



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 31, 2023 – 03:15 AM EDT

PDB ID : 1TEI
Title : STRUCTURE OF CONCAVALIN A COMPLEXED TO BETA-D-GLCNAC (1,2)ALPHA-D-MAN-(1,6)[BETA-D-GLCNAC(1,2)ALPHA-D-MAN (1,6)]ALPHA-D-MAN
Authors : Naismith, J.H.; Moothoo, D.N.
Deposited on : 1997-05-28
Resolution : 2.70 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.34
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.34

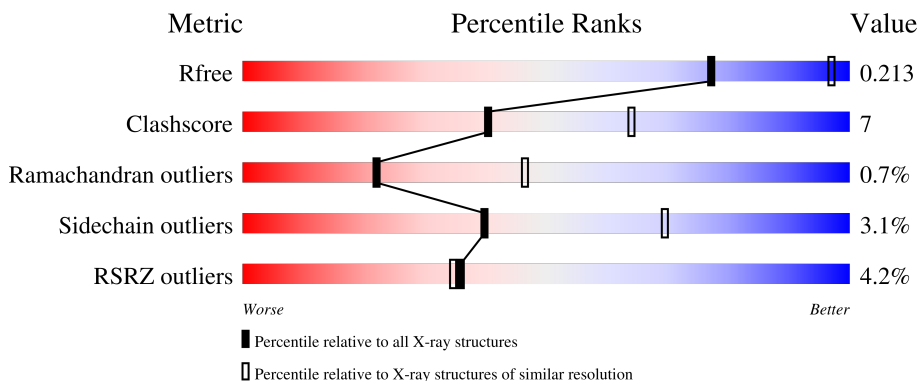
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	237	 3% 85% 14%
1	B	237	 3% 84% 14% .
1	C	237	 % 83% 14% .
1	D	237	 3% 83% 14% .
1	E	237	 5% 82% 16% .

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Mol	Chain	Length	Quality of chain
1	F	237	 9% 81% 17%
1	G	237	 5% 83% 15%
1	H	237	 5% 88% 10%
2	I	5	 20% 80%
2	J	5	 60% 40%
2	K	5	 60% 40%
2	L	5	 40% 60%
2	M	5	 40% 60%
2	N	5	 60% 20% 20%
2	O	5	 60% 40%
2	P	5	 60% 40%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15060 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 CONCANAVALIN A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	237	1809	1141	302	364	2	37	0	0
1	B	237	1809	1141	302	364	2	45	0	0
1	C	237	1809	1141	302	364	2	30	0	0
1	D	237	1809	1141	302	364	2	62	0	0
1	E	237	1809	1141	302	364	2	28	0	0
1	F	237	1809	1141	302	364	2	78	0	0
1	G	237	1809	1141	302	364	2	26	0	0
1	H	237	1809	1141	302	364	2	64	0	0

There are 16 discrepancies between the modelled and reference sequences:

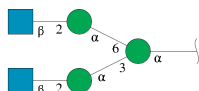
Chain	Residue	Modelled	Actual	Comment	Reference
A	151	ASP	GLU	conflict	UNP P02866
A	155	GLU	ARG	conflict	UNP P02866
B	151	ASP	GLU	conflict	UNP P02866
B	155	GLU	ARG	conflict	UNP P02866
C	151	ASP	GLU	conflict	UNP P02866
C	155	GLU	ARG	conflict	UNP P02866
D	151	ASP	GLU	conflict	UNP P02866
D	155	GLU	ARG	conflict	UNP P02866
E	151	ASP	GLU	conflict	UNP P02866
E	155	GLU	ARG	conflict	UNP P02866
F	151	ASP	GLU	conflict	UNP P02866
F	155	GLU	ARG	conflict	UNP P02866
G	151	ASP	GLU	conflict	UNP P02866

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Chain	Residue	Modelled	Actual	Comment	Reference
G	155	GLU	ARG	conflict	UNP P02866
H	151	ASP	GLU	conflict	UNP P02866
H	155	GLU	ARG	conflict	UNP P02866

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	I	5	62	34	2	26	0	0	0
2	J	5	62	34	2	26	0	0	0
2	K	5	62	34	2	26	0	0	0
2	L	5	62	34	2	26	0	0	0
2	M	5	62	34	2	26	0	0	0
2	N	5	62	34	2	26	11	0	0
2	O	5	62	34	2	26	0	0	0
2	P	5	62	34	2	26	10	0	0

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		
3	B	1	Total	Mn	0	0
			1	1		
3	C	1	Total	Mn	0	0
			1	1		
3	D	1	Total	Mn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	1	Total 1	Mn 1	0	0
3	F	1	Total 1	Mn 1	0	0
3	G	1	Total 1	Mn 1	0	0
3	H	1	Total 1	Mn 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Ca 1	0	0
4	B	1	Total 1	Ca 1	0	0
4	C	1	Total 1	Ca 1	0	0
4	D	1	Total 1	Ca 1	0	0
4	E	1	Total 1	Ca 1	0	0
4	F	1	Total 1	Ca 1	0	0
4	G	1	Total 1	Ca 1	0	0
4	H	1	Total 1	Ca 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	15	Total 15	O 15	0	0
5	B	10	Total 10	O 10	0	0
5	C	13	Total 13	O 13	0	0
5	D	7	Total 7	O 7	0	0
5	E	19	Total 19	O 19	0	0

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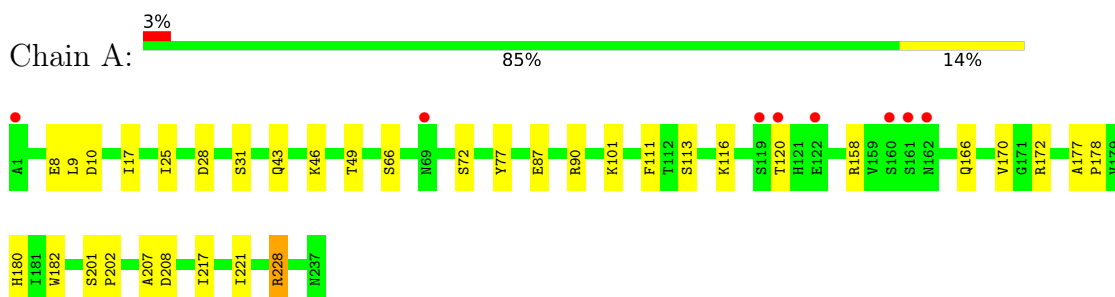
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	F	5	Total O 5 5	0	0
5	G	5	Total O 5 5	0	0
5	H	2	Total O 2 2	0	0

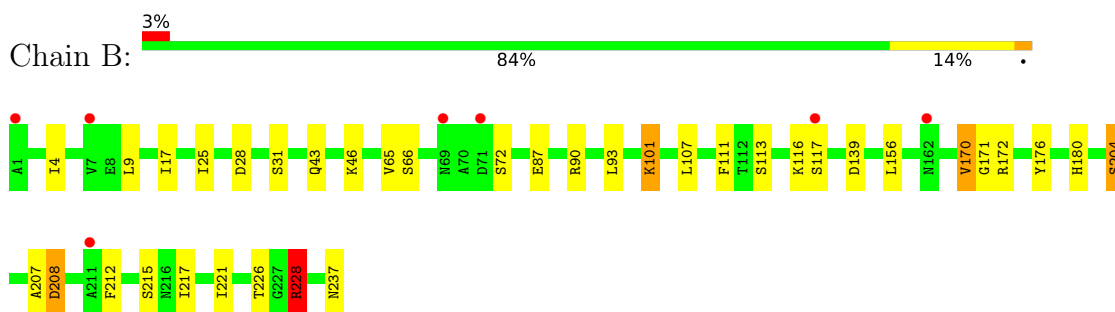
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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.

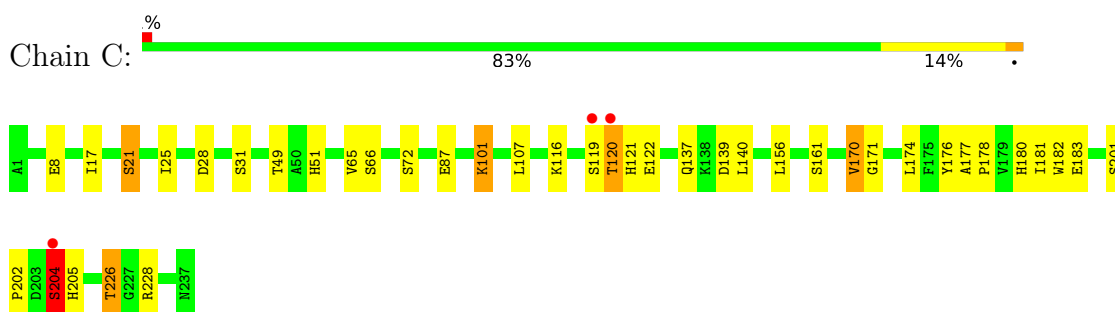
- Molecule 1: CONCANAVALIN A



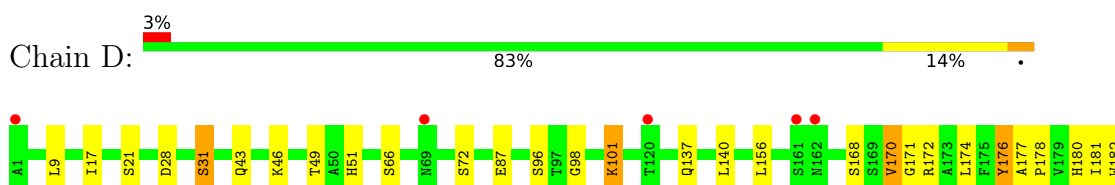
- Molecule 1: CONCANAVALIN A



- Molecule 1: CONCANAVALIN A

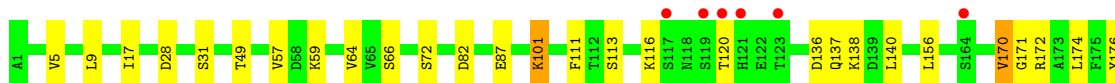
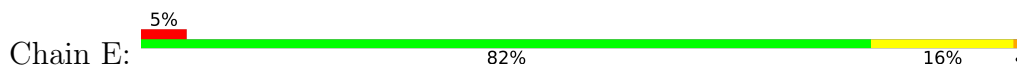


- Molecule 1: CONCANAVALIN A

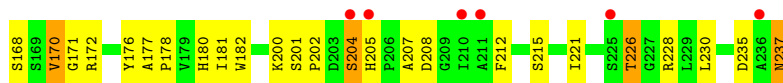
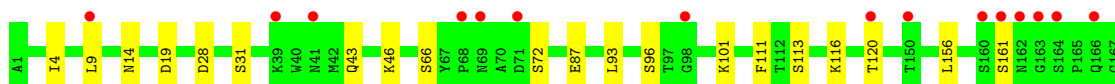
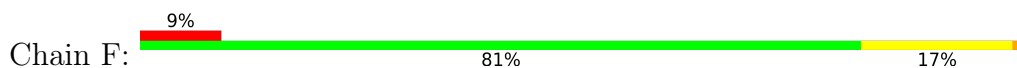




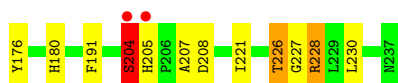
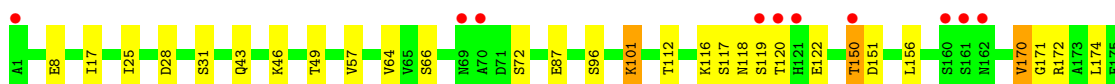
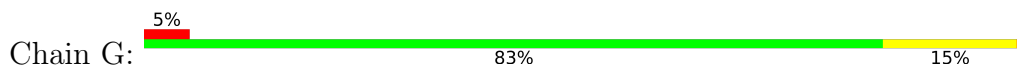
- Molecule 1: CONCANAVALIN A



