



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:57 PM UTC

PDB ID : 5OJM / pdb_00005ojm
Title : Structure of a chimaeric beta3-alpha5 GABAA receptor in complex with nanobody Nb25
Authors : Miller, P.S.; Scott, S.; Masiulis, S.; De Colibus, L.; Pardon, E.; Steyaert, J.; Aricescu, A.R.
Deposited on : 2017-07-21
Resolution : 3.30 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

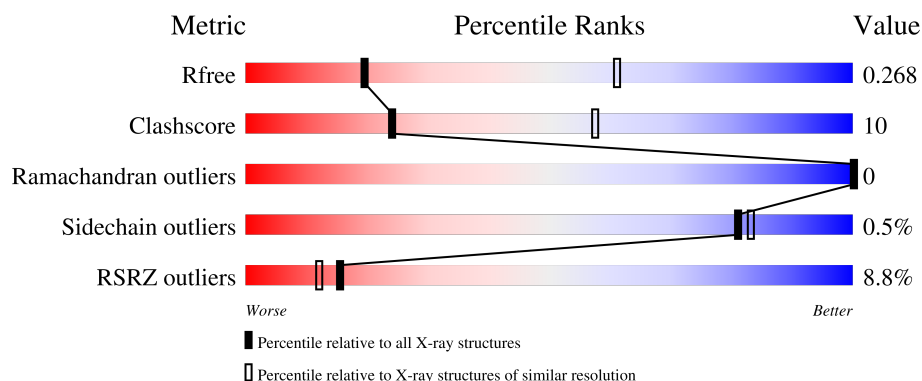
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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.



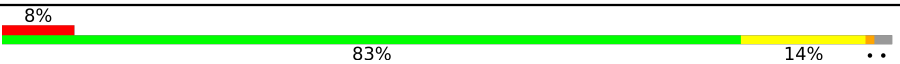
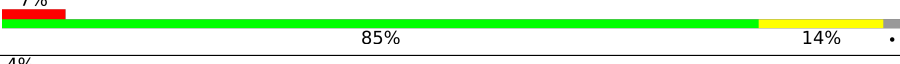
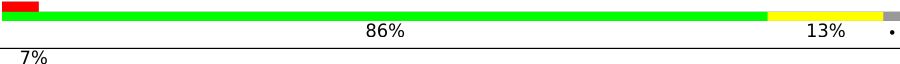
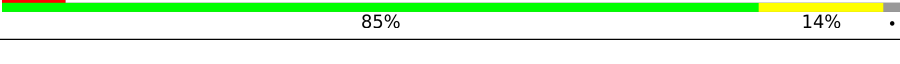
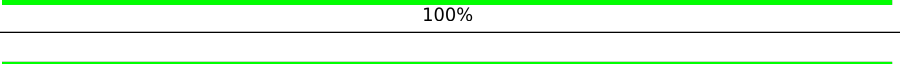
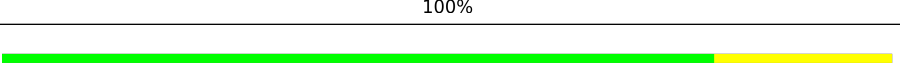
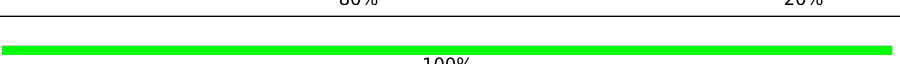
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	395	7% 69% 14% 16%
1	B	395	6% 68% 16% 16%
1	C	395	9% 66% 17% 16%
1	D	395	8% 66% 18% 16%
1	E	395	6% 67% 16% 16%

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Mol	Chain	Length	Quality of chain
2	K	125	
2	L	125	
2	M	125	
2	N	125	
2	O	125	
3	F	5	
3	G	5	
3	H	5	
4	I	6	
4	J	6	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18538 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 Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2670	1736	442	476	16			
1	B	333	Total	C	N	O	S	0	0	0
			2692	1748	446	481	17			
1	C	330	Total	C	N	O	S	0	0	0
			2670	1736	442	476	16			
1	D	332	Total	C	N	O	S	0	0	0
			2684	1744	444	479	17			
1	E	330	Total	C	N	O	S	0	0	0
			2670	1736	442	476	16			

There are 100 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	316	SER	-	linker	UNP P31644
A	317	GLN	-	linker	UNP P31644
A	318	PRO	-	linker	UNP P31644
A	319	ALA	-	linker	UNP P31644
A	320	ARG	-	linker	UNP P31644
A	321	ALA	-	linker	UNP P31644
A	322	ALA	-	linker	UNP P31644
A	404	ILE	VAL	conflict	UNP P31644
A	428	THR	ALA	conflict	UNP P31644
A	432	GLY	-	expression tag	UNP P31644
A	433	THR	-	expression tag	UNP P31644
A	434	THR	-	expression tag	UNP P31644
A	435	GLU	-	expression tag	UNP P31644
A	436	THR	-	expression tag	UNP P31644
A	437	SER	-	expression tag	UNP P31644
A	438	GLN	-	expression tag	UNP P31644
A	439	VAL	-	expression tag	UNP P31644
A	440	ALA	-	expression tag	UNP P31644

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Chain	Residue	Modelled	Actual	Comment	Reference
A	441	PRO	-	expression tag	UNP P31644
A	442	ALA	-	expression tag	UNP P31644
B	316	SER	-	linker	UNP P31644
B	317	GLN	-	linker	UNP P31644
B	318	PRO	-	linker	UNP P31644
B	319	ALA	-	linker	UNP P31644
B	320	ARG	-	linker	UNP P31644
B	321	ALA	-	linker	UNP P31644
B	322	ALA	-	linker	UNP P31644
B	404	ILE	VAL	conflict	UNP P31644
B	428	THR	ALA	conflict	UNP P31644
B	432	GLY	-	expression tag	UNP P31644
B	433	THR	-	expression tag	UNP P31644
B	434	THR	-	expression tag	UNP P31644
B	435	GLU	-	expression tag	UNP P31644
B	436	THR	-	expression tag	UNP P31644
B	437	SER	-	expression tag	UNP P31644
B	438	GLN	-	expression tag	UNP P31644
B	439	VAL	-	expression tag	UNP P31644
B	440	ALA	-	expression tag	UNP P31644
B	441	PRO	-	expression tag	UNP P31644
B	442	ALA	-	expression tag	UNP P31644
C	316	SER	-	linker	UNP P31644
C	317	GLN	-	linker	UNP P31644
C	318	PRO	-	linker	UNP P31644
C	319	ALA	-	linker	UNP P31644
C	320	ARG	-	linker	UNP P31644
C	321	ALA	-	linker	UNP P31644
C	322	ALA	-	linker	UNP P31644
C	404	ILE	VAL	conflict	UNP P31644
C	428	THR	ALA	conflict	UNP P31644
C	432	GLY	-	expression tag	UNP P31644
C	433	THR	-	expression tag	UNP P31644
C	434	THR	-	expression tag	UNP P31644
C	435	GLU	-	expression tag	UNP P31644
C	436	THR	-	expression tag	UNP P31644
C	437	SER	-	expression tag	UNP P31644
C	438	GLN	-	expression tag	UNP P31644
C	439	VAL	-	expression tag	UNP P31644
C	440	ALA	-	expression tag	UNP P31644
C	441	PRO	-	expression tag	UNP P31644
C	442	ALA	-	expression tag	UNP P31644

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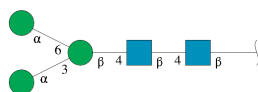
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Chain	Residue	Modelled	Actual	Comment	Reference
D	316	SER	-	linker	UNP P31644
D	317	GLN	-	linker	UNP P31644
D	318	PRO	-	linker	UNP P31644
D	319	ALA	-	linker	UNP P31644
D	320	ARG	-	linker	UNP P31644
D	321	ALA	-	linker	UNP P31644
D	322	ALA	-	linker	UNP P31644
D	404	ILE	VAL	conflict	UNP P31644
D	428	THR	ALA	conflict	UNP P31644
D	432	GLY	-	expression tag	UNP P31644
D	433	THR	-	expression tag	UNP P31644
D	434	THR	-	expression tag	UNP P31644
D	435	GLU	-	expression tag	UNP P31644
D	436	THR	-	expression tag	UNP P31644
D	437	SER	-	expression tag	UNP P31644
D	438	GLN	-	expression tag	UNP P31644
D	439	VAL	-	expression tag	UNP P31644
D	440	ALA	-	expression tag	UNP P31644
D	441	PRO	-	expression tag	UNP P31644
D	442	ALA	-	expression tag	UNP P31644
E	316	SER	-	linker	UNP P31644
E	317	GLN	-	linker	UNP P31644
E	318	PRO	-	linker	UNP P31644
E	319	ALA	-	linker	UNP P31644
E	320	ARG	-	linker	UNP P31644
E	321	ALA	-	linker	UNP P31644
E	322	ALA	-	linker	UNP P31644
E	404	ILE	VAL	conflict	UNP P31644
E	428	THR	ALA	conflict	UNP P31644
E	432	GLY	-	expression tag	UNP P31644
E	433	THR	-	expression tag	UNP P31644
E	434	THR	-	expression tag	UNP P31644
E	435	GLU	-	expression tag	UNP P31644
E	436	THR	-	expression tag	UNP P31644
E	437	SER	-	expression tag	UNP P31644
E	438	GLN	-	expression tag	UNP P31644
E	439	VAL	-	expression tag	UNP P31644
E	440	ALA	-	expression tag	UNP P31644
E	441	PRO	-	expression tag	UNP P31644
E	442	ALA	-	expression tag	UNP P31644

- Molecule 2 is a protein called Nanobody Nb25.

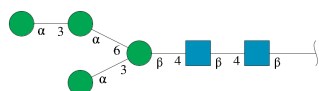
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			
2	L	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			
2	M	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			
2	N	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			
2	O	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			

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



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

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	I	6	Total	C	N	O	0	0	0
			72	40	2	30			
4	J	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

