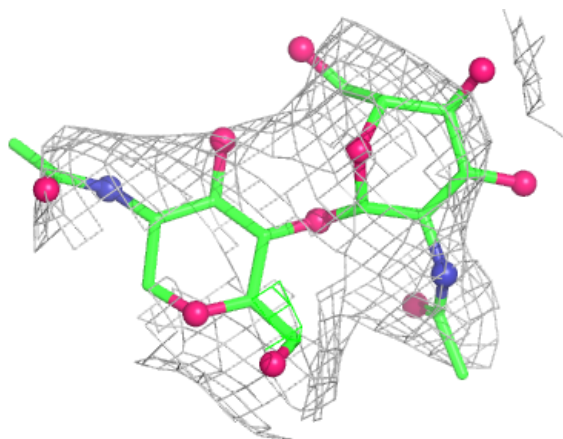
Electron density around Chain S:

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



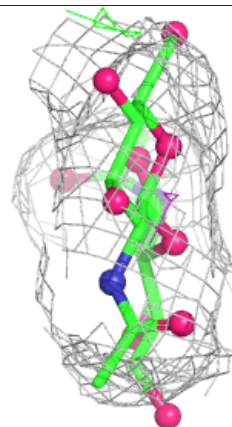
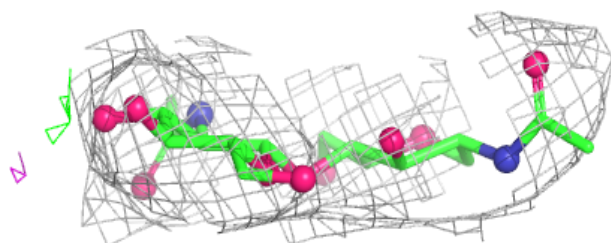
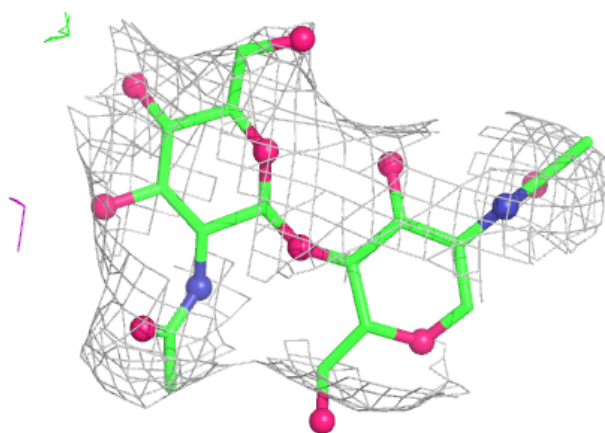
Electron density around Chain U:

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



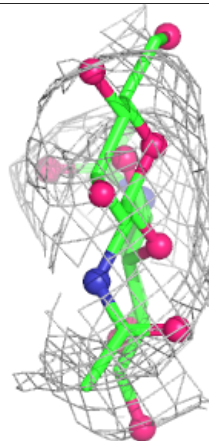
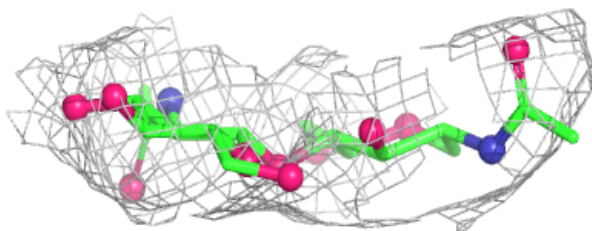
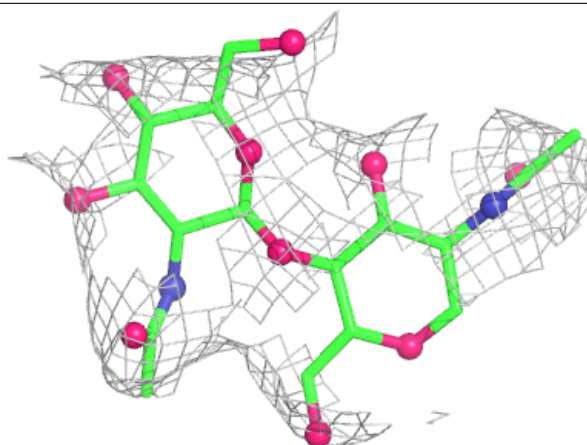
Electron density around Chain V:

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

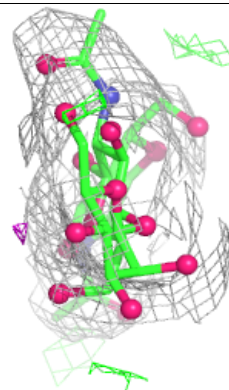
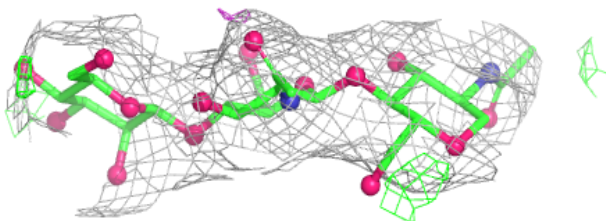
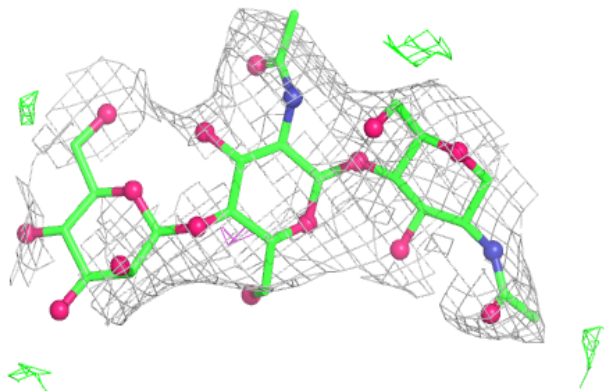


Electron density around Chain X:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

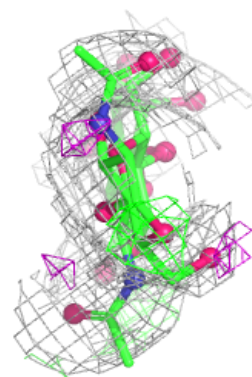
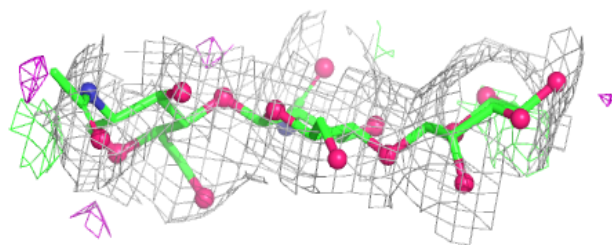
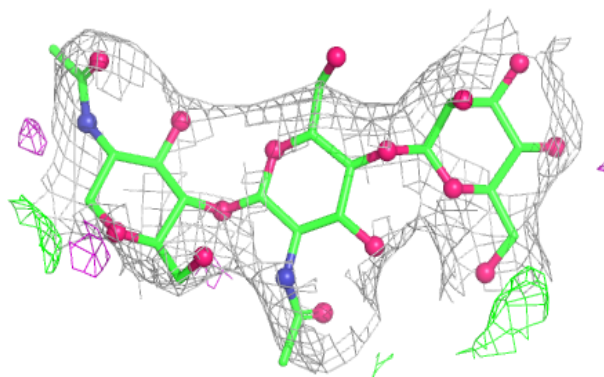
**Electron density around Chain L:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

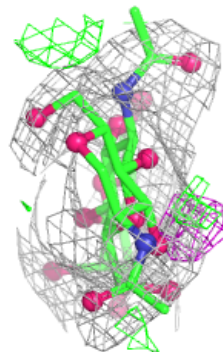
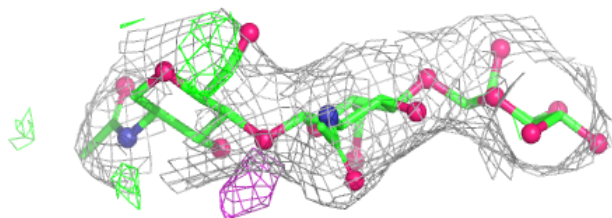
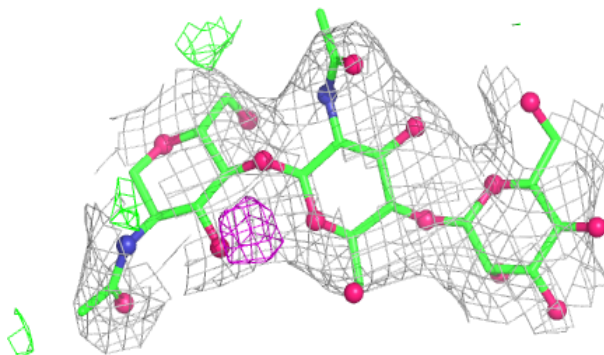