- Molecule 1: CONCANAVALIN A



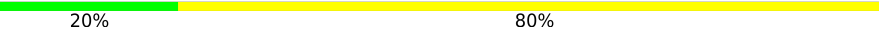
- Molecule 1: CONCANAVALIN A



- Molecule 1: CONCANAVALIN A



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

Chain I:  20% 80%



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

Chain J:  60% 40%



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

Chain K:  60% 40%



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

Chain L:  40% 60%



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

Chain M:  40% 60%



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

Chain N:  60% 20% 20%



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

anose

Chain O:  60% 40%

MAN1
MAN2
MAN3
MAN4
MAN5

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

Chain P:  60% 40%

MAN1
MAN2
MAN3
MAN4
MAN5

4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	176.50Å 122.80Å 124.60Å 90.00° 134.20° 90.00°	Depositor
Resolution (Å)	25.00 – 2.70 24.51 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.1 (25.00-2.70) 97.1 (24.51-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.09 (at 2.71Å)	Xtrriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.179 , 0.219 0.174 , 0.213	Depositor DCC
R_{free} test set	5119 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtrriage
Anisotropy	0.246	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.000 for h+2*k,-h-l 0.000 for k+1,h+1,-l 0.000 for -k+1,-h-l,-l 0.009 for -h+k-l,-l,-k 0.000 for -h-k-l,l,k 0.000 for h-k+1,l,-h-l 0.000 for -k-l,-h-l,k 0.000 for h+k+1,-l,-h-l 0.000 for k-l,h+1,-k 0.007 for h,-k,-h-l 0.021 for -h-2*k,-k,l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15060	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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, CA, MAN, MN

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	0.55	0/1851	0.83	1/2522 (0.0%)
1	B	0.50	0/1851	0.78	1/2522 (0.0%)
1	C	0.53	0/1851	0.81	0/2522
1	D	0.57	1/1851 (0.1%)	0.81	1/2522 (0.0%)
1	E	0.54	0/1851	0.82	2/2522 (0.1%)
1	F	0.50	0/1851	0.79	1/2522 (0.0%)
1	G	0.51	0/1851	0.80	0/2522
1	H	0.54	1/1851 (0.1%)	0.79	0/2522
All	All	0.53	2/14808 (0.0%)	0.80	6/20176 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	176	TYR	CE2-CZ	-5.71	1.31	1.38
1	H	176	TYR	CE1-CZ	-5.28	1.31	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	9	LEU	N-CA-C	-5.17	97.03	111.00
1	F	9	LEU	N-CA-C	-5.17	97.05	111.00
1	A	9	LEU	N-CA-C	-5.15	97.11	111.00
1	B	9	LEU	N-CA-C	-5.08	97.28	111.00
1	D	9	LEU	N-CA-C	-5.06	97.34	111.00
1	E	5	VAL	N-CA-C	-5.04	97.39	111.00

There are no chirality outliers.

There are no planarity outliers.

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	1809	0	1755	26	1
1	B	1809	0	1755	23	0
1	C	1809	0	1755	30	0
1	D	1809	0	1755	28	0
1	E	1809	0	1755	32	0
1	F	1809	0	1755	28	0
1	G	1809	0	1755	37	0
1	H	1809	0	1755	24	0
2	I	62	0	54	1	0
2	J	62	0	54	1	0
2	K	62	0	54	1	0
2	L	62	0	54	2	0
2	M	62	0	54	1	0
2	N	62	0	54	1	1
2	O	62	0	54	1	0
2	P	62	0	54	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
5	A	15	0	0	2	0
5	B	10	0	0	2	0
5	C	13	0	0	1	0
5	D	7	0	0	0	0
5	E	19	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	F	5	0	0	1	0
5	G	5	0	0	0	0
5	H	2	0	0	0	0
All	All	15060	0	14472	206	1

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 7.

All (206) 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:82:ASP:HB2	5:E:258:HOH:O	1.77	0.83
1:F:204:SER:HB2	1:F:205:HIS:CD2	2.16	0.79
1:F:207:ALA:HB1	1:F:208:ASP:OD2	1.83	0.79
1:C:17:ILE:HD13	1:C:228:ARG:HD3	1.65	0.79
1:F:170:VAL:HG22	1:F:226:THR:HG23	1.64	0.79
1:A:17:ILE:HD13	1:A:228:ARG:HD3	1.63	0.79
1:D:17:ILE:HD13	1:D:228:ARG:HD3	1.62	0.79
1:H:172:ARG:HD2	1:H:221:ILE:HG12	1.65	0.78
1:E:228:ARG:HG2	2:M:4:MAN:O3	1.84	0.78
1:E:49:THR:HG21	1:G:116:LYS:NZ	2.00	0.77
1:B:172:ARG:HD2	1:B:221:ILE:HG12	1.68	0.76
5:A:252:HOH:O	1:C:51:HIS:HB2	1.86	0.74
1:A:116:LYS:NZ	1:C:49:THR:HG21	2.03	0.73
1:G:17:ILE:HD13	1:G:228:ARG:HD3	1.71	0.73
1:A:116:LYS:HZ3	1:C:49:THR:HG21	1.53	0.71
5:F:247:HOH:O	1:H:51:HIS:HB2	1.90	0.71
1:F:204:SER:HB2	1:F:205:HIS:HD2	1.56	0.70
1:B:17:ILE:HD13	1:B:228:ARG:HD3	1.73	0.70
1:H:17:ILE:HD13	1:H:228:ARG:HD3	1.72	0.70
1:E:57:VAL:HG22	1:G:64:VAL:HG21	1.74	0.69
1:G:66:SER:HB3	1:G:72:SER:HB3	1.75	0.68
1:H:170:VAL:HG22	1:H:226:THR:HG22	1.75	0.68
1:G:228:ARG:HG2	2:O:4:MAN:O3	1.94	0.68
1:F:116:LYS:HZ1	1:H:49:THR:HG21	1.60	0.66
5:B:249:HOH:O	1:D:51:HIS:HB2	1.95	0.66
1:E:17:ILE:HD13	1:E:228:ARG:HD3	1.76	0.65
1:B:66:SER:HB3	1:B:72:SER:HB3	1.77	0.65
1:B:116:LYS:HZ1	1:D:49:THR:HG21	1.60	0.65
1:B:170:VAL:HG22	1:B:226:THR:HG22	1.77	0.65
1:C:66:SER:HB3	1:C:72:SER:HB3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:ARG:HG2	2:J:4:MAN:O3	1.97	0.64
1:F:205:HIS:CD2	1:F:205:HIS:N	2.66	0.64
1:C:228:ARG:HG2	2:K:4:MAN:O3	1.98	0.64
1:E:49:THR:HG21	1:G:116:LYS:HZ1	1.64	0.62
1:A:66:SER:HB3	1:A:72:SER:HB3	1.81	0.62
1:B:156:LEU:O	1:B:171:GLY:HA3	2.00	0.62
1:E:220:SER:HB2	5:E:255:HOH:O	2.00	0.60
1:H:172:ARG:HD2	1:H:221:ILE:CG1	2.30	0.60
1:B:172:ARG:HD2	1:B:221:ILE:CG1	2.31	0.60
1:D:228:ARG:HG2	2:L:4:MAN:O3	2.01	0.59
1:H:66:SER:HB3	1:H:72:SER:HB3	1.83	0.59
1:D:156:LEU:O	1:D:171:GLY:HA3	2.02	0.59
1:F:116:LYS:NZ	1:H:49:THR:HG21	2.17	0.58
1:C:178:PRO:HD3	1:D:176:TYR:HD1	1.69	0.58
1:D:172:ARG:HD2	1:D:221:ILE:HG12	1.85	0.58
1:E:49:THR:HG21	1:G:116:LYS:HZ3	1.68	0.57
1:E:59:LYS:HE3	5:E:256:HOH:O	2.03	0.56
1:E:204:SER:HB2	1:E:205:HIS:CD2	2.40	0.56
1:F:172:ARG:HD2	1:F:221:ILE:HG12	1.88	0.56
1:G:43:GLN:HE21	1:G:46:LYS:HD2	1.71	0.56
1:E:66:SER:HB3	1:E:72:SER:HB3	1.87	0.56
1:G:204:SER:HB2	1:G:205:HIS:CD2	2.41	0.55
1:F:156:LEU:O	1:F:171:GLY:HA3	2.05	0.55
1:G:156:LEU:O	1:G:171:GLY:HA3	2.07	0.55
1:F:28:ASP:HB3	1:F:31:SER:O	2.07	0.54
1:B:116:LYS:NZ	1:D:49:THR:HG21	2.23	0.54
1:A:43:GLN:HE21	1:A:46:LYS:HD2	1.73	0.54
1:D:170:VAL:HG22	1:D:226:THR:CG2	2.37	0.54
1:C:119:SER:HB2	1:C:122:GLU:HB2	1.90	0.54
1:F:43:GLN:HE21	1:F:46:LYS:HD2	1.72	0.54
1:E:116:LYS:NZ	1:G:49:THR:HG21	2.23	0.54
1:A:116:LYS:NZ	5:A:252:HOH:O	2.39	0.53
1:B:66:SER:HB3	1:B:72:SER:CB	2.38	0.53
1:A:49:THR:HG21	1:C:116:LYS:NZ	2.24	0.53
1:C:170:VAL:HG22	1:C:226:THR:CG2	2.39	0.53
1:H:101:LYS:N	1:H:101:LYS:HD3	2.24	0.52
1:E:156:LEU:O	1:E:171:GLY:HA3	2.09	0.52
1:A:87:GLU:HG3	1:A:182:TRP:O	2.09	0.52
1:G:176:TYR:CE2	1:H:176:TYR:HE1	2.28	0.52
1:E:101:LYS:N	1:E:101:LYS:HD3	2.24	0.52
1:H:228:ARG:HG2	2:P:4:MAN:O3	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:SER:HB2	1:C:205:HIS:CD2	2.45	0.51
1:D:66:SER:HB3	1:D:72:SER:HB3	1.92	0.51
1:A:228:ARG:HG2	2:I:4:MAN:O3	2.10	0.51
1:H:181:ILE:HG23	1:H:182:TRP:HD1	1.75	0.51
1:A:66:SER:HB3	1:A:72:SER:CB	2.40	0.51
1:D:170:VAL:HG22	1:D:226:THR:HG22	1.91	0.51
1:E:235:ASP:OD1	1:E:237:ASN:HB2	2.11	0.50
1:B:111:PHE:CE2	1:B:113:SER:HB2	2.47	0.50
1:F:66:SER:HB3	1:F:72:SER:HB3	1.93	0.50
1:B:43:GLN:HE21	1:B:46:LYS:HD2	1.76	0.50
1:G:66:SER:HB3	1:G:72:SER:CB	2.41	0.50
1:G:101:LYS:N	1:G:101:LYS:HD3	2.26	0.50
1:G:172:ARG:HD2	1:G:221:ILE:CG1	2.41	0.50
1:C:181:ILE:HG23	1:C:182:TRP:HD1	1.77	0.50
1:D:43:GLN:HE21	1:D:46:LYS:HD2	1.77	0.50
1:C:87:GLU:HG2	1:C:180:HIS:NE2	2.27	0.49
1:H:170:VAL:HG22	1:H:226:THR:CG2	2.41	0.49
1:E:172:ARG:HD2	1:E:221:ILE:HG12	1.94	0.49
1:D:101:LYS:N	1:D:101:LYS:HD3	2.27	0.49
1:C:28:ASP:HB3	1:C:31:SER:O	2.11	0.49
1:C:178:PRO:HD3	1:D:176:TYR:CD1	2.47	0.49
1:A:158:ARG:HG3	1:A:166:GLN:HG3	1.94	0.49
1:H:87:GLU:HG2	1:H:180:HIS:NE2	2.28	0.48
1:D:181:ILE:HG23	1:D:182:TRP:HD1	1.78	0.48
1:F:172:ARG:HD2	1:F:221:ILE:CG1	2.44	0.48
1:G:208:ASP:OD1	1:G:227:GLY:HA2	2.13	0.48
1:A:172:ARG:HD2	1:A:221:ILE:HG12	1.94	0.47
1:H:156:LEU:O	1:H:171:GLY:HA3	2.13	0.47
1:G:87:GLU:HG2	1:G:180:HIS:NE2	2.29	0.47
1:H:204:SER:HB2	1:H:205:HIS:CD2	2.49	0.47
1:B:25:ILE:HG21	1:B:65:VAL:HG21	1.97	0.47
1:A:17:ILE:CD1	1:A:228:ARG:HD3	2.41	0.47
1:E:176:TYR:CE2	1:F:176:TYR:CE2	3.02	0.47
1:E:64:VAL:HG21	1:G:57:VAL:HG22	1.97	0.47
1:B:17:ILE:CD1	1:B:228:ARG:HD3	2.42	0.47
1:G:43:GLN:NE2	1:G:46:LYS:HD2	2.30	0.47
1:E:170:VAL:HG22	1:E:226:THR:HG22	1.97	0.47
1:H:170:VAL:HG13	1:H:226:THR:HA	1.95	0.47
1:C:25:ILE:HG21	1:C:65:VAL:HG21	1.96	0.46
1:A:116:LYS:HZ3	1:C:49:THR:CG2	2.25	0.46
1:H:43:GLN:HE21	1:H:46:LYS:HD3	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:137:GLN:HG2	1:C:140:LEU:HD12	1.96	0.46
1:D:137:GLN:HG2	1:D:140:LEU:HD12	1.96	0.46
1:H:8:GLU:O	1:H:25:ILE:HA	2.16	0.46
1:D:96:SER:OG	1:D:230:LEU:HA	2.16	0.46
1:D:181:ILE:HG23	1:D:182:TRP:CD1	2.51	0.46
1:G:150:THR:O	1:G:151:ASP:HB2	2.15	0.46
1:C:87:GLU:HG2	1:C:180:HIS:CD2	2.50	0.46
1:E:174:LEU:HD12	1:E:174:LEU:N	2.31	0.46
1:E:87:GLU:HG3	1:E:182:TRP:O	2.14	0.46
1:E:111:PHE:CE2	1:E:113:SER:HB2	2.51	0.46
1:B:139:ASP:O	1:B:176:TYR:HB2	2.16	0.46
1:E:28:ASP:HB3	1:E:31:SER:O	2.16	0.46
1:C:170:VAL:HG22	1:C:226:THR:HG22	1.98	0.45
1:G:8:GLU:O	1:G:25:ILE:HA	2.16	0.45
1:C:177:ALA:HA	1:C:178:PRO:HD3	1.87	0.45
1:F:96:SER:OG	1:F:230:LEU:HA	2.17	0.45
1:B:107:LEU:HD12	1:B:107:LEU:N	2.32	0.45
1:E:87:GLU:HG2	1:E:180:HIS:CD2	2.52	0.45
1:D:98:GLY:HA2	2:L:5:NAG:O7	2.17	0.45
1:F:205:HIS:HD2	1:F:205:HIS:N	2.13	0.45
1:G:172:ARG:HD2	1:G:221:ILE:HG12	1.98	0.45
1:C:8:GLU:O	1:C:25:ILE:HA	2.17	0.45
1:G:176:TYR:CE2	1:H:176:TYR:CE1	3.05	0.45
1:C:156:LEU:O	1:C:171:GLY:HA3	2.18	0.44
1:E:116:LYS:HZ1	1:G:49:THR:HG21	1.80	0.44
1:G:96:SER:OG	1:G:230:LEU:HA	2.17	0.44
1:A:177:ALA:HA	1:A:178:PRO:HD3	1.83	0.44
1:H:87:GLU:HG2	1:H:180:HIS:CD2	2.52	0.44
1:A:87:GLU:HG2	1:A:180:HIS:CD2	2.52	0.44
1:A:172:ARG:HD2	1:A:221:ILE:CG1	2.47	0.44
1:E:87:GLU:HG2	1:E:180:HIS:NE2	2.32	0.44
1:A:90:ARG:NH1	1:A:217:ILE:HG23	2.32	0.44
1:D:17:ILE:CD1	1:D:228:ARG:HD3	2.41	0.44
1:F:87:GLU:HG2	1:F:180:HIS:CD2	2.53	0.44
1:D:28:ASP:HB3	1:D:31:SER:O	2.17	0.44
1:D:201:SER:HA	1:D:202:PRO:HD3	1.86	0.44
1:G:170:VAL:HG22	1:G:226:THR:CG2	2.48	0.44
1:E:207:ALA:HA	1:E:208:ASP:HA	1.81	0.43
1:E:64:VAL:HG21	1:G:57:VAL:CG2	2.48	0.43
1:B:93:LEU:HD23	1:B:212:PHE:HA	2.00	0.43
1:E:136:ASP:OD2	1:E:138:LYS:HE3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:43:GLN:NE2	1:F:46:LYS:HD2	2.33	0.43
1:D:208:ASP:OD1	1:D:227:GLY:HA2	2.18	0.43
1:F:111:PHE:CE2	1:F:113:SER:HB2	2.54	0.43
1:B:87:GLU:HG2	1:B:180:HIS:NE2	2.34	0.43
1:C:183:GLU:HG2	5:C:256:HOH:O	2.17	0.43
1:E:210:ILE:HG22	1:E:211:ALA:N	2.34	0.43
1:C:174:LEU:HD12	1:C:174:LEU:N	2.34	0.43
1:F:200:LYS:HB2	1:F:200:LYS:HE3	1.88	0.42
1:A:43:GLN:NE2	1:A:46:LYS:HD2	2.34	0.42
1:A:87:GLU:HG2	1:A:180:HIS:NE2	2.34	0.42
1:C:120:THR:O	1:C:122:GLU:N	2.52	0.42
1:E:137:GLN:HG2	1:E:140:LEU:HD12	2.02	0.42
1:F:201:SER:HA	1:F:202:PRO:HD3	1.84	0.42
1:F:235:ASP:OD1	1:F:237:ASN:HB3	2.18	0.42
1:G:174:LEU:HD12	1:G:174:LEU:N	2.34	0.42
1:B:90:ARG:NH1	1:B:217:ILE:HG23	2.34	0.42
1:A:28:ASP:HB3	1:A:31:SER:O	2.18	0.42
1:B:207:ALA:HA	1:B:208:ASP:HA	1.82	0.42
1:D:174:LEU:N	1:D:174:LEU:HD12	2.34	0.42
1:F:177:ALA:HA	1:F:178:PRO:HD3	1.86	0.42
1:G:119:SER:O	1:G:122:GLU:HB2	2.20	0.42
1:B:101:LYS:N	1:B:101:LYS:HD3	2.35	0.42
1:G:87:GLU:HG2	1:G:180:HIS:CD2	2.55	0.42
1:B:28:ASP:HB3	1:B:31:SER:O	2.20	0.42
1:D:87:GLU:HG2	1:D:180:HIS:CD2	2.54	0.42
1:F:14:ASN:C	1:F:19:ASP:HB2	2.40	0.42
1:F:168:SER:HG	2:N:5:NAG:C7	2.33	0.42
1:H:181:ILE:HG23	1:H:182:TRP:CD1	2.52	0.42
1:A:111:PHE:CE2	1:A:113:SER:HB2	2.54	0.42
1:C:107:LEU:HD12	1:C:107:LEU:N	2.35	0.42
1:F:4:ILE:HD13	1:F:215:SER:HB3	2.02	0.42
1:E:57:VAL:HG13	1:G:64:VAL:HG22	2.02	0.42
1:G:117:SER:HB3	1:G:118:ASN:H	1.68	0.42
1:A:201:SER:HA	1:A:202:PRO:HD3	1.83	0.41
1:C:139:ASP:O	1:C:176:TYR:HB2	2.20	0.41
1:B:4:ILE:HD13	1:B:215:SER:HB3	2.01	0.41
1:G:43:GLN:NE2	1:G:46:LYS:NZ	2.68	0.41
5:B:249:HOH:O	1:D:51:HIS:CG	2.73	0.41
1:G:112:THR:O	1:G:191:PHE:HA	2.21	0.41
1:G:170:VAL:HG22	1:G:226:THR:HG23	2.01	0.41
1:C:101:LYS:N	1:C:101:LYS:HD3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:SER:HA	1:C:202:PRO:HD3	1.87	0.41
1:D:237:ASN:HD22	1:D:237:ASN:HA	1.64	0.41
1:G:28:ASP:HB3	1:G:31:SER:O	2.20	0.41
1:H:87:GLU:HG3	1:H:182:TRP:O	2.20	0.41
1:F:181:ILE:HG23	1:F:182:TRP:HD1	1.85	0.41
1:A:207:ALA:HA	1:A:208:ASP:HA	1.83	0.41
1:A:8:GLU:O	1:A:25:ILE:HA	2.20	0.40
1:A:8:GLU:OE2	1:A:10:ASP:HB2	2.21	0.40
1:H:177:ALA:HA	1:H:178:PRO:HD3	1.91	0.40
1:D:177:ALA:HA	1:D:178:PRO:HD3	1.92	0.40
1:F:93:LEU:HD23	1:F:212:PHE:HA	2.03	0.40
1:G:207:ALA:HA	1:G:208:ASP:HA	1.80	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:TYR:OH	2:N:3:NAG:O7[3_445]	2.11	0.09