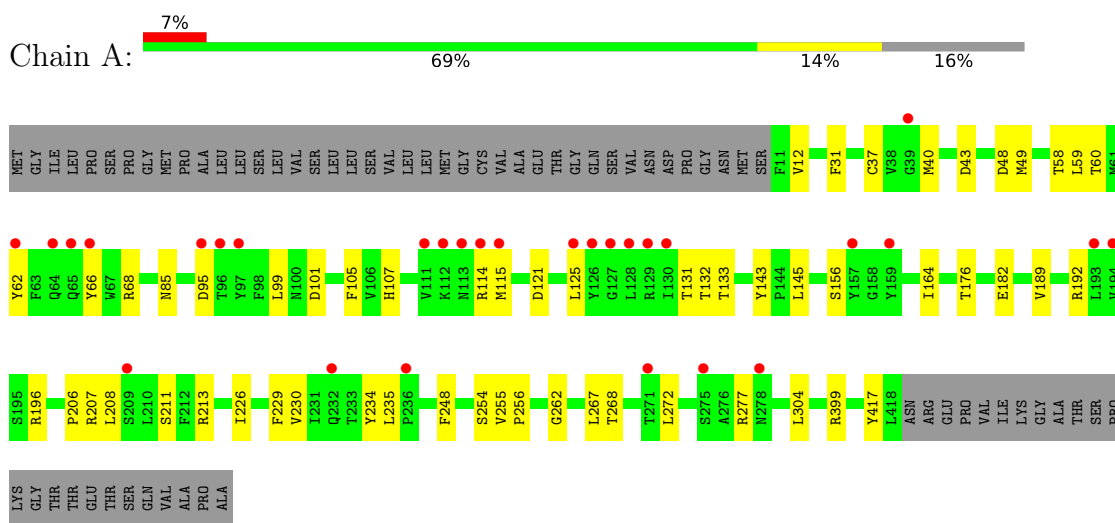


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		

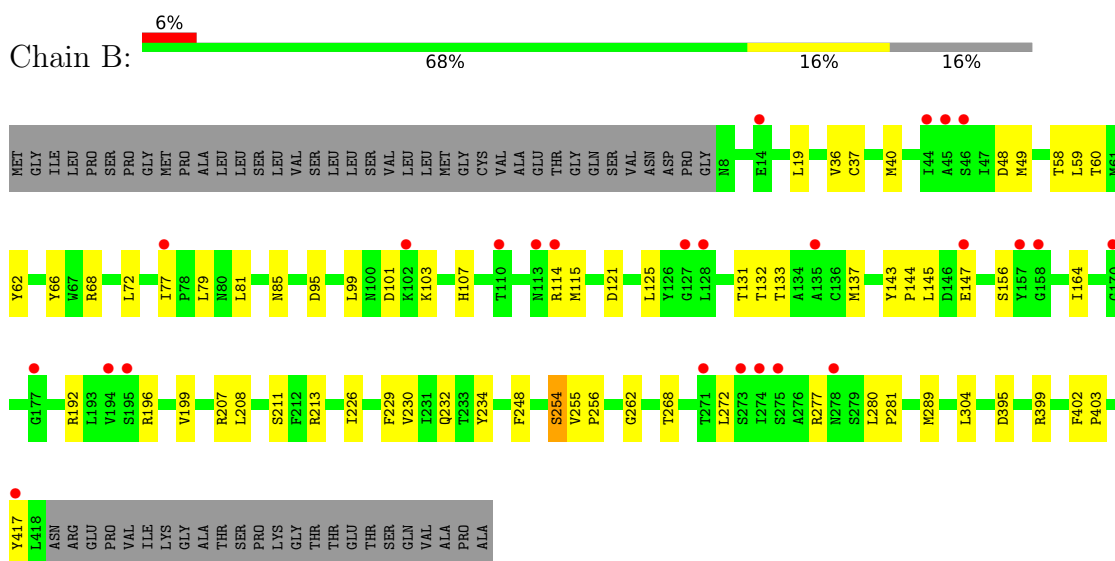
3 Residue-property plots

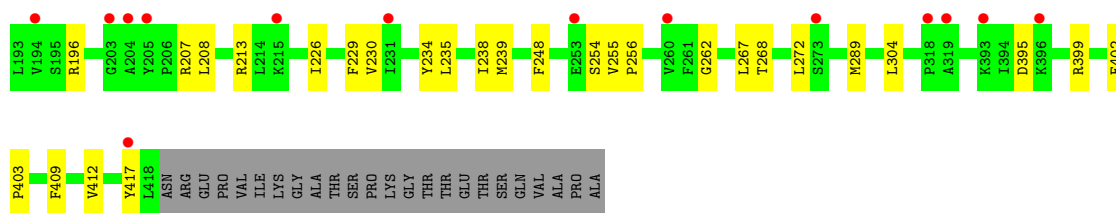
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5

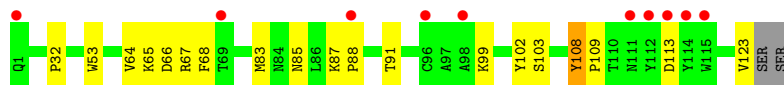
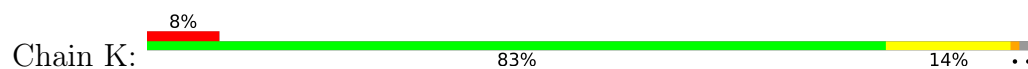


- Molecule 1: Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5

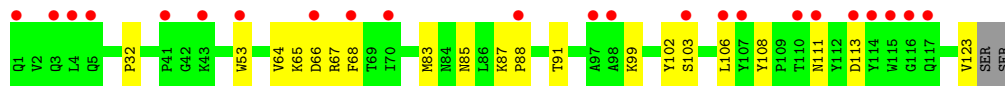
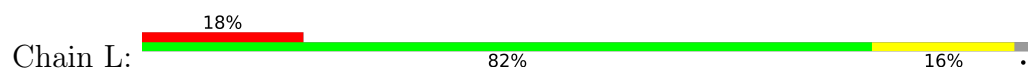




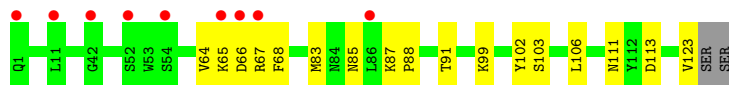
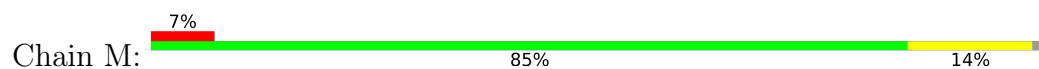
● Molecule 2: Nanobody Nb25



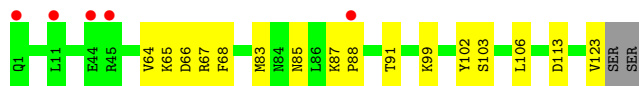
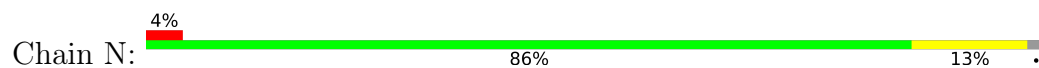
● Molecule 2: Nanobody Nb25



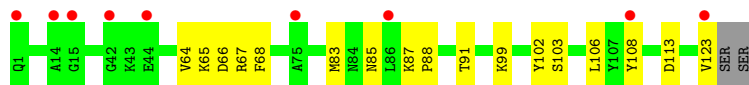
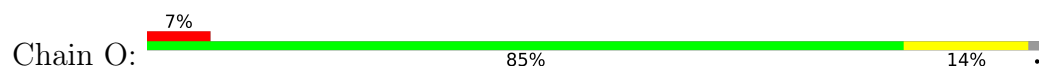
● Molecule 2: Nanobody Nb25



● Molecule 2: Nanobody Nb25

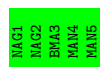


● Molecule 2: Nanobody Nb25



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





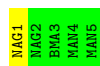
- Molecule 3: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G: 100%



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

Chain H: 80% 20%



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

Chain I: 100%



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

Chain J: 50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.35Å 139.94Å 191.50Å 90.00° 102.25° 90.00°	Depositor
Resolution (Å)	48.88 – 3.30 48.88 – 3.30	Depositor EDS
% Data completeness (in resolution range)	73.6 (48.88-3.30) 100.0 (48.88-3.30)	Depositor EDS
R_{merge}	0.35	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 3.25Å)	Xtriage
Refinement program	REFMAC 5.8.0158, PHENIX (1.12rc)-2787-000)	Depositor
R, R_{free}	0.234 , 0.250 0.251 , 0.268	Depositor DCC
R_{free} test set	3412 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	85.8	Xtriage
Anisotropy	0.424	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 11.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	18538	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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: NAG, MAN, BMA

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	0.68	0/2737	0.62	0/3724
1	B	0.67	1/2759 (0.0%)	0.61	0/3753
1	C	0.67	0/2737	0.62	0/3724
1	D	0.70	0/2751	0.61	0/3742
1	E	0.66	0/2737	0.61	0/3724
2	K	0.55	0/975	0.64	2/1321 (0.2%)
2	L	0.53	0/975	0.63	2/1321 (0.2%)
2	M	0.54	0/975	0.62	0/1321
2	N	0.55	0/975	0.62	0/1321
2	O	0.53	0/975	0.63	2/1321 (0.2%)
All	All	0.64	1/18596 (0.0%)	0.62	6/25272 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	254	SER	CA-C	-5.04	1.47	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	108	TYR	CA-C-N	5.55	125.08	119.19
2	K	108	TYR	C-N-CA	5.55	125.08	119.19
2	L	108	TYR	CA-C-N	5.30	124.81	119.19
2	L	108	TYR	C-N-CA	5.30	124.81	119.19
2	O	108	TYR	CA-C-N	5.18	124.68	119.19
2	O	108	TYR	C-N-CA	5.18	124.68	119.19

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	2670	0	2678	62	0
1	B	2692	0	2698	73	0
1	C	2670	0	2678	69	0
1	D	2684	0	2692	70	0
1	E	2670	0	2678	65	0
2	K	951	0	895	15	0
2	L	951	0	895	16	0
2	M	951	0	895	16	0
2	N	951	0	895	14	0
2	O	951	0	895	14	0
3	F	61	0	52	0	0
3	G	61	0	52	0	0
3	H	61	0	52	1	0
4	I	72	0	61	0	0
4	J	72	0	61	2	0
5	A	14	0	13	0	0
5	B	14	0	13	0	0
5	C	14	0	13	0	0
5	D	14	0	13	0	0
5	E	14	0	13	0	0
All	All	18538	0	18242	366	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 10.