Electron density around Chain N:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

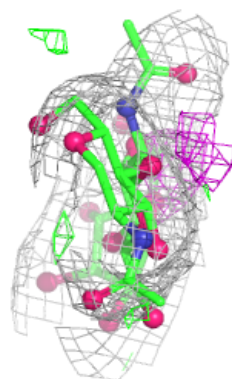
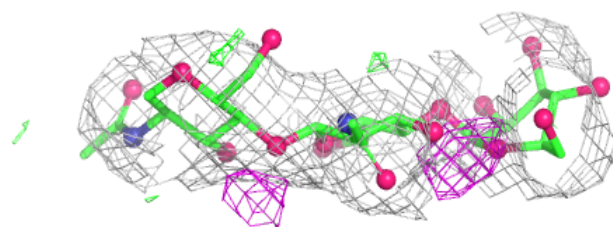
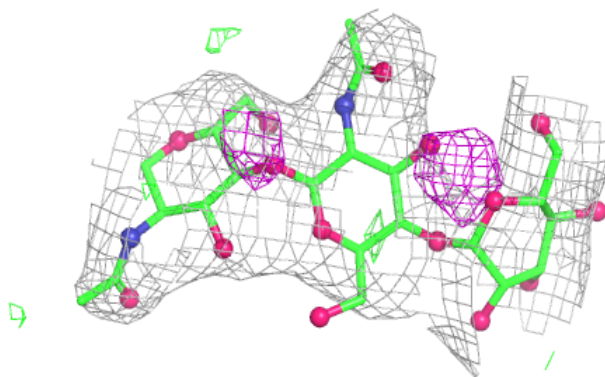
**Electron density around Chain Q:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

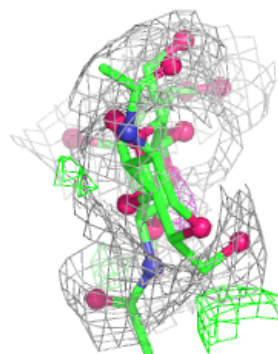
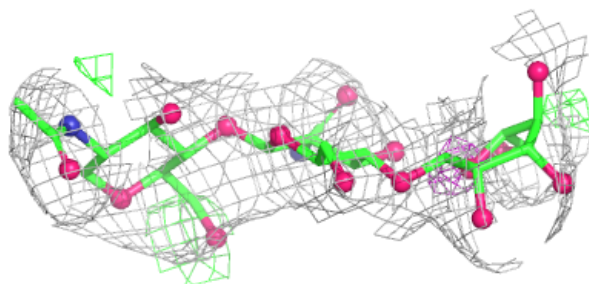
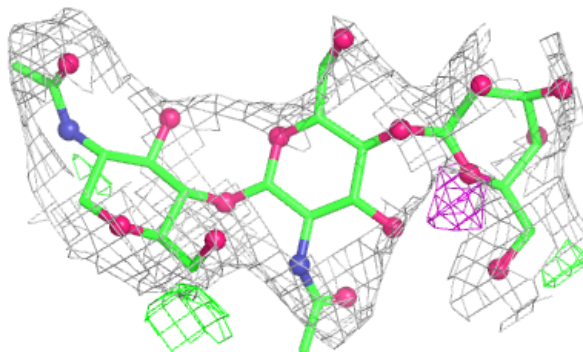


Electron density around Chain T:

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

**Electron density around Chain W:**

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



6.4 Ligands

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
7	NAG	E	504	14/15	0.62	0.14	136,174,190,191	0
6	SO4	H	301	5/5	0.79	0.12	154,161,191,193	0
6	SO4	D	508	5/5	0.79	0.08	169,186,200,203	0
6	SO4	H	302	5/5	0.88	0.08	125,141,156,163	0
5	CL	A	508	1/1	0.89	0.05	114,114,114,114	0
7	NAG	E	505	14/15	0.91	0.07	117,132,149,151	0
5	CL	E	508	1/1	0.96	0.06	87,87,87,87	0
5	CL	C	508	1/1	0.97	0.04	70,70,70,70	0
5	CL	D	509	1/1	0.97	0.03	88,88,88,88	0
5	CL	B	508	1/1	0.99	0.06	84,84,84,84	0

6.5 Other polymers

There are no such residues in this entry.