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	235/237 (99%)	220 (94%)	14 (6%)	1 (0%)	34 60
1	B	235/237 (99%)	219 (93%)	14 (6%)	2 (1%)	17 40
1	C	235/237 (99%)	221 (94%)	11 (5%)	3 (1%)	12 30
1	D	235/237 (99%)	220 (94%)	14 (6%)	1 (0%)	34 60
1	E	235/237 (99%)	215 (92%)	19 (8%)	1 (0%)	34 60
1	F	235/237 (99%)	218 (93%)	15 (6%)	2 (1%)	17 40
1	G	235/237 (99%)	216 (92%)	17 (7%)	2 (1%)	17 40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	235/237 (99%)	216 (92%)	17 (7%)	2 (1%)	17	40
All	All	1880/1896 (99%)	1745 (93%)	121 (6%)	14 (1%)	22	46

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	204	SER
1	H	150	THR
1	B	204	SER
1	B	228	ARG
1	C	121	HIS
1	D	204	SER
1	E	228	ARG
1	G	150	THR
1	G	204	SER
1	H	228	ARG
1	C	21	SER
1	F	228	ARG
1	A	228	ARG
1	C	204	SER

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	203/203 (100%)	200 (98%)	3 (2%)	65	86
1	B	203/203 (100%)	196 (97%)	7 (3%)	37	66
1	C	203/203 (100%)	196 (97%)	7 (3%)	37	66
1	D	203/203 (100%)	194 (96%)	9 (4%)	28	56
1	E	203/203 (100%)	197 (97%)	6 (3%)	41	70
1	F	203/203 (100%)	197 (97%)	6 (3%)	41	70
1	G	203/203 (100%)	197 (97%)	6 (3%)	41	70
1	H	203/203 (100%)	197 (97%)	6 (3%)	41	70

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1624/1624 (100%)	1574 (97%)	50 (3%)	40 69

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	LYS
1	A	120	THR
1	A	170	VAL
1	B	101	LYS
1	B	117	SER
1	B	170	VAL
1	B	204	SER
1	B	208	ASP
1	B	228	ARG
1	B	237	ASN
1	C	21	SER
1	C	101	LYS
1	C	120	THR
1	C	161	SER
1	C	170	VAL
1	C	204	SER
1	C	226	THR
1	D	21	SER
1	D	31	SER
1	D	101	LYS
1	D	168	SER
1	D	170	VAL
1	D	185	SER
1	D	204	SER
1	D	228	ARG
1	D	237	ASN
1	E	101	LYS
1	E	120	THR
1	E	170	VAL
1	E	204	SER
1	E	223	SER
1	E	228	ARG
1	F	101	LYS
1	F	120	THR
1	F	161	SER
1	F	170	VAL
1	F	226	THR

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Mol	Chain	Res	Type
1	F	237	ASN
1	G	101	LYS
1	G	120	THR
1	G	170	VAL
1	G	204	SER
1	G	226	THR
1	G	228	ARG
1	H	101	LYS
1	H	158	ARG
1	H	168	SER
1	H	170	VAL
1	H	204	SER
1	H	228	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	237	ASN
1	B	43	GLN
1	B	205	HIS
1	B	237	ASN
1	C	166	GLN
1	C	205	HIS
1	C	237	ASN
1	D	43	GLN
1	D	237	ASN
1	E	205	HIS
1	E	237	ASN
1	F	43	GLN
1	F	205	HIS
1	F	237	ASN
1	G	43	GLN
1	G	69	ASN
1	G	205	HIS
1	G	237	ASN
1	H	43	GLN
1	H	205	HIS
1	H	237	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