All (366) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:TYR:OH	1:E:99:LEU:CD1	2.22	0.87
1:D:207:ARG:NH1	2:N:102:TYR:O	2.07	0.87
1:A:99:LEU:CD1	1:B:62:TYR:OH	2.24	0.86
1:B:143:TYR:OH	1:B:229:PHE:HE2	1.57	0.85
1:C:196:ARG:NH1	2:M:113:ASP:OD1	2.10	0.83
1:B:196:ARG:NH1	2:L:113:ASP:OD1	2.10	0.83
1:A:99:LEU:HD12	1:B:62:TYR:OH	1.79	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:99:LEU:CD1	1:E:62:TYR:OH	2.28	0.81
1:B:99:LEU:CD1	1:C:62:TYR:OH	2.28	0.81
1:C:99:LEU:CD1	1:D:62:TYR:OH	2.31	0.79
1:A:207:ARG:NH1	2:K:102:TYR:O	2.15	0.79
2:O:64:VAL:HG13	2:O:68:PHE:CD1	2.18	0.79
1:B:207:ARG:NH1	2:L:102:TYR:O	2.15	0.78
2:L:64:VAL:HG13	2:L:68:PHE:CD1	2.18	0.78
2:M:64:VAL:HG13	2:M:68:PHE:CD1	2.18	0.78
1:E:196:ARG:NH1	2:O:113:ASP:OD1	2.15	0.78
2:N:64:VAL:HG13	2:N:68:PHE:CD1	2.18	0.78
1:E:99:LEU:HD22	1:E:207:ARG:NH2	1.99	0.78
2:K:64:VAL:HG13	2:K:68:PHE:CD1	2.18	0.77
1:B:99:LEU:HD22	1:B:207:ARG:NH2	1.99	0.77
1:C:99:LEU:HD22	1:C:207:ARG:NH2	2.00	0.77
1:C:207:ARG:NH1	2:M:102:TYR:O	2.19	0.76
1:A:99:LEU:HD22	1:A:207:ARG:NH2	2.00	0.76
1:E:207:ARG:NH1	2:O:102:TYR:O	2.18	0.76
1:D:49:MET:HB3	1:D:58:THR:HG22	1.68	0.76
1:B:49:MET:HB3	1:B:58:THR:HG22	1.68	0.76
1:A:62:TYR:OH	1:E:99:LEU:HD12	1.87	0.75
1:D:99:LEU:HD22	1:D:207:ARG:NH2	2.02	0.75
1:D:99:LEU:HD12	1:E:62:TYR:OH	1.88	0.73
1:C:49:MET:HB3	1:C:58:THR:HG22	1.71	0.72
1:E:49:MET:HB3	1:E:58:THR:HG22	1.71	0.72
1:E:192:ARG:NH1	4:J:1:NAG:O7	2.25	0.70
1:D:196:ARG:NH1	2:N:113:ASP:OD1	2.23	0.70
1:A:49:MET:HB3	1:A:58:THR:HG22	1.71	0.70
1:B:99:LEU:HD12	1:C:62:TYR:OH	1.92	0.70
1:B:145:LEU:CD2	1:B:226:ILE:CD1	2.72	0.67
1:E:60:THR:OG1	1:E:131:THR:HG22	1.95	0.67
1:A:145:LEU:CD2	1:A:226:ILE:CD1	2.72	0.67
1:D:145:LEU:CD2	1:D:226:ILE:CD1	2.73	0.66
1:A:60:THR:OG1	1:A:131:THR:HG22	1.95	0.66
1:A:248:PHE:O	1:A:399:ARG:NE	2.28	0.66
1:C:145:LEU:CD2	1:C:226:ILE:CD1	2.73	0.66
1:C:99:LEU:HD12	1:D:62:TYR:OH	1.94	0.66
1:C:60:THR:OG1	1:C:131:THR:HG22	1.96	0.66
1:E:248:PHE:O	1:E:399:ARG:NE	2.29	0.66
1:C:248:PHE:O	1:C:399:ARG:NE	2.29	0.65
1:D:19:LEU:HD22	1:D:72:LEU:HA	1.79	0.65
1:B:60:THR:OG1	1:B:131:THR:HG22	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:153:GLU:OE1	1:C:209:SER:OG	2.14	0.65
1:D:248:PHE:O	1:D:399:ARG:NE	2.30	0.65
1:D:60:THR:OG1	1:D:131:THR:HG22	1.96	0.65
1:D:58:THR:HB	1:D:133:THR:HG22	1.79	0.64
1:E:145:LEU:CD2	1:E:226:ILE:CD1	2.74	0.64
1:D:40:MET:HE1	1:D:96:THR:HG21	1.80	0.64
2:O:91:THR:HG22	2:O:123:VAL:H	1.63	0.64
1:C:58:THR:CB	1:C:133:THR:HG22	2.29	0.63
1:D:40:MET:HG3	1:D:208:LEU:HD12	1.80	0.63
2:K:91:THR:HG22	2:K:123:VAL:H	1.63	0.63
1:C:58:THR:HB	1:C:133:THR:HG22	1.80	0.63
2:L:91:THR:HG22	2:L:123:VAL:H	1.63	0.63
2:M:64:VAL:HG13	2:M:68:PHE:CG	2.33	0.63
1:E:40:MET:HG3	1:E:208:LEU:HD12	1.81	0.63
1:D:58:THR:CB	1:D:133:THR:HG22	2.29	0.62
1:A:58:THR:CB	1:A:133:THR:HG22	2.30	0.62
1:A:58:THR:HB	1:A:133:THR:HG22	1.80	0.62
1:B:143:TYR:OH	1:B:229:PHE:CE2	2.41	0.62
2:M:91:THR:HG22	2:M:123:VAL:H	1.65	0.62
2:O:64:VAL:HG13	2:O:68:PHE:CG	2.34	0.62
2:K:64:VAL:HG13	2:K:68:PHE:CG	2.34	0.61
1:B:58:THR:HB	1:B:133:THR:HG22	1.83	0.61
2:L:64:VAL:HG13	2:L:68:PHE:CG	2.36	0.61
1:B:49:MET:HB3	1:B:58:THR:CG2	2.31	0.60
2:N:64:VAL:HG13	2:N:68:PHE:CG	2.35	0.60
2:N:91:THR:HG22	2:N:123:VAL:H	1.65	0.60
1:E:58:THR:CB	1:E:133:THR:HG22	2.32	0.60
1:A:164:ILE:HG23	1:A:206:PRO:HG3	1.81	0.60
1:D:68:ARG:NH1	1:D:121:ASP:OD1	2.34	0.60
1:E:68:ARG:NH1	1:E:121:ASP:OD1	2.35	0.60
1:C:68:ARG:NH1	1:C:121:ASP:OD1	2.34	0.60
1:B:68:ARG:NH1	1:B:121:ASP:OD1	2.34	0.59
1:A:68:ARG:NH1	1:A:121:ASP:OD1	2.35	0.59
1:C:145:LEU:CD2	1:C:226:ILE:HD13	2.31	0.59
1:A:62:TYR:OH	1:E:99:LEU:HD11	2.03	0.59
1:B:40:MET:HG3	1:B:208:LEU:HD12	1.85	0.59
1:C:232:GLN:O	1:C:236:PRO:HG2	2.03	0.59
1:D:49:MET:HB3	1:D:58:THR:CG2	2.33	0.58
1:B:58:THR:CB	1:B:133:THR:HG22	2.32	0.58
1:B:19:LEU:HD22	1:B:72:LEU:HA	1.84	0.58
1:C:40:MET:HG3	1:C:208:LEU:HD12	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:58:THR:HB	1:E:133:THR:HG22	1.84	0.58
1:E:49:MET:HB3	1:E:58:THR:CG2	2.34	0.58
1:E:99:LEU:HD22	1:E:207:ARG:CZ	2.33	0.58
1:E:145:LEU:CD2	1:E:226:ILE:HD13	2.33	0.58
1:C:99:LEU:HD22	1:C:207:ARG:CZ	2.33	0.58
1:A:196:ARG:NH1	2:K:113:ASP:OD1	2.37	0.57
2:O:67:ARG:NH2	2:O:85:ASN:O	2.37	0.57
1:C:255:VAL:N	1:C:256:PRO:HD2	2.19	0.57
1:B:255:VAL:N	1:B:256:PRO:HD2	2.19	0.57
1:E:255:VAL:N	1:E:256:PRO:HD2	2.19	0.57
1:C:19:LEU:HD22	1:C:72:LEU:HA	1.86	0.57
1:B:145:LEU:HD23	1:B:226:ILE:CD1	2.35	0.57
1:B:137:MET:HE2	1:C:182:GLU:HB3	1.86	0.57
1:D:255:VAL:N	1:D:256:PRO:HD2	2.19	0.57
1:E:19:LEU:HD22	1:E:72:LEU:HA	1.86	0.57
1:A:255:VAL:N	1:A:256:PRO:HD2	2.19	0.57
2:L:67:ARG:NH2	2:L:85:ASN:O	2.38	0.57
1:A:99:LEU:HD22	1:A:207:ARG:CZ	2.34	0.57
1:D:88:ALA:HB2	1:D:116:ILE:HD11	1.87	0.57
1:D:99:LEU:HD22	1:D:207:ARG:CZ	2.35	0.57
1:D:145:LEU:CD2	1:D:226:ILE:HD13	2.35	0.57
1:C:288:ALA:HB2	1:C:417:TYR:CE2	2.40	0.56
1:A:40:MET:HG3	1:A:208:LEU:HD12	1.86	0.56
1:E:235:LEU:HD23	1:E:238:ILE:HD11	1.86	0.56
2:M:67:ARG:NH2	2:M:85:ASN:O	2.38	0.56
2:N:67:ARG:NH2	2:N:85:ASN:O	2.38	0.56
2:K:67:ARG:NH2	2:K:85:ASN:O	2.38	0.56
1:E:143:TYR:HH	1:E:229:PHE:HE1	1.55	0.55
1:E:226:ILE:HG22	1:E:226:ILE:O	2.06	0.55
1:C:226:ILE:O	1:C:226:ILE:HG22	2.06	0.55
1:B:277:ARG:HB3	1:C:232:GLN:HE21	1.71	0.55
1:B:145:LEU:CD2	1:B:226:ILE:HD13	2.37	0.54
1:B:145:LEU:CD2	1:B:226:ILE:HD11	2.37	0.54
1:D:68:ARG:NH1	1:D:121:ASP:O	2.37	0.54
1:A:145:LEU:HD23	1:A:226:ILE:CD1	2.37	0.54
1:A:226:ILE:O	1:A:226:ILE:HG22	2.07	0.54
2:N:65:LYS:O	2:N:66:ASP:HB2	2.08	0.54
1:A:248:PHE:O	1:A:399:ARG:NH2	2.40	0.54
1:C:145:LEU:HD23	1:C:226:ILE:CD1	2.38	0.54
1:A:49:MET:HB3	1:A:58:THR:CG2	2.36	0.54
1:D:143:TYR:HH	1:D:229:PHE:HE1	1.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:248:PHE:O	1:C:399:ARG:NH2	2.41	0.54
2:M:65:LYS:O	2:M:66:ASP:HB2	2.07	0.54
1:B:115:MET:HE3	1:B:125:LEU:HG	1.91	0.53
1:A:145:LEU:CD2	1:A:226:ILE:HD13	2.38	0.53
1:B:226:ILE:HG22	1:B:226:ILE:O	2.07	0.53
1:C:277:ARG:HB3	1:D:232:GLN:HE21	1.73	0.53
1:B:99:LEU:HD22	1:B:207:ARG:CZ	2.38	0.53
2:N:68:PHE:HE2	2:N:83:MET:SD	2.31	0.53
1:B:103:LYS:HE3	1:C:105:PHE:CE2	2.44	0.53
2:K:68:PHE:CE2	2:K:83:MET:SD	3.02	0.53
1:A:62:TYR:HH	1:E:99:LEU:HD12	1.74	0.53
1:C:49:MET:HB3	1:C:58:THR:CG2	2.36	0.53
1:D:226:ILE:HG22	1:D:226:ILE:O	2.07	0.53
1:A:115:MET:HE3	1:A:125:LEU:HG	1.91	0.53
1:D:248:PHE:O	1:D:399:ARG:NH2	2.41	0.53
2:O:68:PHE:HE2	2:O:83:MET:SD	2.32	0.53
2:L:68:PHE:CE2	2:L:83:MET:SD	3.03	0.53
2:N:68:PHE:CE2	2:N:83:MET:SD	3.03	0.53
2:K:68:PHE:HE2	2:K:83:MET:SD	2.31	0.52
2:L:68:PHE:HE2	2:L:83:MET:SD	2.31	0.52
1:C:230:VAL:HA	1:C:234:TYR:HB2	1.91	0.52
2:K:65:LYS:O	2:K:66:ASP:HB2	2.07	0.52
2:M:68:PHE:HE2	2:M:83:MET:SD	2.32	0.52
2:O:68:PHE:CE2	2:O:83:MET:SD	3.03	0.52
1:A:145:LEU:CD2	1:A:226:ILE:HD11	2.37	0.52
1:B:143:TYR:HH	1:B:229:PHE:HE2	0.74	0.52
2:M:68:PHE:CE2	2:M:83:MET:SD	3.03	0.52
2:O:65:LYS:O	2:O:66:ASP:HB2	2.08	0.52
2:L:65:LYS:O	2:L:66:ASP:HB2	2.09	0.52
1:E:48:ASP:HB2	1:E:58:THR:HG23	1.92	0.52
1:B:277:ARG:HB3	1:C:232:GLN:NE2	2.24	0.52
1:C:235:LEU:HD23	1:C:238:ILE:HD11	1.91	0.51
1:D:48:ASP:HB2	1:D:58:THR:CG2	2.40	0.51
1:D:145:LEU:HD23	1:D:226:ILE:CD1	2.40	0.51
1:E:248:PHE:O	1:E:399:ARG:NH2	2.42	0.51
2:O:91:THR:HG22	2:O:123:VAL:N	2.25	0.51
1:C:48:ASP:HB2	1:C:58:THR:CG2	2.41	0.51
1:E:68:ARG:NH1	1:E:121:ASP:O	2.38	0.51
2:L:91:THR:HG22	2:L:123:VAL:N	2.25	0.51
1:B:48:ASP:HB2	1:B:58:THR:HG23	1.93	0.51
1:B:68:ARG:NH1	1:B:121:ASP:O	2.38	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:91:THR:HG22	2:M:123:VAL:N	2.26	0.51
2:K:91:THR:HG22	2:K:123:VAL:N	2.26	0.51
1:B:48:ASP:HB2	1:B:58:THR:CG2	2.41	0.51
1:E:114:ARG:HB2	1:E:128:LEU:HD23	1.93	0.51
1:D:48:ASP:HB2	1:D:58:THR:HG23	1.92	0.50
1:D:230:VAL:HA	1:D:234:TYR:HB2	1.94	0.50
1:A:230:VAL:HA	1:A:234:TYR:HB2	1.92	0.50
1:E:145:LEU:HD23	1:E:226:ILE:CD1	2.40	0.50
1:A:48:ASP:HB2	1:A:58:THR:CG2	2.42	0.50
1:D:66:TYR:CZ	1:D:125:LEU:HD13	2.47	0.50
1:C:48:ASP:HB2	1:C:58:THR:HG23	1.93	0.50
1:E:115:MET:HE3	1:E:125:LEU:HG	1.92	0.50
1:A:277:ARG:HB3	1:B:232:GLN:HE21	1.76	0.49
1:D:96:THR:HG22	1:D:166:PHE:CZ	2.47	0.49
1:D:115:MET:HE3	1:D:125:LEU:HG	1.95	0.49
1:A:277:ARG:HB3	1:B:232:GLN:NE2	2.28	0.49
1:C:59:LEU:O	1:C:131:THR:HA	2.13	0.49
2:N:91:THR:HG22	2:N:123:VAL:N	2.26	0.49
1:D:226:ILE:HG22	1:D:230:VAL:HG23	1.94	0.49
1:B:230:VAL:HA	1:B:234:TYR:HB2	1.95	0.49
1:B:77:ILE:HG21	1:B:81:LEU:HD21	1.94	0.49
1:A:48:ASP:HB2	1:A:58:THR:HG23	1.93	0.49
1:E:48:ASP:HB2	1:E:58:THR:CG2	2.41	0.49
2:N:87:LYS:HB3	2:N:88:PRO:HD2	1.94	0.49