40 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAN	I	1	2	12,12,12	0.54	0	17,17,17	0.75	1 (5%)
2	MAN	I	2	2	11,11,12	0.41	0	15,15,17	0.84	0
2	NAG	I	3	2	14,14,15	0.53	0	17,19,21	0.88	1 (5%)
2	MAN	I	4	2	11,11,12	0.52	0	15,15,17	0.48	0
2	NAG	I	5	2	14,14,15	0.49	0	17,19,21	1.07	2 (11%)
2	MAN	J	1	2	12,12,12	0.45	0	17,17,17	0.73	0
2	MAN	J	2	2	11,11,12	0.55	0	15,15,17	0.67	0
2	NAG	J	3	2	14,14,15	0.59	0	17,19,21	0.88	1 (5%)
2	MAN	J	4	2	11,11,12	0.33	0	15,15,17	0.66	0
2	NAG	J	5	2	14,14,15	0.40	0	17,19,21	0.98	0
2	MAN	K	1	2	12,12,12	0.33	0	17,17,17	0.70	0
2	MAN	K	2	2	11,11,12	0.65	0	15,15,17	0.51	0
2	NAG	K	3	2	14,14,15	0.48	0	17,19,21	0.94	1 (5%)
2	MAN	K	4	2	11,11,12	0.54	0	15,15,17	0.45	0
2	NAG	K	5	2	14,14,15	0.41	0	17,19,21	1.05	0
2	MAN	L	1	2	12,12,12	0.49	0	17,17,17	0.68	0
2	MAN	L	2	2	11,11,12	0.45	0	15,15,17	0.70	0
2	NAG	L	3	2	14,14,15	0.56	0	17,19,21	1.09	2 (11%)
2	MAN	L	4	2	11,11,12	0.28	0	15,15,17	0.56	0
2	NAG	L	5	2	14,14,15	0.42	0	17,19,21	0.90	0
2	MAN	M	1	2	12,12,12	0.40	0	17,17,17	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAN	M	2	2	11,11,12	0.41	0	15,15,17	0.72	0
2	NAG	M	3	2	14,14,15	0.38	0	17,19,21	0.82	1 (5%)
2	MAN	M	4	2	11,11,12	0.41	0	15,15,17	0.48	0
2	NAG	M	5	2	14,14,15	0.46	0	17,19,21	1.16	2 (11%)
2	MAN	N	1	2	12,12,12	0.43	0	17,17,17	0.73	0
2	MAN	N	2	2	11,11,12	0.71	0	15,15,17	0.58	0
2	NAG	N	3	2	14,14,15	0.60	0	17,19,21	1.09	2 (11%)
2	MAN	N	4	2	11,11,12	0.39	0	15,15,17	0.52	0
2	NAG	N	5	2	14,14,15	0.58	0	17,19,21	0.99	0
2	MAN	O	1	2	12,12,12	0.50	0	17,17,17	0.68	0
2	MAN	O	2	2	11,11,12	0.45	0	15,15,17	0.76	0
2	NAG	O	3	2	14,14,15	0.51	0	17,19,21	0.65	0
2	MAN	O	4	2	11,11,12	0.40	0	15,15,17	0.53	0
2	NAG	O	5	2	14,14,15	0.39	0	17,19,21	1.07	1 (5%)
2	MAN	P	1	2	12,12,12	0.53	0	17,17,17	0.66	0
2	MAN	P	2	2	11,11,12	0.50	0	15,15,17	0.82	0
2	NAG	P	3	2	14,14,15	0.76	1 (7%)	17,19,21	0.80	1 (5%)
2	MAN	P	4	2	11,11,12	0.52	0	15,15,17	0.57	0
2	NAG	P	5	2	14,14,15	0.65	0	17,19,21	0.76	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
2	MAN	I	1	2	-	1/2/22/22	0/1/1/1
2	MAN	I	2	2	-	0/2/19/22	0/1/1/1
2	NAG	I	3	2	-	0/6/23/26	0/1/1/1
2	MAN	I	4	2	-	0/2/19/22	0/1/1/1
2	NAG	I	5	2	-	0/6/23/26	0/1/1/1
2	MAN	J	1	2	-	0/2/22/22	0/1/1/1
2	MAN	J	2	2	-	2/2/19/22	0/1/1/1
2	NAG	J	3	2	-	0/6/23/26	0/1/1/1
2	MAN	J	4	2	-	2/2/19/22	0/1/1/1
2	NAG	J	5	2	-	2/6/23/26	0/1/1/1
2	MAN	K	1	2	-	0/2/22/22	0/1/1/1
2	MAN	K	2	2	-	2/2/19/22	0/1/1/1
2	NAG	K	3	2	-	0/6/23/26	0/1/1/1
2	MAN	K	4	2	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	K	5	2	-	2/6/23/26	0/1/1/1
2	MAN	L	1	2	-	0/2/22/22	0/1/1/1
2	MAN	L	2	2	-	2/2/19/22	0/1/1/1
2	NAG	L	3	2	-	0/6/23/26	0/1/1/1
2	MAN	L	4	2	-	2/2/19/22	0/1/1/1
2	NAG	L	5	2	-	0/6/23/26	0/1/1/1
2	MAN	M	1	2	-	0/2/22/22	0/1/1/1
2	MAN	M	2	2	-	2/2/19/22	0/1/1/1
2	NAG	M	3	2	-	0/6/23/26	0/1/1/1
2	MAN	M	4	2	-	2/2/19/22	0/1/1/1
2	NAG	M	5	2	-	2/6/23/26	0/1/1/1
2	MAN	N	1	2	-	0/2/22/22	0/1/1/1
2	MAN	N	2	2	-	2/2/19/22	0/1/1/1
2	NAG	N	3	2	-	0/6/23/26	0/1/1/1
2	MAN	N	4	2	-	2/2/19/22	0/1/1/1
2	NAG	N	5	2	-	2/6/23/26	0/1/1/1
2	MAN	O	1	2	-	0/2/22/22	0/1/1/1
2	MAN	O	2	2	-	1/2/19/22	0/1/1/1
2	NAG	O	3	2	-	0/6/23/26	0/1/1/1
2	MAN	O	4	2	-	0/2/19/22	0/1/1/1
2	NAG	O	5	2	-	0/6/23/26	0/1/1/1
2	MAN	P	1	2	-	0/2/22/22	0/1/1/1
2	MAN	P	2	2	-	2/2/19/22	0/1/1/1
2	NAG	P	3	2	-	0/6/23/26	0/1/1/1
2	MAN	P	4	2	-	2/2/19/22	0/1/1/1
2	NAG	P	5	2	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	3	NAG	C1-C2	2.15	1.55	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	3	NAG	C1-O5-C5	3.22	116.55	112.19
2	J	3	NAG	C1-O5-C5	2.85	116.05	112.19
2	K	3	NAG	C1-O5-C5	2.71	115.86	112.19
2	M	5	NAG	C6-C5-C4	2.64	119.18	113.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	5	NAG	C3-C4-C5	-2.43	105.90	110.24
2	I	5	NAG	C1-C2-N2	-2.43	106.34	110.49
2	I	5	NAG	C6-C5-C4	2.33	118.47	113.00
2	M	5	NAG	C3-C4-C5	-2.33	106.08	110.24
2	N	3	NAG	C2-N2-C7	-2.28	119.65	122.90
2	M	3	NAG	C1-O5-C5	2.21	115.19	112.19
2	L	3	NAG	C2-N2-C7	-2.14	119.86	122.90
2	I	3	NAG	C1-O5-C5	2.10	115.03	112.19
2	N	3	NAG	C1-O5-C5	2.09	115.03	112.19
2	P	3	NAG	C1-O5-C5	2.02	114.93	112.19
2	I	1	MAN	C6-C5-C4	-2.00	108.32	113.00

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	M	5	NAG	C4-C5-C6-O6
2	K	2	MAN	O5-C5-C6-O6
2	K	2	MAN	C4-C5-C6-O6
2	M	5	NAG	O5-C5-C6-O6
2	L	4	MAN	C4-C5-C6-O6
2	N	5	NAG	C4-C5-C6-O6
2	N	4	MAN	C4-C5-C6-O6
2	N	4	MAN	O5-C5-C6-O6
2	N	5	NAG	O5-C5-C6-O6
2	L	4	MAN	O5-C5-C6-O6
2	P	4	MAN	C4-C5-C6-O6
2	N	2	MAN	C4-C5-C6-O6
2	N	2	MAN	O5-C5-C6-O6
2	L	2	MAN	C4-C5-C6-O6
2	J	2	MAN	C4-C5-C6-O6
2	P	4	MAN	O5-C5-C6-O6
2	J	5	NAG	C4-C5-C6-O6
2	M	4	MAN	C4-C5-C6-O6
2	P	5	NAG	C4-C5-C6-O6
2	J	4	MAN	C4-C5-C6-O6
2	M	4	MAN	O5-C5-C6-O6
2	L	2	MAN	O5-C5-C6-O6
2	K	4	MAN	C4-C5-C6-O6
2	K	5	NAG	C4-C5-C6-O6
2	J	2	MAN	O5-C5-C6-O6
2	P	5	NAG	O5-C5-C6-O6

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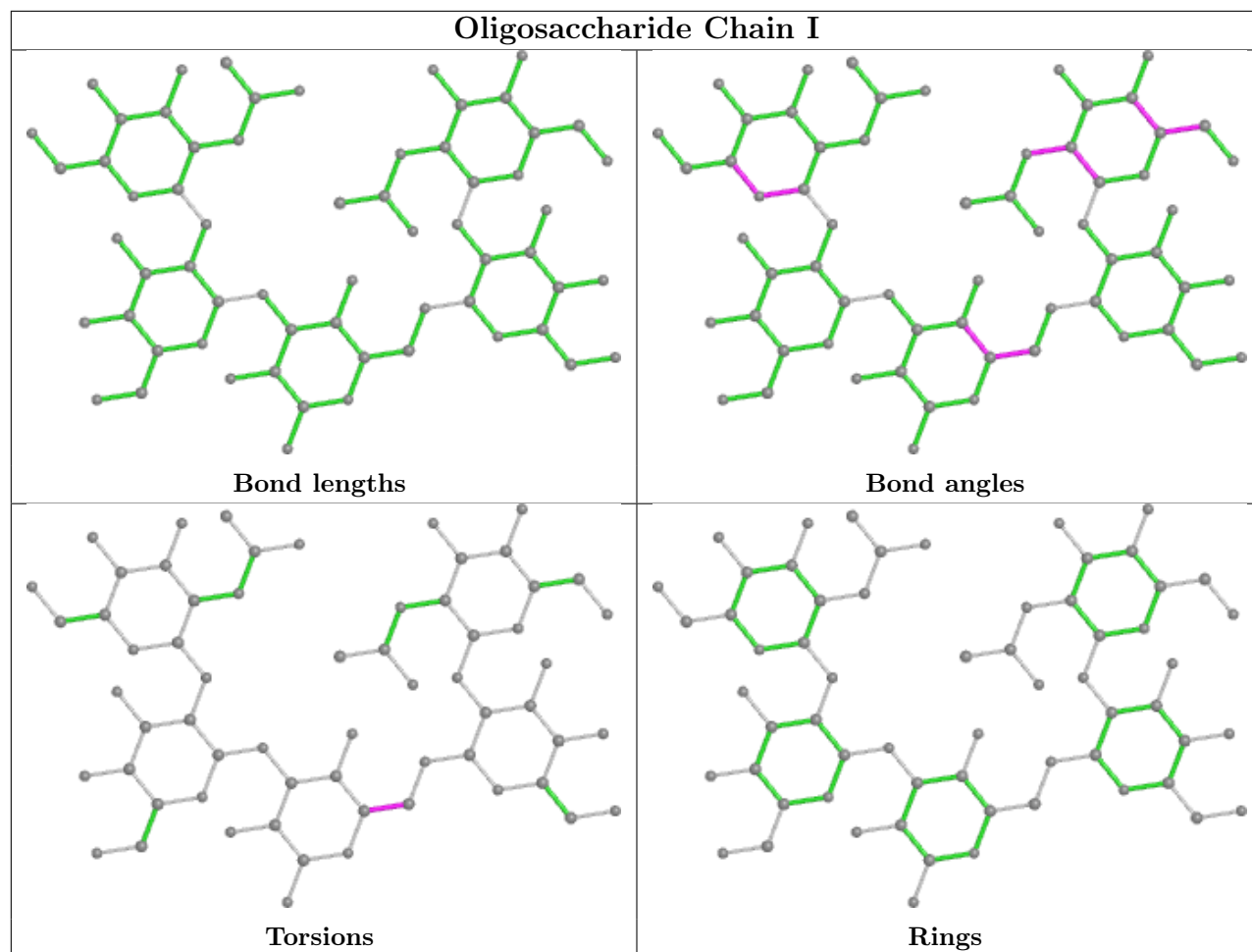
Mol	Chain	Res	Type	Atoms
2	J	5	NAG	O5-C5-C6-O6
2	P	2	MAN	C4-C5-C6-O6
2	I	1	MAN	O5-C5-C6-O6
2	K	4	MAN	O5-C5-C6-O6
2	J	4	MAN	O5-C5-C6-O6
2	M	2	MAN	C4-C5-C6-O6
2	K	5	NAG	O5-C5-C6-O6
2	M	2	MAN	O5-C5-C6-O6
2	O	2	MAN	C4-C5-C6-O6
2	P	2	MAN	O5-C5-C6-O6