1:A:189:VAL:HG12	1:A:213:ARG:HB3	1.94	0.48
1:C:66:TYR:CZ	1:C:125:LEU:HD13	2.48	0.48
1:C:235:LEU:HA	1:C:238:ILE:HG12	1.94	0.48
1:D:59:LEU:O	1:D:131:THR:HA	2.13	0.48
2:K:87:LYS:HB3	2:K:88:PRO:HD2	1.94	0.48
1:C:251:ASN:OD1	1:C:253:GLU:HB2	2.13	0.48
1:B:59:LEU:O	1:B:131:THR:HA	2.14	0.48
1:C:115:MET:HE3	1:C:125:LEU:HG	1.96	0.48
1:E:59:LEU:O	1:E:131:THR:HA	2.14	0.48
1:E:230:VAL:HA	1:E:234:TYR:HB2	1.94	0.48
1:C:85:ASN:HB2	1:C:114:ARG:HG3	1.95	0.48
1:C:103:LYS:HE3	1:D:105:PHE:CE2	2.48	0.48
1:B:268:THR:O	1:B:272:LEU:HG	2.14	0.48
1:C:31:PHE:CE2	1:D:12:VAL:HG11	2.49	0.48
1:C:68:ARG:NH1	1:C:121:ASP:O	2.39	0.48
1:D:99:LEU:HD11	1:E:62:TYR:OH	2.14	0.48
2:O:87:LYS:HB3	2:O:88:PRO:HD2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:LEU:CD1	1:D:417:TYR:HB3	2.44	0.47
2:M:87:LYS:HB3	2:M:88:PRO:HD2	1.95	0.47
1:A:182:GLU:HB3	1:E:137:MET:HE2	1.96	0.47
1:D:145:LEU:CD2	1:D:226:ILE:HD11	2.43	0.47
2:L:87:LYS:HB3	2:L:88:PRO:HD2	1.96	0.47
1:A:59:LEU:O	1:A:131:THR:HA	2.14	0.47
1:A:12:VAL:HG11	1:E:31:PHE:CE2	2.50	0.47
1:C:145:LEU:CD2	1:C:226:ILE:HD11	2.45	0.47
1:C:268:THR:O	1:C:272:LEU:HG	2.15	0.46
1:D:145:LEU:HD23	1:D:226:ILE:HG12	1.97	0.46
1:D:251:ASN:OD1	1:D:253:GLU:HB2	2.15	0.46
1:C:288:ALA:CB	1:C:417:TYR:CE2	2.99	0.46
1:E:235:LEU:HA	1:E:238:ILE:HG12	1.96	0.46
1:B:36:VAL:HB	1:B:164:ILE:HD13	1.97	0.46
1:D:107:HIS:NE2	1:D:131:THR:HG23	2.30	0.46
1:B:79:LEU:HD12	1:B:81:LEU:HD23	1.97	0.46
1:B:143:TYR:CZ	1:B:229:PHE:CE2	3.03	0.46
1:B:145:LEU:CD1	1:B:417:TYR:HB3	2.46	0.46
1:A:66:TYR:CZ	1:A:125:LEU:HD13	2.51	0.46
1:A:31:PHE:CE2	1:B:81:LEU:HD22	2.51	0.46
1:A:143:TYR:HH	1:A:229:PHE:HE1	1.63	0.46
1:A:68:ARG:NH1	1:A:121:ASP:O	2.38	0.46
1:A:255:VAL:HB	1:B:254:SER:OG	2.15	0.46
1:B:107:HIS:NE2	1:B:131:THR:HG23	2.31	0.46
1:D:99:LEU:HD12	1:E:62:TYR:HH	1.77	0.45
1:A:145:LEU:CD1	1:A:417:TYR:HB3	2.46	0.45
1:E:66:TYR:CZ	1:E:125:LEU:HD13	2.51	0.45
2:K:99:LYS:HG2	2:K:103:SER:HB2	1.99	0.45
2:O:68:PHE:CE2	2:O:83:MET:HB3	2.52	0.45
1:A:85:ASN:HB2	1:A:114:ARG:CG	2.47	0.45
1:C:26:ARG:NH1	1:D:20:LEU:HD12	2.31	0.45
1:E:268:THR:O	1:E:272:LEU:HG	2.16	0.45
1:E:145:LEU:CD1	1:E:417:TYR:HB3	2.46	0.45
1:B:66:TYR:CZ	1:B:125:LEU:HD13	2.51	0.45
2:K:68:PHE:CE2	2:K:83:MET:HB3	2.51	0.45
4:J:3:BMA:H62	4:J:4:MAN:H2	1.77	0.45
1:D:85:ASN:HB2	1:D:114:ARG:CG	2.46	0.45
2:N:68:PHE:CE2	2:N:83:MET:HB3	2.51	0.45
1:E:145:LEU:HD23	1:E:226:ILE:HG12	1.98	0.45
1:C:95:ASP:OD1	1:C:156:SER:HB2	2.16	0.45
1:D:85:ASN:HB2	1:D:114:ARG:HG3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:68:PHE:CE2	2:M:83:MET:HB3	2.52	0.45
1:A:107:HIS:NE2	1:A:131:THR:HG23	2.31	0.44
1:D:192:ARG:HB2	1:D:211:SER:OG	2.17	0.44
1:E:226:ILE:HG22	1:E:230:VAL:HG23	2.00	0.44
1:B:144:PRO:HD3	1:B:289:MET:HB2	1.99	0.44
1:E:95:ASP:OD1	1:E:156:SER:HB2	2.17	0.44
1:E:262:GLY:HA3	1:E:304:LEU:HD13	2.00	0.44
1:B:66:TYR:CE2	1:B:125:LEU:HD13	2.53	0.44
1:B:95:ASP:OD1	1:B:156:SER:HB2	2.17	0.44
1:C:226:ILE:HG22	1:C:230:VAL:HG23	1.99	0.44
1:C:248:PHE:O	1:C:399:ARG:CZ	2.65	0.44
1:A:268:THR:O	1:A:272:LEU:HG	2.17	0.44
1:D:36:VAL:HB	1:D:164:ILE:HD13	1.98	0.44
2:M:99:LYS:HG2	2:M:103:SER:HB2	1.99	0.44
1:A:99:LEU:HD12	1:B:62:TYR:HH	1.77	0.44
1:A:248:PHE:O	1:A:399:ARG:CZ	2.66	0.44
1:B:226:ILE:HG22	1:B:230:VAL:HG23	2.00	0.44
1:D:66:TYR:CE2	1:D:125:LEU:HD13	2.52	0.44
1:D:95:ASP:OD1	1:D:156:SER:HB2	2.17	0.44
1:D:268:THR:O	1:D:272:LEU:HG	2.18	0.44
1:E:107:HIS:NE2	1:E:131:THR:HG23	2.33	0.44
2:O:99:LYS:HG2	2:O:103:SER:HB2	1.99	0.44
1:B:395:ASP:O	1:B:399:ARG:HG3	2.18	0.44
2:L:68:PHE:CE2	2:L:83:MET:HB3	2.52	0.44
1:A:58:THR:HG23	1:A:58:THR:O	2.18	0.44
1:A:145:LEU:HD23	1:A:226:ILE:HG12	1.99	0.44
1:D:31:PHE:CE2	1:E:12:VAL:HG11	2.53	0.44
2:L:99:LYS:HG2	2:L:103:SER:HB2	2.00	0.44
1:A:95:ASP:OD1	1:A:156:SER:HB2	2.17	0.43
1:A:235:LEU:HD23	1:A:235:LEU:HA	1.89	0.43
1:B:145:LEU:HD23	1:B:226:ILE:HD11	2.00	0.43
1:B:147:GLU:OE1	1:B:213:ARG:NH1	2.50	0.43
1:E:145:LEU:CD2	1:E:226:ILE:HD11	2.46	0.43
1:C:145:LEU:CD1	1:C:417:TYR:HB3	2.48	0.43
1:A:226:ILE:HG22	1:A:230:VAL:HG23	2.00	0.43
1:C:402:PHE:HB2	1:C:403:PRO:HD3	2.00	0.43
1:E:402:PHE:HB2	1:E:403:PRO:HD3	2.00	0.43
1:C:58:THR:HG23	1:C:58:THR:O	2.18	0.43
2:M:99:LYS:HE2	2:M:106:LEU:HD23	2.00	0.43
1:B:192:ARG:HB2	1:B:211:SER:OG	2.17	0.43
2:N:99:LYS:HG2	2:N:103:SER:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:LEU:CD1	1:B:81:LEU:HD23	2.49	0.43
1:B:101:ASP:HB2	1:B:132:THR:HG23	2.01	0.43
1:D:248:PHE:O	1:D:399:ARG:CZ	2.67	0.43
2:L:99:LYS:HE2	2:L:106:LEU:HD23	2.01	0.43
2:O:99:LYS:HE2	2:O:106:LEU:HD23	2.01	0.43
1:A:99:LEU:HD11	1:B:62:TYR:OH	2.15	0.43
1:B:85:ASN:HB2	1:B:114:ARG:HG3	2.00	0.43
1:A:262:GLY:HA3	1:A:304:LEU:HD13	2.01	0.43
1:B:199:VAL:HB	2:L:111:ASN:HD21	1.84	0.43
1:C:107:HIS:NE2	1:C:131:THR:HG23	2.33	0.43
1:E:66:TYR:CE2	1:E:125:LEU:HD13	2.54	0.43
1:A:192:ARG:HB2	1:A:211:SER:OG	2.19	0.42
1:D:144:PRO:HD3	1:D:289:MET:HB2	2.00	0.42
1:B:248:PHE:HB3	1:B:399:ARG:HG2	1.99	0.42
1:E:238:ILE:HG13	1:E:239:MET:N	2.33	0.42
1:A:101:ASP:HB2	1:A:132:THR:HG23	2.01	0.42
1:C:66:TYR:CE2	1:C:125:LEU:HD13	2.55	0.42
1:D:83:LEU:HD11	1:D:118:LEU:HD12	2.01	0.42
1:B:145:LEU:HD23	1:B:226:ILE:HG12	2.02	0.42
1:C:31:PHE:CE2	1:D:81:LEU:HD22	2.54	0.42
1:D:262:GLY:HA3	1:D:304:LEU:HD13	2.02	0.42
1:D:58:THR:HG23	1:D:58:THR:O	2.19	0.42
1:B:402:PHE:HB2	1:B:403:PRO:HD3	2.01	0.42
1:E:43:ASP:HA	1:E:176:THR:OG1	2.20	0.42
1:E:248:PHE:O	1:E:399:ARG:CZ	2.67	0.41
1:B:280:LEU:HB3	1:B:281:PRO:HD2	2.02	0.41
1:C:88:ALA:HB2	1:C:116:ILE:HD11	2.02	0.41
1:C:255:VAL:N	1:C:256:PRO:CD	2.83	0.41
1:D:145:LEU:HD22	1:D:226:ILE:HD11	2.02	0.41
1:A:66:TYR:CE2	1:A:125:LEU:HD13	2.54	0.41
1:C:192:ARG:HB2	1:C:211:SER:OG	2.20	0.41
1:A:255:VAL:N	1:A:256:PRO:CD	2.83	0.41
1:C:254:SER:HB3	1:C:257:ALA:HB3	2.02	0.41
1:E:144:PRO:HD3	1:E:289:MET:HB2	2.03	0.41
1:B:85:ASN:HB2	1:B:114:ARG:CG	2.49	0.41
1:E:101:ASP:HB2	1:E:132:THR:HG23	2.02	0.41
1:E:189:VAL:CG1	1:E:213:ARG:HB3	2.50	0.41
1:C:36:VAL:HB	1:C:164:ILE:HD13	2.02	0.41
1:D:405:LEU:HD23	1:D:405:LEU:HA	1.90	0.41
1:A:267:LEU:HD22	1:B:268:THR:HG21	2.01	0.41
1:B:289:MET:HE2	1:B:289:MET:HB3	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:145:LEU:HD23	1:C:226:ILE:HG12	2.01	0.41
1:D:26:ARG:NH2	1:E:17:ASP:OD1	2.48	0.41
2:K:108:TYR:HA	2:K:109:PRO:HD3	1.86	0.41
1:C:85:ASN:HB2	1:C:114:ARG:CG	2.50	0.41
1:E:58:THR:HG23	1:E:58:THR:O	2.20	0.41
1:A:43:ASP:HA	1:A:176:THR:OG1	2.21	0.41
1:A:268:THR:HG21	1:E:267:LEU:HD22	2.03	0.41
1:B:58:THR:HG23	1:B:58:THR:O	2.19	0.41
1:B:255:VAL:N	1:B:256:PRO:CD	2.84	0.41
1:D:145:LEU:HD12	1:D:417:TYR:HB3	2.03	0.41
1:D:179:GLU:CD	2:M:106:LEU:HD11	2.46	0.41
1:D:402:PHE:HB2	1:D:403:PRO:HD3	2.02	0.41
1:E:395:ASP:O	1:E:399:ARG:HG3	2.21	0.41
2:L:32:PRO:HB2	2:L:53:TRP:HB2	2.03	0.41
1:B:262:GLY:HA3	1:B:304:LEU:HD13	2.03	0.41
1:C:202:THR:OG1	1:C:205:TYR:OH	2.20	0.41
1:D:254:SER:HB3	1:D:257:ALA:HB3	2.02	0.41
1:D:409:PHE:O	1:D:412:VAL:HG22	2.20	0.41
1:E:235:LEU:HD23	1:E:235:LEU:HA	1.91	0.41
2:N:99:LYS:HE2	2:N:106:LEU:HD23	2.03	0.41
1:B:145:LEU:HD13	1:B:417:TYR:HB3	2.03	0.40
1:B:99:LEU:HD11	1:C:62:TYR:OH	2.15	0.40
1:D:43:ASP:HA	1:D:176:THR:OG1	2.21	0.40
1:D:295:VAL:HG11	1:D:409:PHE:CZ	2.57	0.40
1:A:105:PHE:CE2	1:E:103:LYS:HE3	2.56	0.40
1:C:43:ASP:HA	1:C:176:THR:OG1	2.21	0.40
1:C:238:ILE:HG13	1:C:239:MET:N	2.36	0.40
2:M:103:SER:O	2:M:111:ASN:HB3	2.22	0.40
1:C:213:ARG:HB2	3:H:1:NAG:H81	2.02	0.40
1:E:409:PHE:O	1:E:412:VAL:HG22	2.21	0.40
2:K:32:PRO:HB2	2:K:53:TRP:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	328/395 (83%)	319 (97%)	9 (3%)	0	100	100
1	B	331/395 (84%)	322 (97%)	9 (3%)	0	100	100
1	C	328/395 (83%)	320 (98%)	8 (2%)	0	100	100
1	D	330/395 (84%)	320 (97%)	10 (3%)	0	100	100
1	E	328/395 (83%)	318 (97%)	10 (3%)	0	100	100
2	K	121/125 (97%)	115 (95%)	6 (5%)	0	100	100
2	L	121/125 (97%)	115 (95%)	6 (5%)	0	100	100
2	M	121/125 (97%)	116 (96%)	5 (4%)	0	100	100
2	N	121/125 (97%)	116 (96%)	5 (4%)	0	100	100
2	O	121/125 (97%)	116 (96%)	5 (4%)	0	100	100
All	All	2250/2600 (86%)	2177 (97%)	73 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	293/346 (85%)	291 (99%)	2 (1%)	76	80
1	B	296/346 (86%)	295 (100%)	1 (0%)	86	86
1	C	293/346 (85%)	291 (99%)	2 (1%)	76	80
1	D	295/346 (85%)	293 (99%)	2 (1%)	76	80
1	E	293/346 (85%)	291 (99%)	2 (1%)	76	80
2	K	97/99 (98%)	97 (100%)	0	100	100
2	L	97/99 (98%)	97 (100%)	0	100	100
2	M	97/99 (98%)	97 (100%)	0	100	100
2	N	97/99 (98%)	97 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	O	97/99 (98%)	97 (100%)	0	100	100
All	All	1955/2225 (88%)	1946 (100%)	9 (0%)	81	83