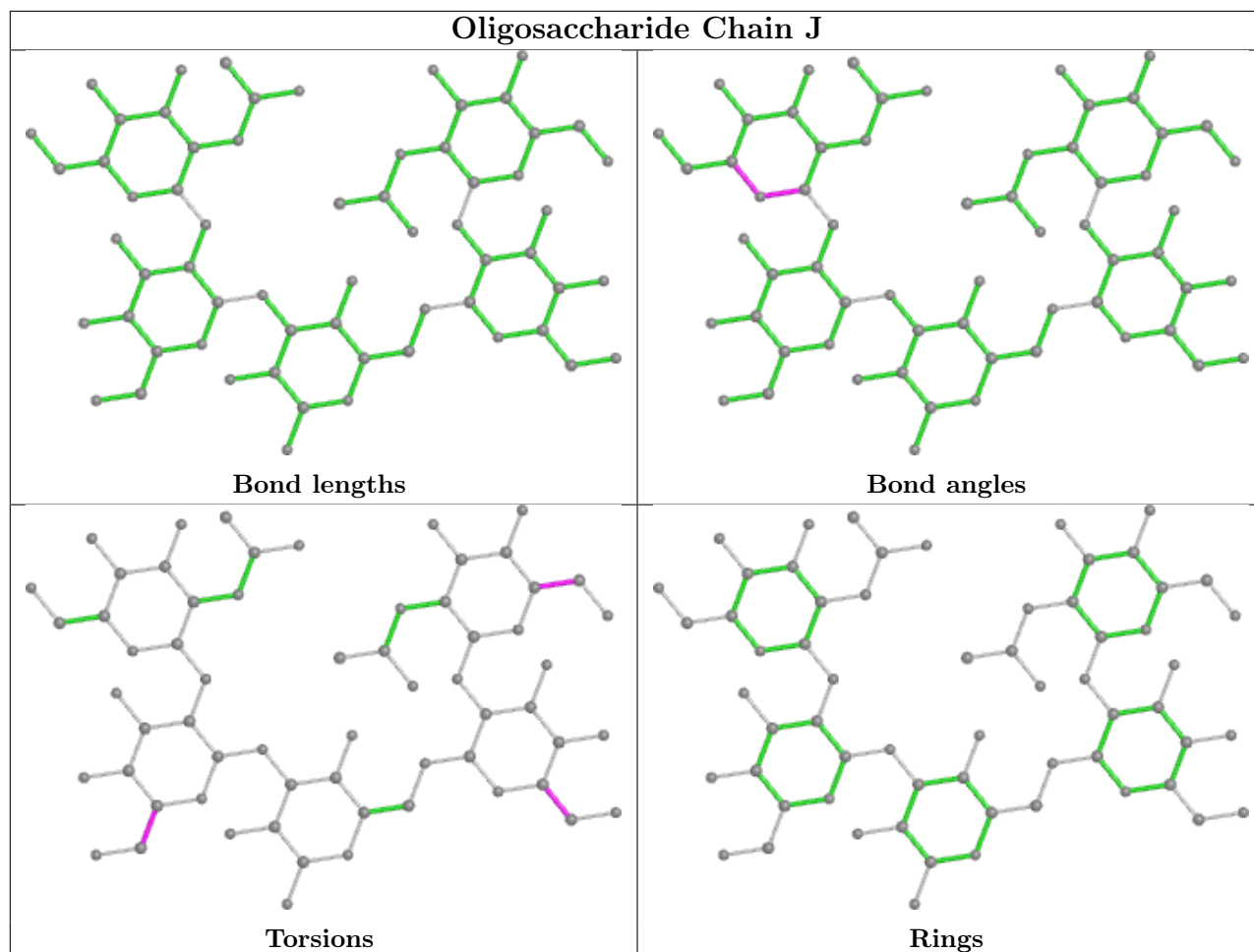
There are no ring outliers.

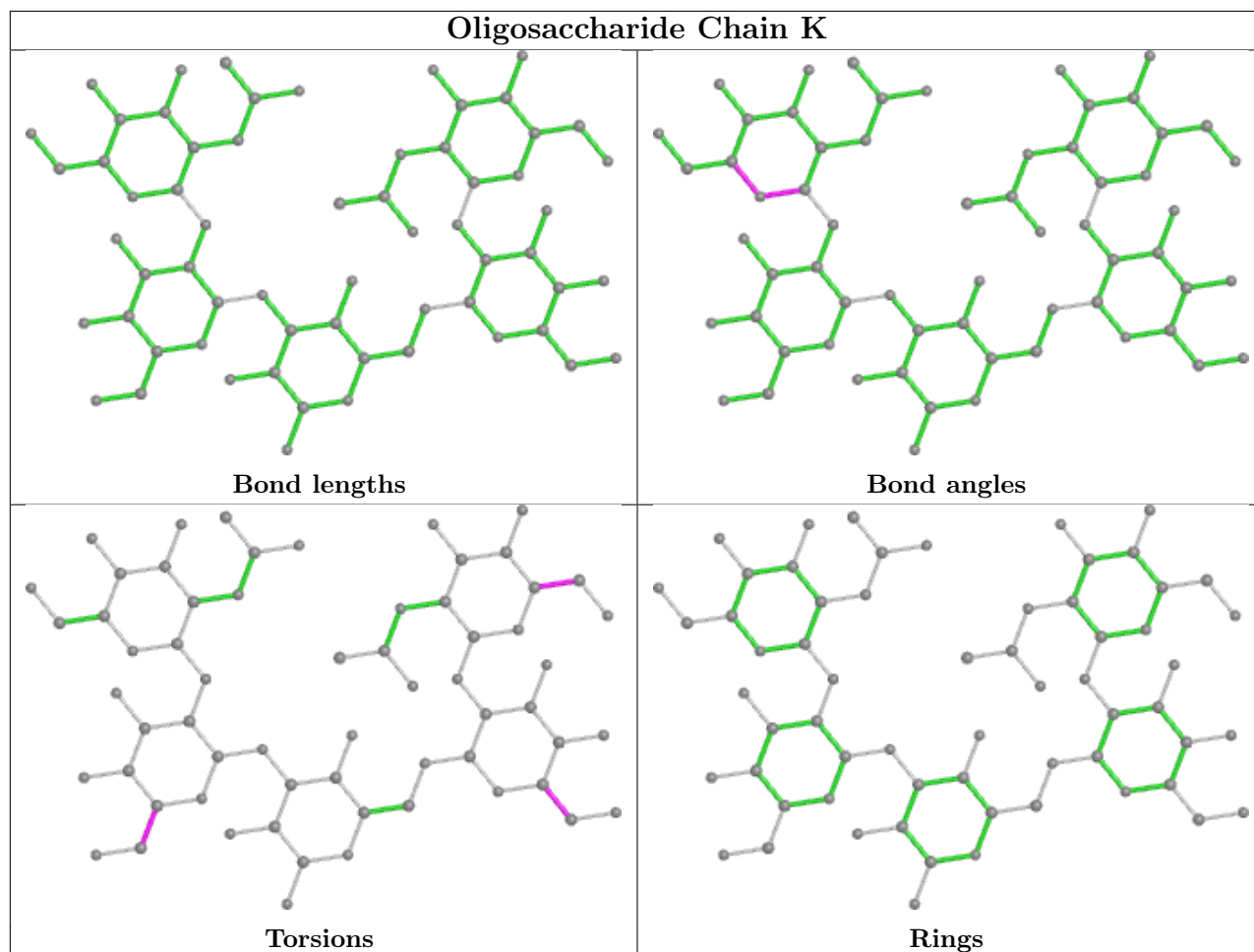
10 monomers are involved in 10 short contacts:

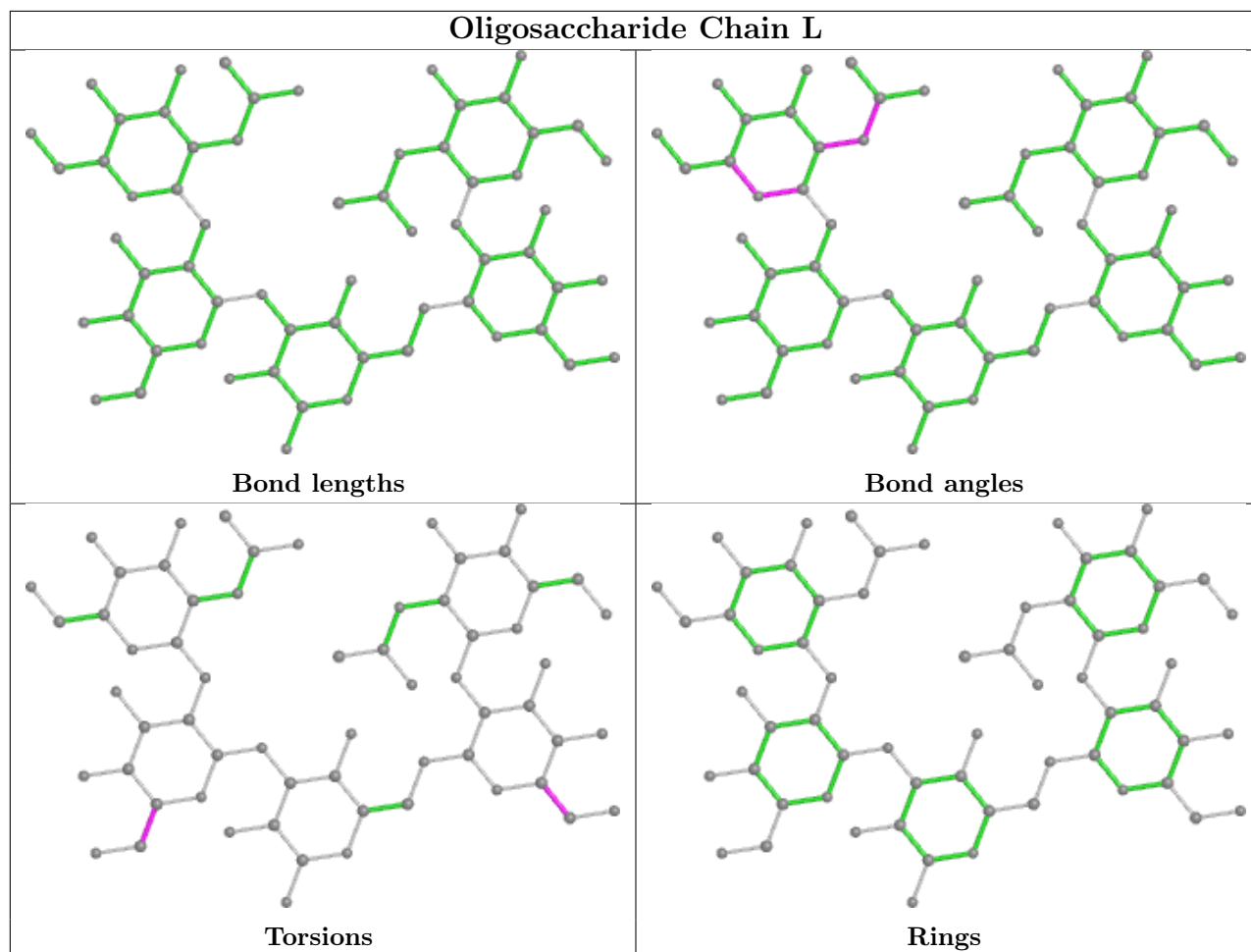
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	4	MAN	1	0
2	N	3	NAG	0	1
2	N	5	NAG	1	0
2	J	4	MAN	1	0
2	I	4	MAN	1	0
2	L	5	NAG	1	0
2	M	4	MAN	1	0
2	O	4	MAN	1	0
2	L	4	MAN	1	0
2	P	4	MAN	1	0

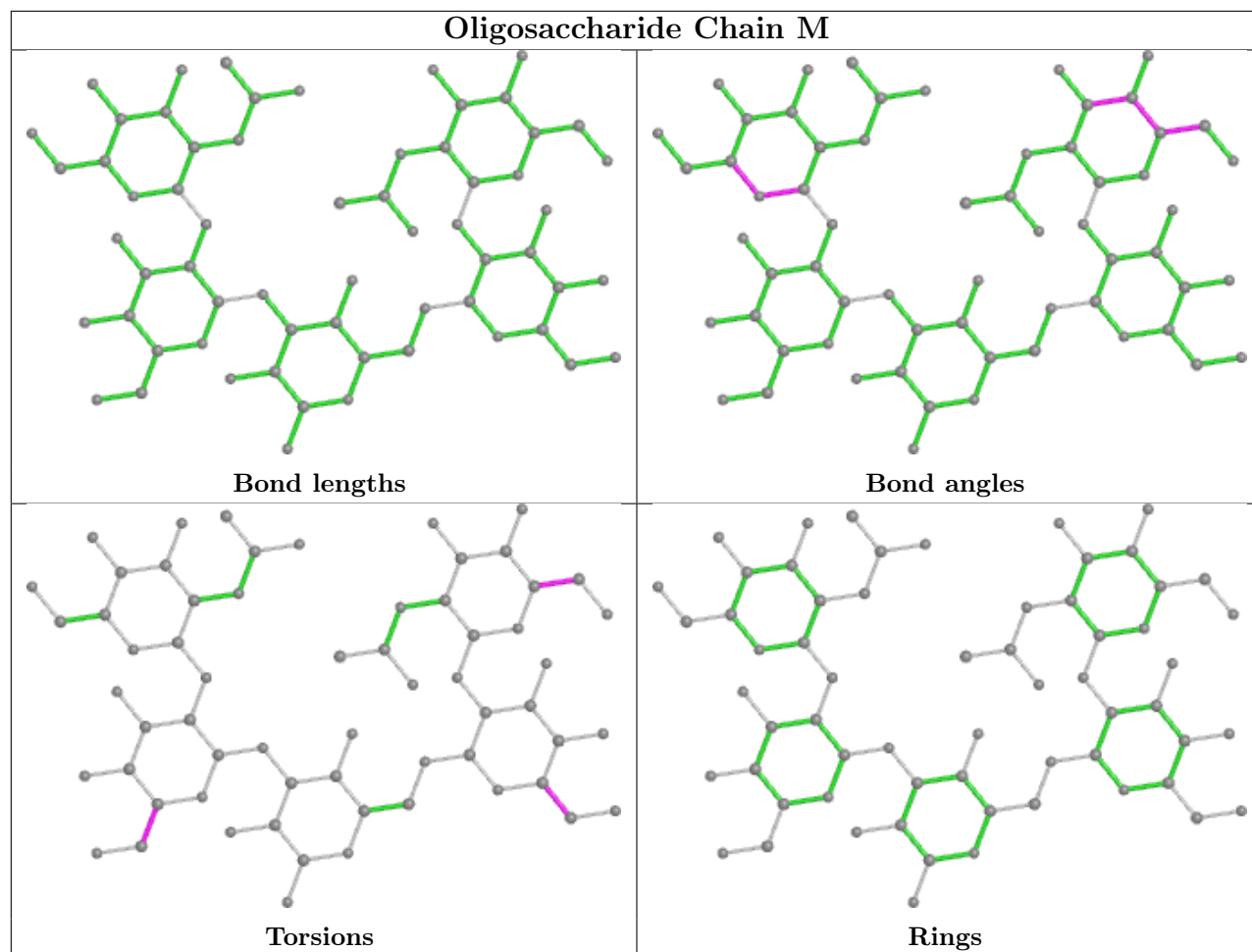
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