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	CYS
1	A	254	SER
1	B	37	CYS
1	C	37	CYS
1	C	254	SER
1	D	37	CYS
1	D	254	SER
1	E	37	CYS
1	E	254	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	191	HIS
1	A	245	GLN
1	A	410	ASN
1	B	54	ASN
1	B	85	ASN
1	B	185	GLN
1	B	191	HIS
1	B	245	GLN
1	B	410	ASN
1	C	54	ASN
1	C	85	ASN
1	C	191	HIS
1	C	245	GLN
1	C	410	ASN
1	D	54	ASN
1	D	85	ASN
1	D	191	HIS
1	D	245	GLN
1	D	410	ASN
1	E	54	ASN
1	E	85	ASN

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Mol	Chain	Res	Type
1	E	191	HIS
1	E	410	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 ⓘ

27 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	F	1	3,1	14,14,15	0.27	0	17,19,21	0.48	0
3	NAG	F	2	3	14,14,15	0.20	0	17,19,21	0.50	0
3	BMA	F	3	3	11,11,12	0.19	0	15,15,17	0.37	0
3	MAN	F	4	3	11,11,12	0.27	0	15,15,17	0.32	0
3	MAN	F	5	3	11,11,12	0.25	0	15,15,17	0.33	0
3	NAG	G	1	3,1	14,14,15	0.29	0	17,19,21	0.62	0
3	NAG	G	2	3	14,14,15	0.19	0	17,19,21	0.52	0
3	BMA	G	3	3	11,11,12	0.39	0	15,15,17	0.71	0
3	MAN	G	4	3	11,11,12	0.28	0	15,15,17	0.31	0
3	MAN	G	5	3	11,11,12	0.35	0	15,15,17	0.30	0
3	NAG	H	1	3,1	14,14,15	0.21	0	17,19,21	0.54	0
3	NAG	H	2	3	14,14,15	0.24	0	17,19,21	0.48	0
3	BMA	H	3	3	11,11,12	0.36	0	15,15,17	0.55	0
3	MAN	H	4	3	11,11,12	0.17	0	15,15,17	0.26	0
3	MAN	H	5	3	11,11,12	0.24	0	15,15,17	0.50	0
4	NAG	I	1	4,1	14,14,15	0.26	0	17,19,21	0.53	0
4	NAG	I	2	4	14,14,15	0.29	0	17,19,21	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	BMA	I	3	4	11,11,12	0.25	0	15,15,17	0.32	0
4	MAN	I	4	4	11,11,12	0.43	0	15,15,17	0.33	0
4	MAN	I	5	4	11,11,12	0.50	0	15,15,17	0.55	0
4	MAN	I	6	4	11,11,12	0.25	0	15,15,17	0.26	0
4	NAG	J	1	4,1	14,14,15	0.24	0	17,19,21	0.54	0
4	NAG	J	2	4	14,14,15	0.24	0	17,19,21	0.40	0
4	BMA	J	3	4	11,11,12	0.32	0	15,15,17	0.35	0
4	MAN	J	4	4	11,11,12	0.32	0	15,15,17	0.44	0
4	MAN	J	5	4	11,11,12	0.30	0	15,15,17	0.34	0
4	MAN	J	6	4	11,11,12	0.43	0	15,15,17	0.45	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	F	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	BMA	F	3	3	-	1/2/19/22	0/1/1/1
3	MAN	F	4	3	-	0/2/19/22	0/1/1/1
3	MAN	F	5	3	-	0/2/19/22	0/1/1/1
3	NAG	G	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	BMA	G	3	3	-	0/2/19/22	0/1/1/1
3	MAN	G	4	3	-	0/2/19/22	0/1/1/1
3	MAN	G	5	3	-	0/2/19/22	0/1/1/1
3	NAG	H	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
3	BMA	H	3	3	-	0/2/19/22	0/1/1/1
3	MAN	H	4	3	-	0/2/19/22	0/1/1/1
3	MAN	H	5	3	-	0/2/19/22	0/1/1/1
4	NAG	I	1	4,1	-	1/6/23/26	0/1/1/1
4	NAG	I	2	4	-	0/6/23/26	0/1/1/1
4	BMA	I	3	4	-	0/2/19/22	0/1/1/1
4	MAN	I	4	4	-	0/2/19/22	0/1/1/1
4	MAN	I	5	4	-	0/2/19/22	0/1/1/1
4	MAN	I	6	4	-	0/2/19/22	0/1/1/1
4	NAG	J	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1
4	BMA	J	3	4	-	1/2/19/22	0/1/1/1
4	MAN	J	4	4	-	0/2/19/22	0/1/1/1
4	MAN	J	5	4	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MAN	J	6	4	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