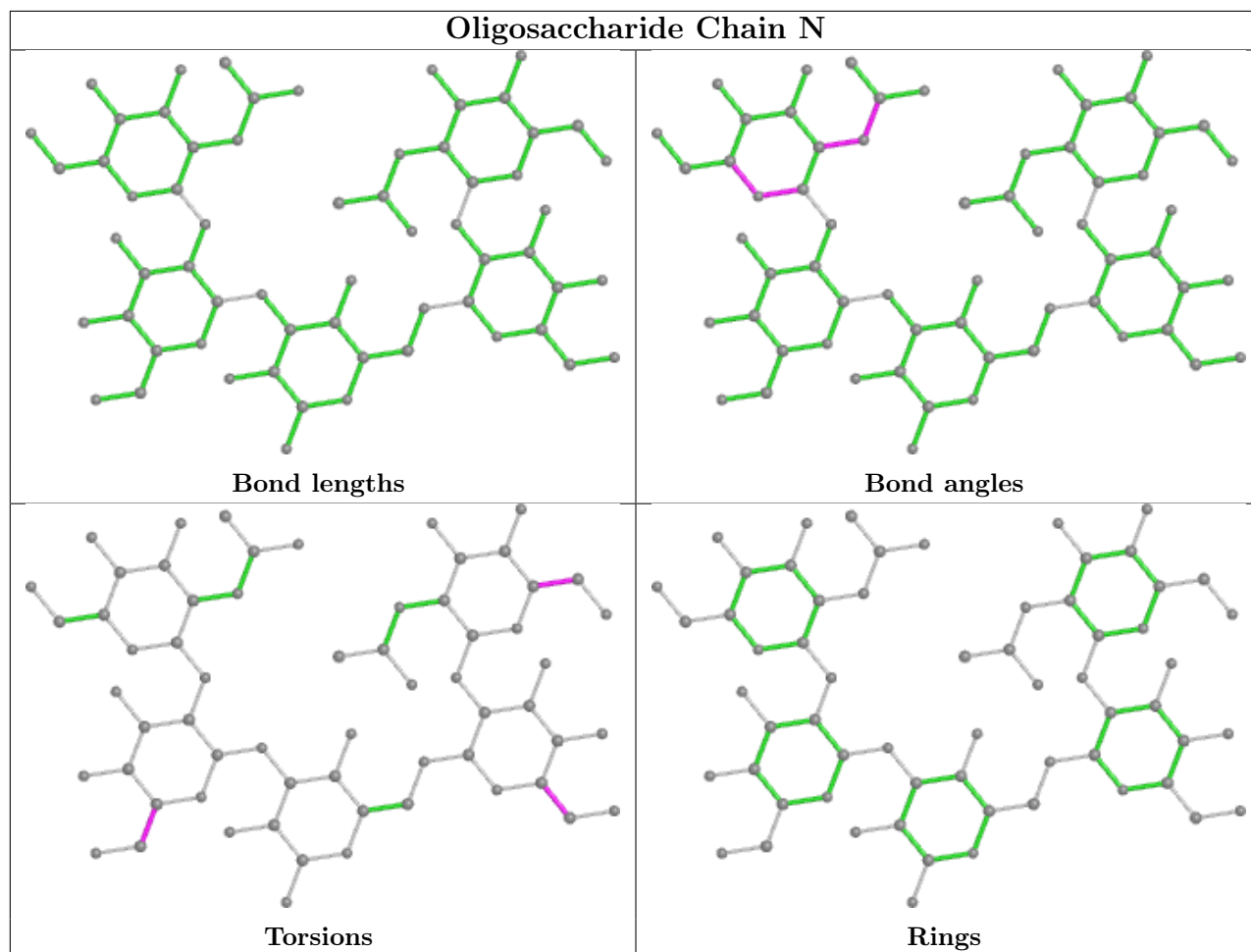


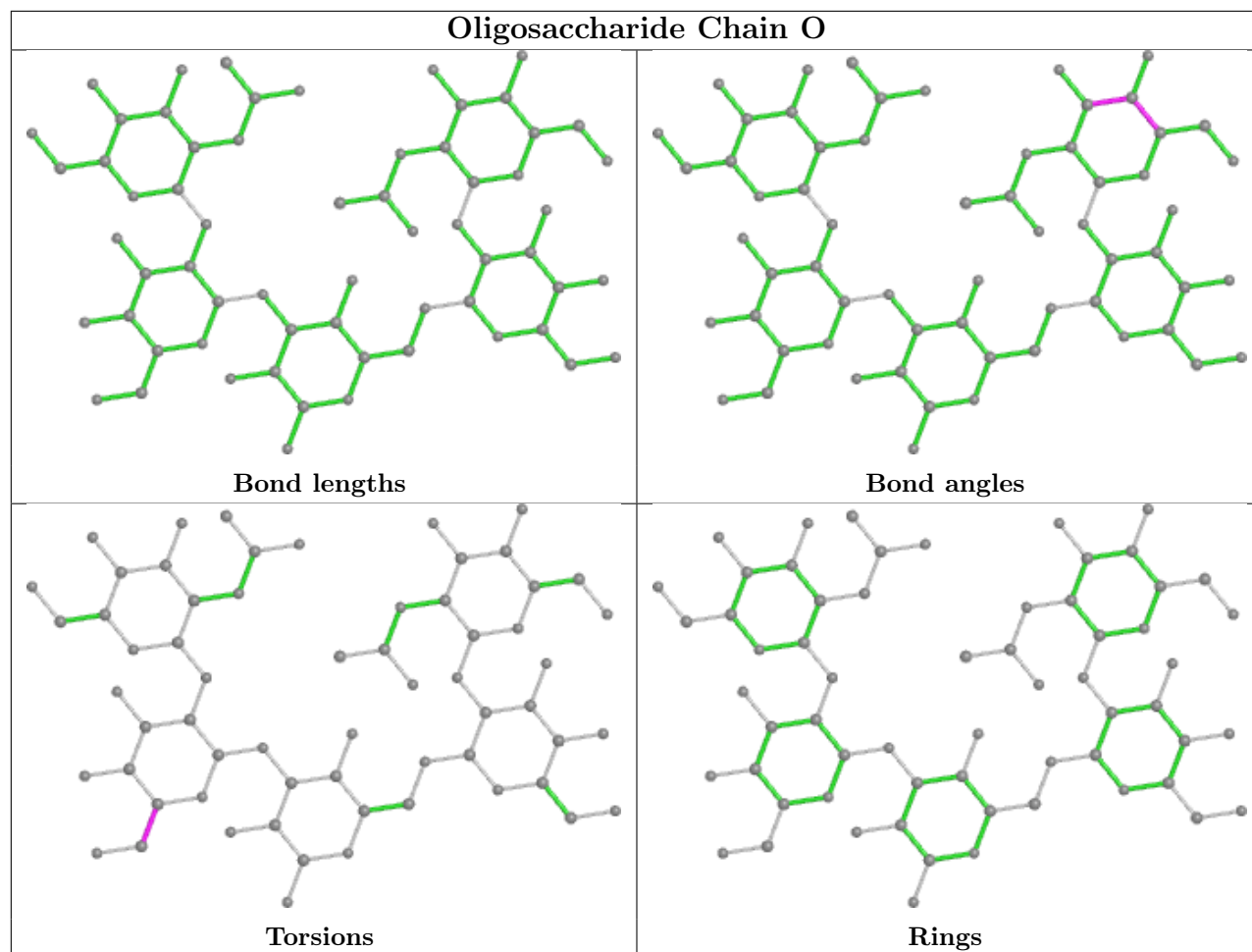


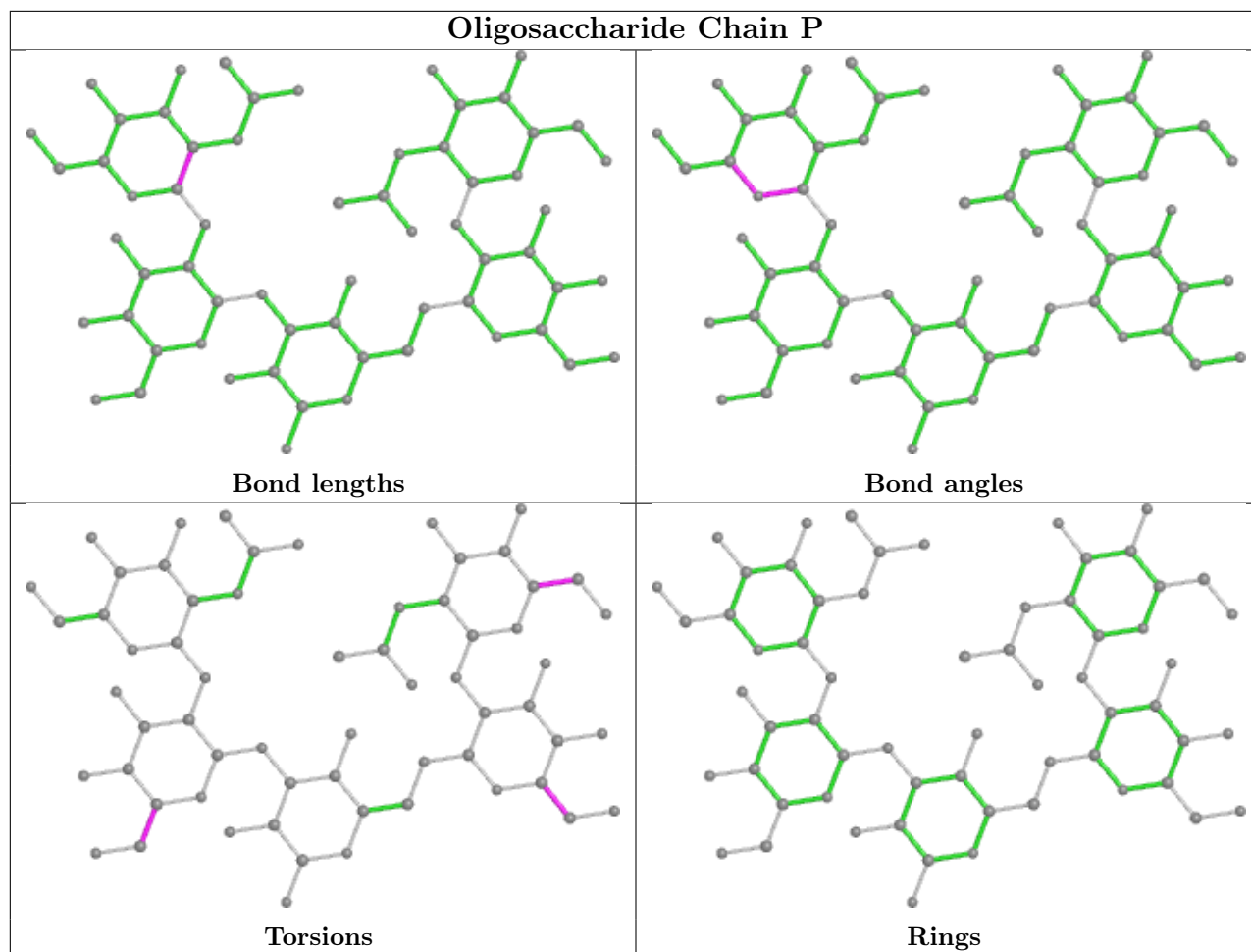












5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 16 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

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	235/237 (99%)	-0.29	8 (3%) 45 45	6, 21, 49, 83	5 (2%)
1	B	232/237 (97%)	-0.23	7 (3%) 50 51	8, 27, 54, 70	3 (1%)
1	C	235/237 (99%)	-0.35	3 (1%) 77 78	6, 24, 55, 91	3 (1%)
1	D	233/237 (98%)	-0.18	6 (2%) 56 57	9, 26, 52, 68	7 (3%)
1	E	236/237 (99%)	-0.17	11 (4%) 31 30	7, 23, 54, 87	5 (2%)
1	F	234/237 (98%)	0.18	21 (8%) 9 7	13, 34, 61, 84	11 (4%)
1	G	237/237 (100%)	-0.13	12 (5%) 28 26	9, 27, 55, 70	5 (2%)
1	H	232/237 (97%)	0.13	11 (4%) 31 30	10, 32, 58, 73	8 (3%)
All	All	1874/1896 (98%)	-0.13	79 (4%) 36 35	6, 27, 56, 91	47 (2%)

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	121	HIS	9.6
1	F	120	THR	9.6
1	C	120	THR	9.2
1	E	120	THR	8.3
1	G	120	THR	7.4
1	A	120	THR	7.2
1	E	121	HIS	6.0
1	F	150	THR	5.0
1	B	1	ALA	4.9
1	E	185	SER	4.7
1	A	1	ALA	4.6
1	F	98	GLY	4.6
1	A	122	GLU	4.3
1	H	122	GLU	4.1
1	D	120	THR	4.0
1	E	119	SER	4.0

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Mol	Chain	Res	Type	RSRZ
1	G	69	ASN	3.9
1	D	204	SER	3.9
1	E	187	VAL	3.8
1	F	163	GLY	3.7
1	F	205	HIS	3.7
1	F	69	ASN	3.6
1	F	161	SER	3.5
1	H	69	ASN	3.5
1	H	162	ASN	3.4
1	F	162	ASN	3.2
1	F	71	ASP	3.1
1	H	160	SER	3.0
1	F	204	SER	3.0
1	G	150	THR	3.0
1	H	21	SER	3.0
1	C	119	SER	2.9
1	D	161	SER	2.9
1	G	161	SER	2.8
1	E	117	SER	2.8
1	G	205	HIS	2.7
1	A	69	ASN	2.7
1	C	204	SER	2.7
1	H	185	SER	2.7
1	G	1	ALA	2.6
1	D	69	ASN	2.6
1	H	71	ASP	2.6
1	F	160	SER	2.6
1	A	119	SER	2.6
1	B	7	VAL	2.5
1	B	71	ASP	2.5
1	G	70	ALA	2.5
1	A	160	SER	2.5
1	B	162	ASN	2.5
1	B	211	ALA	2.4
1	F	9	LEU	2.4
1	H	83	ASN	2.4
1	H	211	ALA	2.4
1	H	13	PRO	2.3
1	D	1	ALA	2.3
1	A	162	ASN	2.3
1	E	188	VAL	2.3
1	F	236	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	G	160	SER	2.2
1	F	166	GLN	2.2
1	F	68	PRO	2.2
1	E	204	SER	2.2
1	A	161	SER	2.2
1	H	16	ASP	2.2
1	G	162	ASN	2.2
1	F	39	LYS	2.2
1	E	164	SER	2.1
1	E	123	THR	2.1
1	G	119	SER	2.1
1	F	210	ILE	2.1
1	B	69	ASN	2.1
1	D	162	ASN	2.1
1	F	164	SER	2.1
1	F	41	ASN	2.1
1	E	186	ALA	2.1
1	F	225	SER	2.1
1	G	204	SER	2.1
1	B	117	SER	2.0
1	F	211	ALA	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(Å ²)	Q<0.9
2	NAG	P	3	14/15	0.49	0.25	59,64,65,66	10
2	MAN	N	2	11/12	0.71	0.33	23,27,31,32	0
2	NAG	N	3	14/15	0.81	0.30	22,28,31,32	11
2	NAG	L	3	14/15	0.83	0.27	46,51,55,56	0
2	MAN	P	2	11/12	0.84	0.36	53,55,58,62	0
2	MAN	N	1	12/12	0.85	0.27	47,50,53,56	0
2	NAG	O	3	14/15	0.85	0.37	51,54,60,63	0

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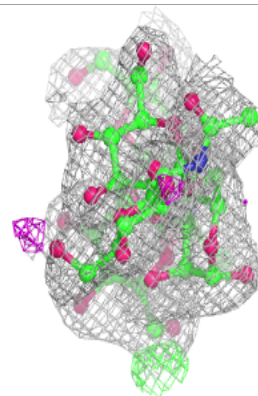
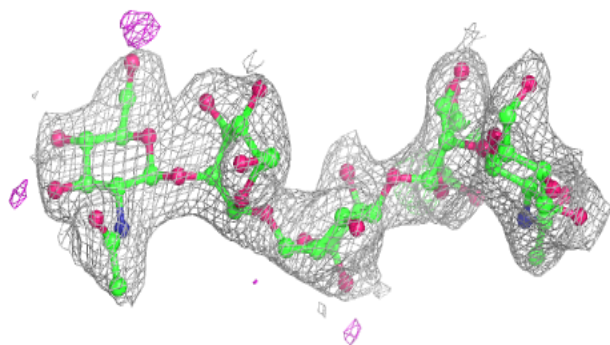
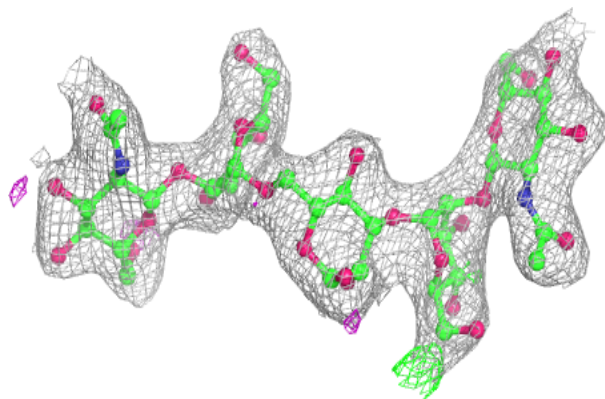
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MAN	P	1	12/12	0.86	0.30	46,49,55,55	0
2	MAN	N	4	11/12	0.87	0.20	37,41,46,48	0
2	NAG	J	3	14/15	0.88	0.23	41,45,50,53	0
2	NAG	N	5	14/15	0.88	0.28	35,42,48,52	0
2	NAG	K	3	14/15	0.88	0.30	47,49,51,51	0
2	MAN	P	4	11/12	0.89	0.22	36,41,46,49	0
2	NAG	M	3	14/15	0.91	0.18	33,37,41,42	0
2	NAG	P	5	14/15	0.91	0.25	29,35,50,50	0
2	MAN	J	1	12/12	0.92	0.16	35,39,45,46	0
2	MAN	L	1	12/12	0.93	0.15	37,38,40,46	0
2	MAN	K	2	11/12	0.93	0.19	42,44,47,52	0
2	MAN	O	1	12/12	0.93	0.22	34,39,44,47	0
2	NAG	O	5	14/15	0.94	0.14	26,30,35,36	0
2	NAG	J	5	14/15	0.94	0.18	23,30,38,38	0
2	NAG	I	3	14/15	0.94	0.12	26,32,35,35	0
2	MAN	O	2	11/12	0.94	0.27	41,43,48,54	0
2	MAN	L	2	11/12	0.94	0.17	30,35,41,47	0
2	MAN	O	4	11/12	0.94	0.13	25,31,32,33	0
2	MAN	J	2	11/12	0.95	0.13	34,35,38,44	0
2	NAG	K	5	14/15	0.95	0.15	20,23,33,41	0
2	NAG	L	5	14/15	0.95	0.14	19,22,32,34	0
2	MAN	M	2	11/12	0.95	0.15	24,26,32,39	0
2	MAN	J	4	11/12	0.95	0.16	27,32,36,40	0
2	MAN	K	1	12/12	0.96	0.13	32,35,36,39	0
2	MAN	M	4	11/12	0.96	0.10	21,23,26,27	0
2	NAG	M	5	14/15	0.96	0.13	11,16,21,22	0
2	MAN	I	1	12/12	0.96	0.09	16,26,29,29	0
2	MAN	M	1	12/12	0.96	0.11	27,28,32,33	0
2	MAN	I	2	11/12	0.96	0.12	17,18,23,36	0
2	MAN	L	4	11/12	0.97	0.09	19,24,27,29	0
2	NAG	I	5	14/15	0.97	0.14	14,19,27,28	0
2	MAN	K	4	11/12	0.98	0.11	19,24,25,25	0
2	MAN	I	4	11/12	0.99	0.10	4,12,17,20	0

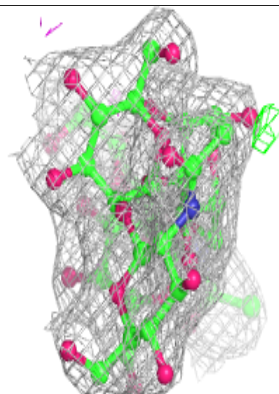
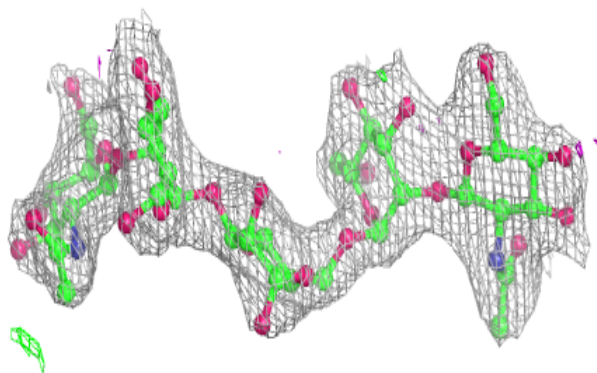
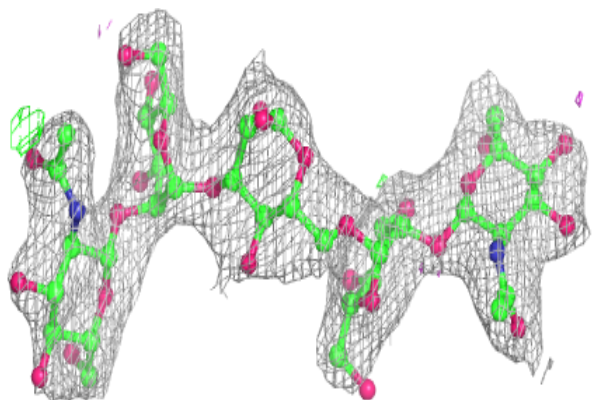
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain I:

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

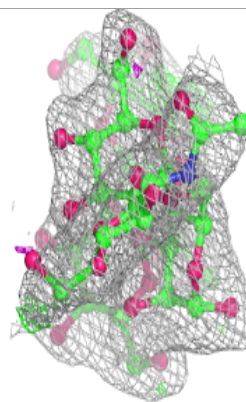
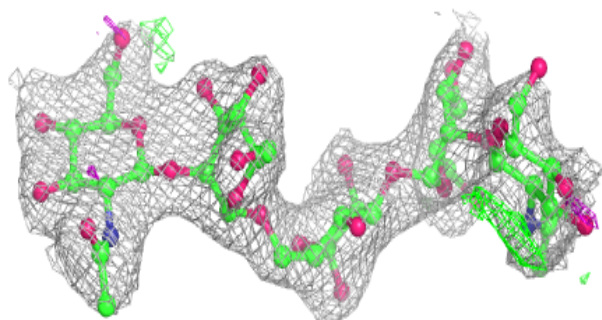
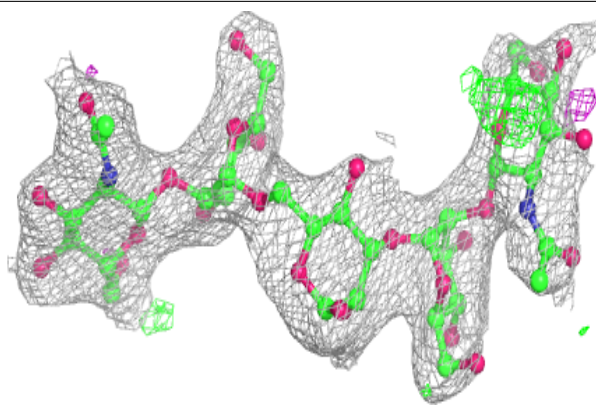
**Electron density around Chain J:**

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

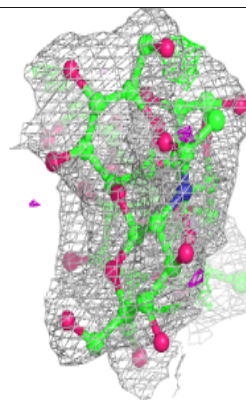
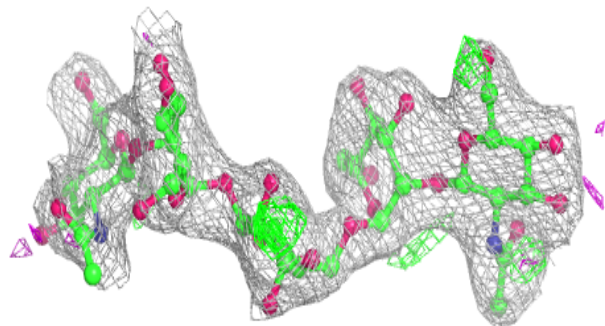
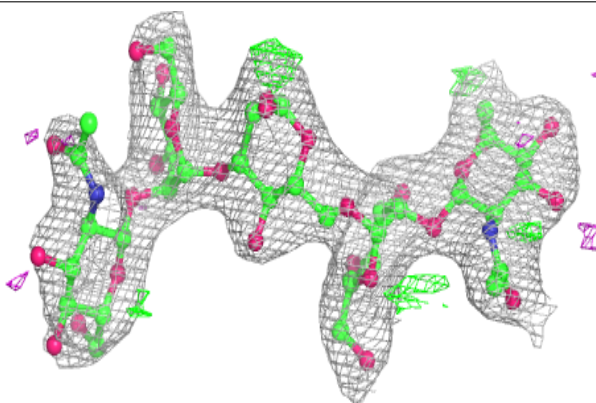


Electron density around Chain K:

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

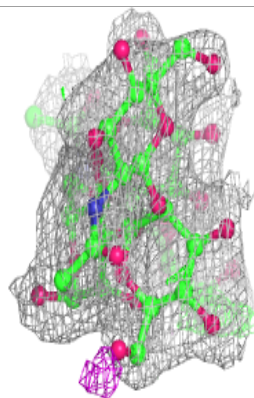
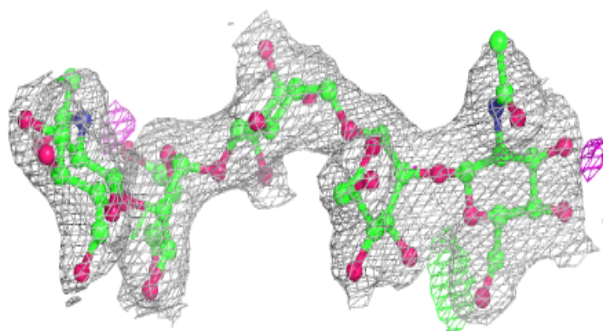
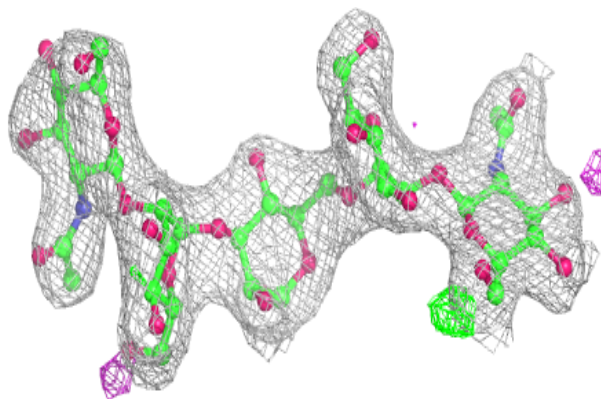
**Electron density around Chain L:**

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

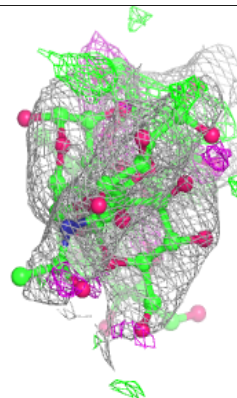
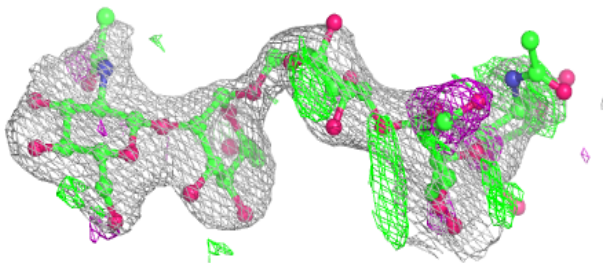
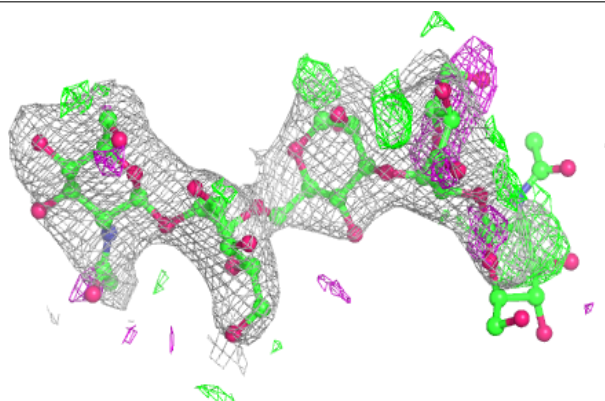


Electron density around Chain M:

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

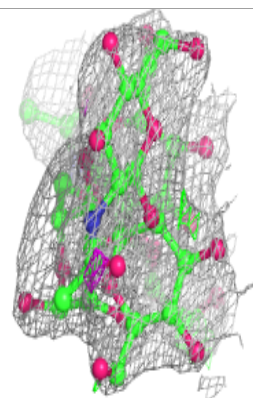
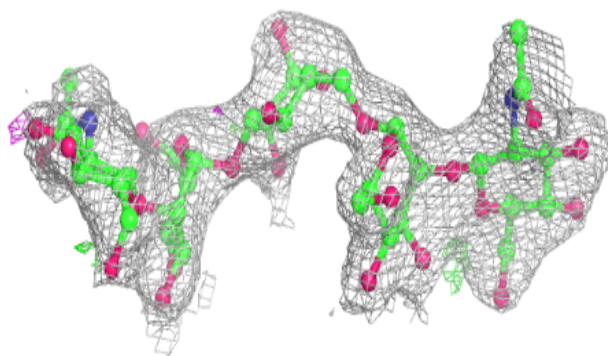
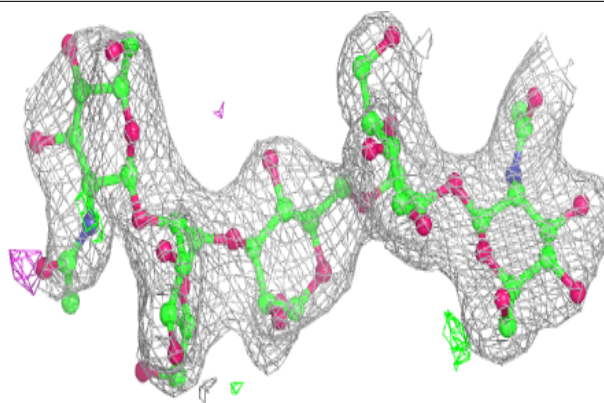
**Electron density around Chain N:**

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

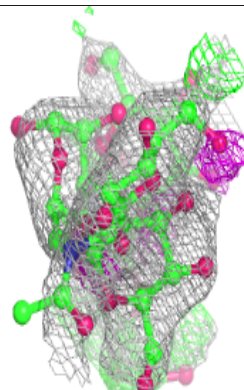
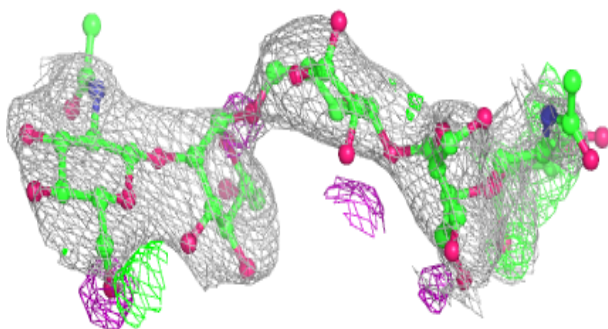
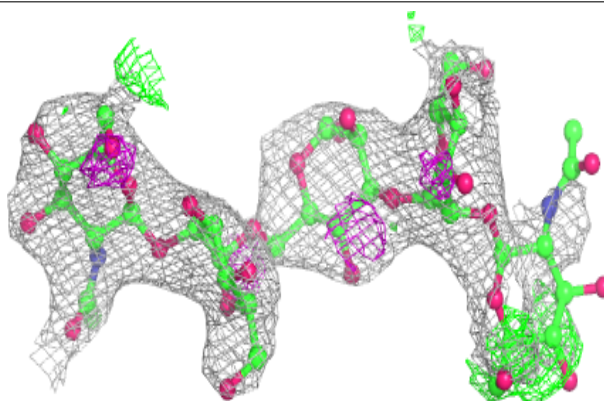


Electron density around Chain O:

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

**Electron density around Chain P:**

$2mF_o-DF_c$ (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(Å ²)	Q<0.9
4	CA	F	239	1/1	0.72	0.18	52,52,52,52	0
4	CA	H	246	1/1	0.77	0.11	47,47,47,47	0
3	MN	H	245	1/1	0.93	0.11	42,42,42,42	0
4	CA	A	239	1/1	0.94	0.09	18,18,18,18	0
4	CA	D	239	1/1	0.94	0.16	33,33,33,33	0
3	MN	F	238	1/1	0.97	0.03	50,50,50,50	0
4	CA	B	239	1/1	0.97	0.06	27,27,27,27	0
4	CA	G	239	1/1	0.97	0.12	31,31,31,31	0
4	CA	C	239	1/1	0.97	0.12	22,22,22,22	0
3	MN	D	238	1/1	0.98	0.06	32,32,32,32	0
3	MN	B	238	1/1	0.99	0.04	24,24,24,24	0
3	MN	G	238	1/1	0.99	0.07	32,32,32,32	0
4	CA	E	239	1/1	0.99	0.07	13,13,13,13	0
3	MN	C	238	1/1	0.99	0.08	24,24,24,24	0
3	MN	A	238	1/1	0.99	0.09	28,28,28,28	0
3	MN	E	238	1/1	0.99	0.08	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.