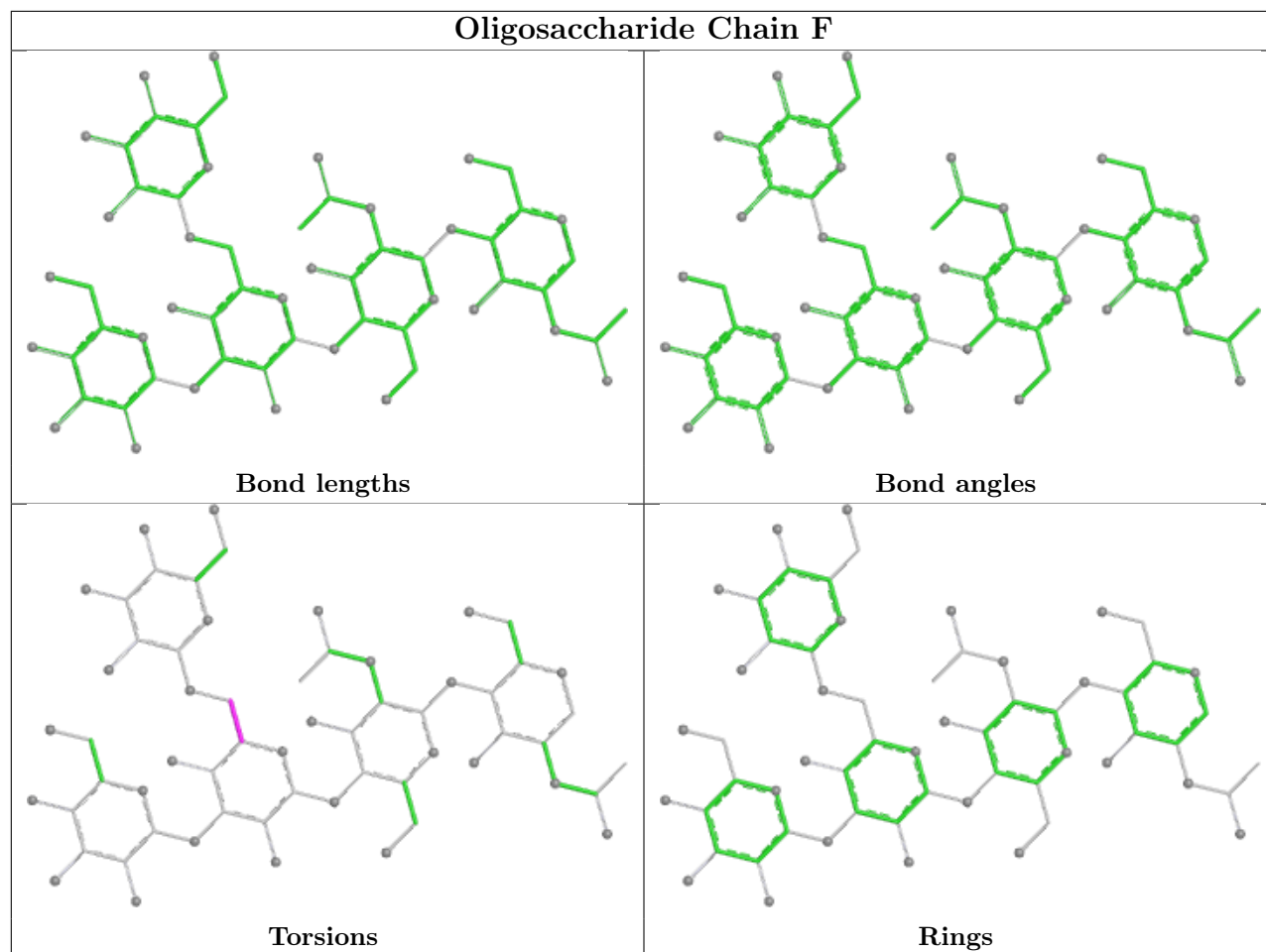
Mol	Chain	Res	Type	Atoms
4	J	3	BMA	O5-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6
3	F	3	BMA	C4-C5-C6-O6

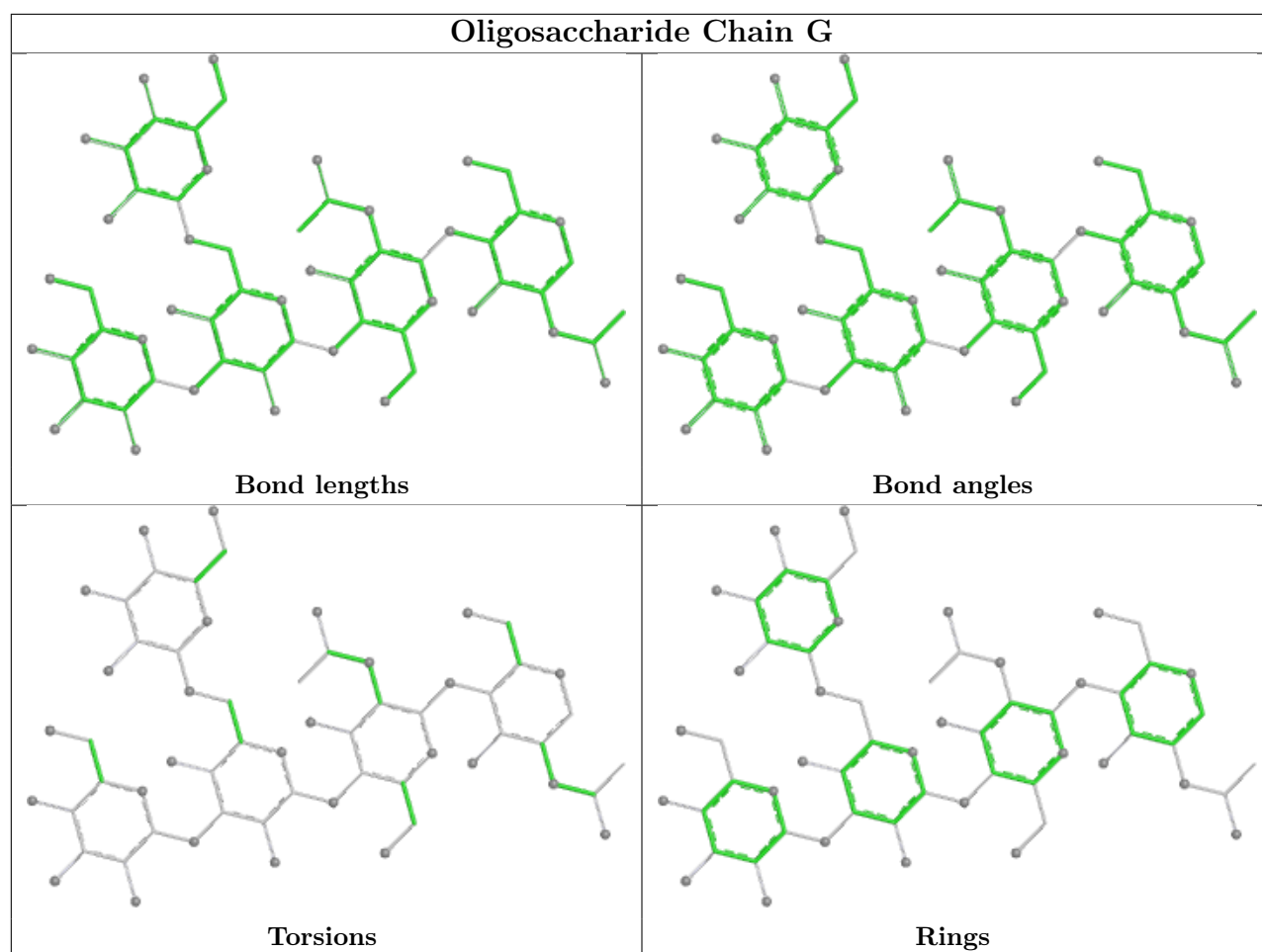
There are no ring outliers.

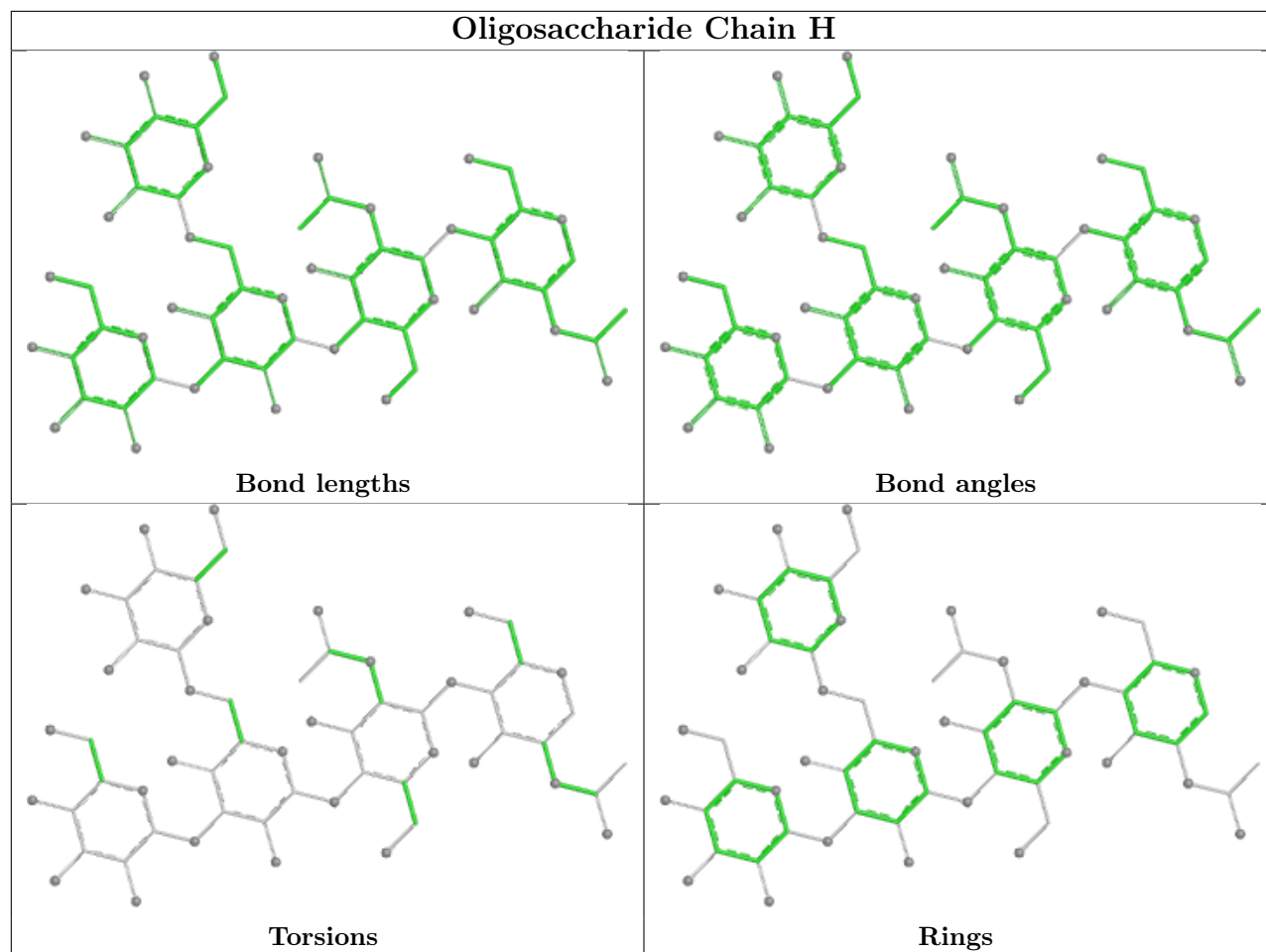
4 monomers are involved in 3 short contacts:

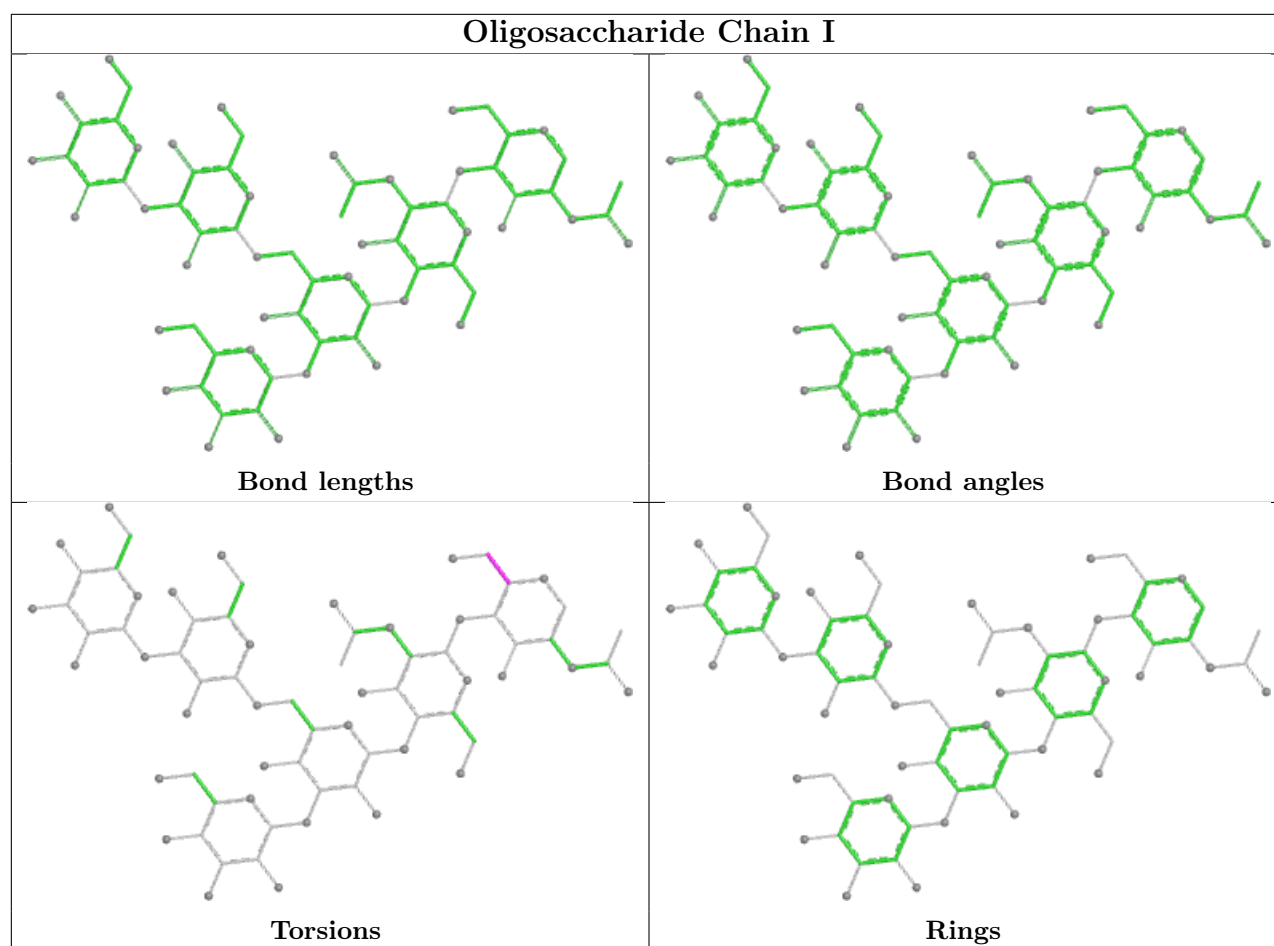
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	4	MAN	1	0
4	J	3	BMA	1	0
4	J	1	NAG	1	0
3	H	1	NAG	1	0

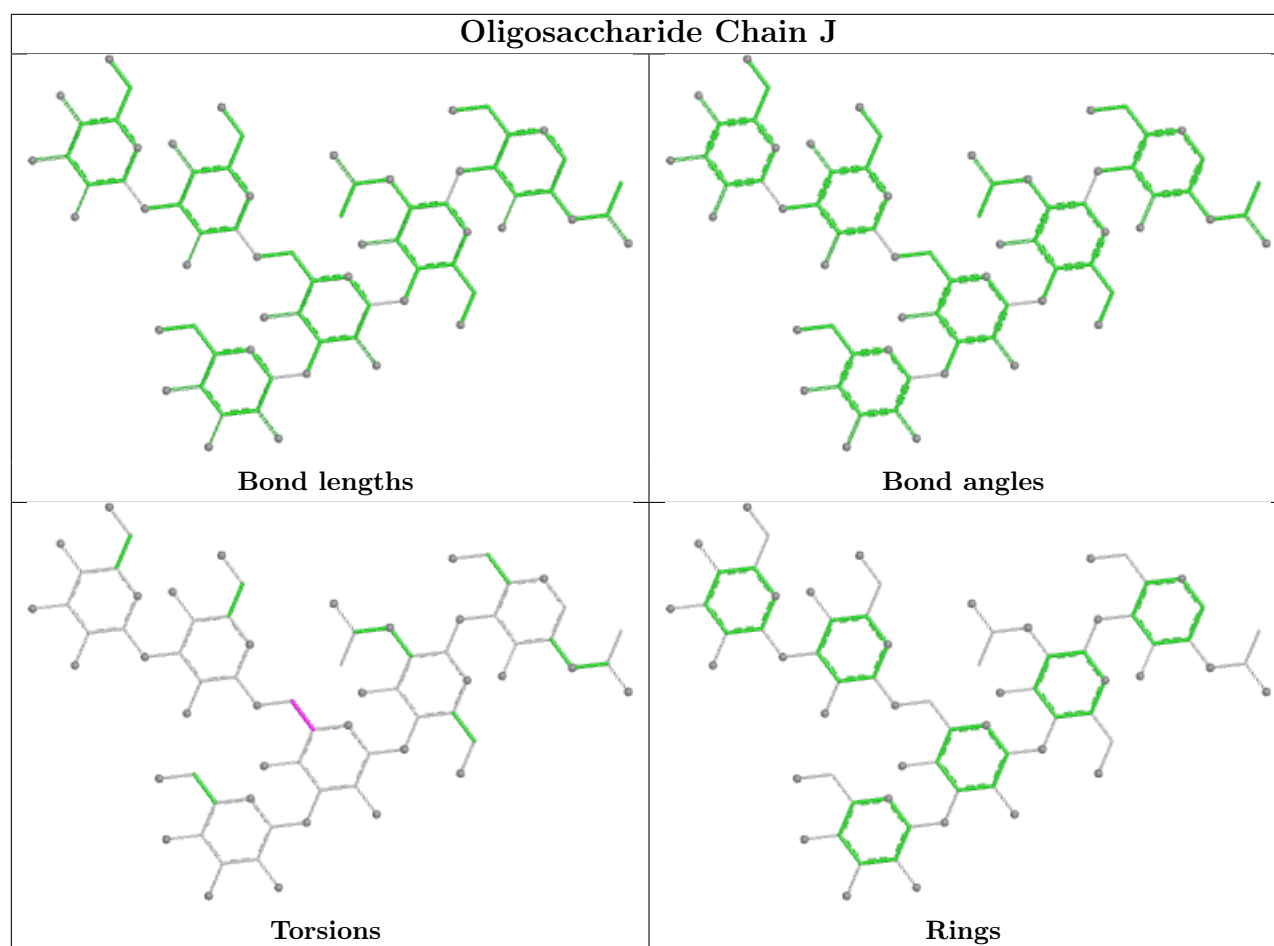
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	E	501	1	14,14,15	0.23	0	17,19,21	0.65	0
5	NAG	B	501	1	14,14,15	0.27	0	17,19,21	0.51	0
5	NAG	A	501	1	14,14,15	0.36	0	17,19,21	0.43	0
5	NAG	C	501	1	14,14,15	0.27	0	17,19,21	0.49	0
5	NAG	D	501	1	14,14,15	0.36	0	17,19,21	0.51	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
5	NAG	E	501	1	-	2/6/23/26	0/1/1/1
5	NAG	B	501	1	-	2/6/23/26	0/1/1/1
5	NAG	A	501	1	-	0/6/23/26	0/1/1/1
5	NAG	C	501	1	-	1/6/23/26	0/1/1/1
5	NAG	D	501	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 (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	501	NAG	C4-C5-C6-O6
5	E	501	NAG	O5-C5-C6-O6
5	D	501	NAG	O5-C5-C6-O6
5	D	501	NAG	C4-C5-C6-O6
5	B	501	NAG	C4-C5-C6-O6
5	B	501	NAG	O5-C5-C6-O6
5	C	501	NAG	C4-C5-C6-O6

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 [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	330/395 (83%)	0.41	29 (8%)	15 12	30, 58, 101, 133	0
1	B	333/395 (84%)	0.34	25 (7%)	20 15	31, 62, 94, 138	0
1	C	330/395 (83%)	0.45	34 (10%)	12 10	31, 61, 104, 137	0
1	D	332/395 (84%)	0.38	30 (9%)	15 11	28, 51, 85, 155	0
1	E	330/395 (83%)	0.33	25 (7%)	20 14	26, 54, 89, 141	0
2	K	123/125 (98%)	0.64	10 (8%)	18 13	54, 101, 132, 151	0
2	L	123/125 (98%)	1.10	23 (18%)	3 3	78, 125, 168, 182	0
2	M	123/125 (98%)	0.41	9 (7%)	21 15	58, 97, 129, 150	0
2	N	123/125 (98%)	0.19	5 (4%)	41 27	52, 102, 141, 153	0
2	O	123/125 (98%)	0.54	9 (7%)	21 15	60, 91, 117, 123	0
All	All	2270/2600 (87%)	0.43	199 (8%)	15 12	26, 65, 129, 182	0

All (199) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	278	ASN	9.9
2	L	117	GLN	9.7
1	C	278	ASN	9.5
1	A	127	GLY	8.2
2	O	44	GLU	7.9
2	L	115	TRP	7.9
1	C	232	GLN	7.7
2	L	1	GLN	6.0
1	A	278	ASN	6.0
2	L	116	GLY	5.9
2	O	14	ALA	5.8
2	M	66	ASP	5.6
2	N	88	PRO	5.2

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Mol	Chain	Res	Type	RSRZ
1	E	204	ALA	5.1
1	A	66	TYR	5.0
1	A	126	TYR	4.8
1	A	97	TYR	4.7
2	O	42	GLY	4.6
1	A	64	GLN	4.6
2	M	1	GLN	4.5
1	D	417	TYR	4.5
1	D	24	ASP	4.4
2	K	115	TRP	4.4
2	K	113	ASP	4.3
1	A	275	SER	4.2
1	A	125	LEU	4.1
1	A	115	MET	4.1
1	A	65	GLN	4.1
1	E	13	LYS	4.1
1	C	275	SER	4.0
1	E	253	GLU	4.0
1	C	170	GLY	3.9
1	C	215	LYS	3.9
1	A	157	TYR	3.9
2	L	5	GLN	3.9
1	B	45	ALA	3.8
1	D	142	ARG	3.7
1	B	417	TYR	3.7
1	A	159	TYR	3.7
1	C	85	ASN	3.7
1	B	273	SER	3.7
1	A	113	ASN	3.6
1	E	319	ALA	3.6
2	L	98	ALA	3.5
2	L	4	LEU	3.5
1	E	161	THR	3.5
1	E	396	LYS	3.5
2	L	114	TYR	3.5
1	C	417	TYR	3.5
1	D	23	TYR	3.5
1	C	89	ASP	3.5
1	D	27	LEU	3.5
1	C	233	THR	3.5
2	K	88	PRO	3.4
1	D	228	TYR	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	44	ILE	3.4
2	L	97	ALA	3.4
2	K	111	ASN	3.4
1	D	172	ASP	3.4
1	B	113	ASN	3.4
2	O	75	ALA	3.4
1	B	275	SER	3.3
1	C	279	SER	3.3
2	O	108	TYR	3.3
1	D	232	GLN	3.2
1	E	165	GLU	3.2
1	E	417	TYR	3.2
1	A	112	LYS	3.2
1	B	157	TYR	3.2
1	C	173	LYS	3.2
1	D	303	ALA	3.2
1	E	231	ILE	3.2
1	C	228	TYR	3.2
1	C	97	TYR	3.1
1	C	145	LEU	3.1
2	N	44	GLU	3.1
1	C	187	SER	3.1
2	M	65	LYS	3.1
1	B	274	ILE	3.1
2	N	1	GLN	3.1
1	A	128	LEU	3.0
2	K	114	TYR	3.0
1	A	193	LEU	3.0
2	L	3	GLN	3.0
2	L	110	THR	3.0
1	C	147	GLU	3.0
1	E	393	LYS	3.0
1	D	127	GLY	2.9
1	D	89	ASP	2.9
1	E	25	ILE	2.9
1	E	162	ASP	2.9
1	E	147	GLU	2.9
1	A	114	ARG	2.9
1	B	114	ARG	2.8
1	C	142	ARG	2.8
1	A	271	THR	2.8
1	B	271	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	126	TYR	2.8
2	O	1	GLN	2.8
1	C	247	SER	2.8
2	K	69	THR	2.8
1	E	142	ARG	2.7
1	E	194	VAL	2.7
1	B	195	SER	2.7
1	B	110	THR	2.7
1	E	205	TYR	2.7
1	E	318	PRO	2.7
1	E	215	LYS	2.7
2	L	68	PHE	2.7
1	B	102	LYS	2.6
1	C	171	GLY	2.6
1	B	77	ILE	2.6
1	D	278	ASN	2.6
1	D	316	SER	2.6
1	E	12	VAL	2.6
2	N	45	ARG	2.6
1	C	126	TYR	2.6
1	C	146	ASP	2.6
1	E	260	VAL	2.6
1	D	31	PHE	2.5
1	D	279	SER	2.5
1	C	114	ARG	2.5
1	E	141	ARG	2.5
1	B	158	GLY	2.5
1	C	195	SER	2.5
2	O	86	LEU	2.5
2	K	96	CYS	2.5
1	E	273	SER	2.5
2	L	111	ASN	2.5
1	C	172	ASP	2.4
1	C	217	ASN	2.4
1	C	229	PHE	2.4
2	L	106	LEU	2.4
1	C	194	VAL	2.4
1	D	66	TYR	2.4
1	D	299	PHE	2.4
2	L	43	LYS	2.4
2	M	67	ARG	2.3
1	D	145	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	112	LYS	2.3
1	C	115	MET	2.3
1	D	82	THR	2.3
1	D	416	THR	2.3
2	K	112	TYR	2.3
1	B	147	GLU	2.3
1	B	128	LEU	2.3
2	L	70	ILE	2.3
1	E	203	GLY	2.3
2	M	86	LEU	2.3
2	K	1	GLN	2.3
1	A	194	VAL	2.3
1	B	194	VAL	2.3
1	D	76	GLY	2.2
2	K	98	ALA	2.2
1	C	193	LEU	2.2
1	A	209	SER	2.2
1	B	46	SER	2.2
2	L	66	ASP	2.2
1	B	177	GLY	2.2
1	D	78	PRO	2.2
2	N	11	LEU	2.2
1	C	271	THR	2.2
1	B	127	GLY	2.2
1	B	135	ALA	2.2
2	O	15	GLY	2.2
1	D	402	PHE	2.2
1	A	111	VAL	2.2
1	D	260	VAL	2.2
1	D	300	VAL	2.2
1	A	130	ILE	2.2
2	L	103	SER	2.2
1	E	14	GLU	2.2
1	A	96	THR	2.1
1	D	93	VAL	2.1
2	M	11	LEU	2.1
1	C	174	ALA	2.1
1	A	39	GLY	2.1
1	A	129	ARG	2.1
2	O	123	VAL	2.1
1	A	232	GLN	2.1
2	L	88	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	95	ASP	2.1
2	L	41	PRO	2.1
1	E	189	VAL	2.1
1	C	110	THR	2.1
1	B	170	GLY	2.1
2	M	42	GLY	2.1
1	A	62	TYR	2.1
1	B	14	GLU	2.1
1	C	143	TYR	2.0
2	L	107	TYR	2.0
2	L	113	ASP	2.0
1	D	147	GLU	2.0
2	M	52	SER	2.0
2	M	54	SER	2.0
2	L	53	TRP	2.0
1	D	257	ALA	2.0
1	C	127	GLY	2.0
1	A	236	PRO	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.

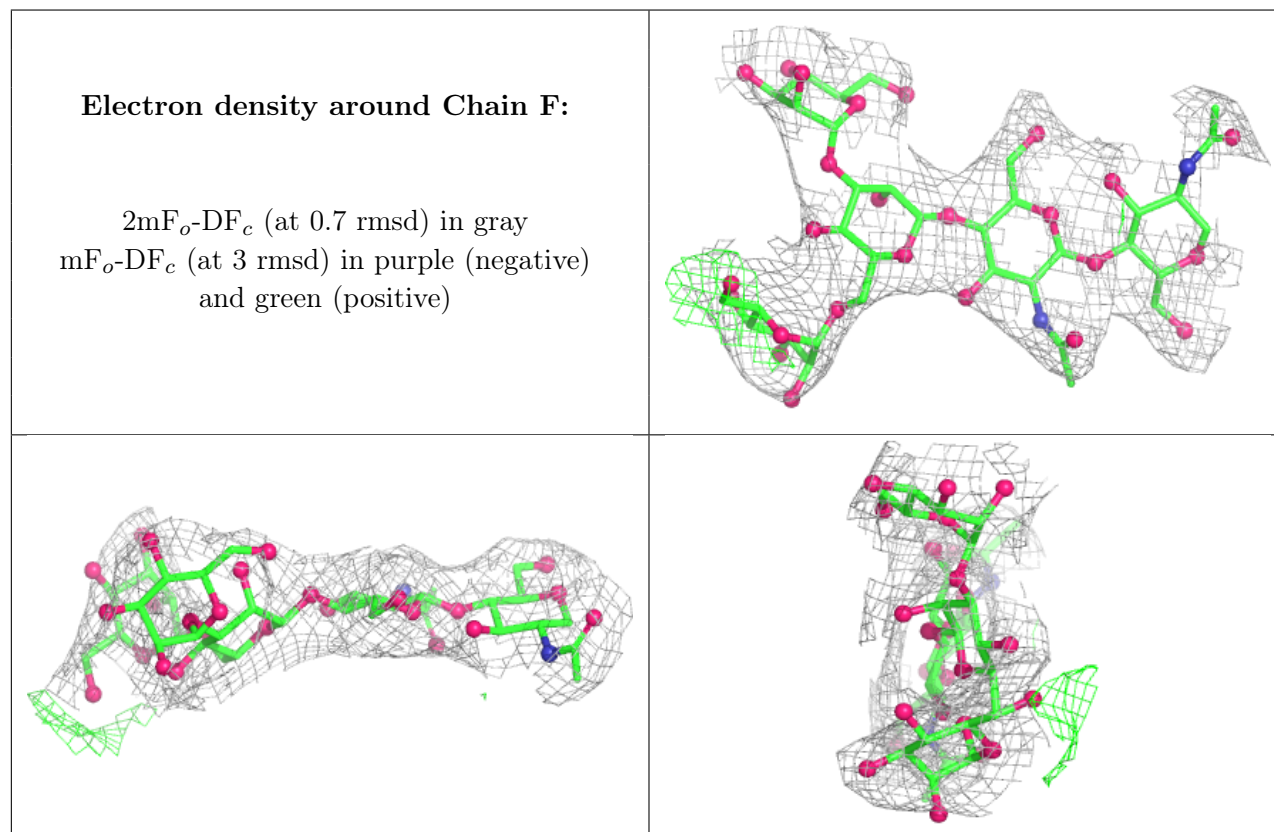
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	F	1	14/15	-	-	54,59,63,71	0
3	NAG	F	2	14/15	-	-	65,69,73,79	0
3	BMA	F	3	11/12	-	-	86,96,110,122	0
3	MAN	F	4	11/12	-	-	93,99,112,112	0
3	MAN	F	5	11/12	-	-	102,109,116,117	0
3	NAG	G	1	14/15	-	-	71,77,81,84	0
3	NAG	G	2	14/15	-	-	84,88,98,101	0
3	BMA	G	3	11/12	-	-	101,108,113,121	0
3	MAN	G	4	11/12	-	-	98,108,114,118	0
3	MAN	G	5	11/12	-	-	98,114,118,120	0

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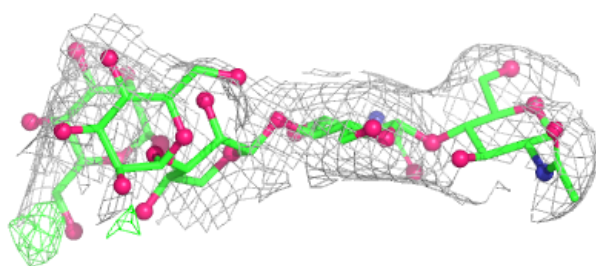
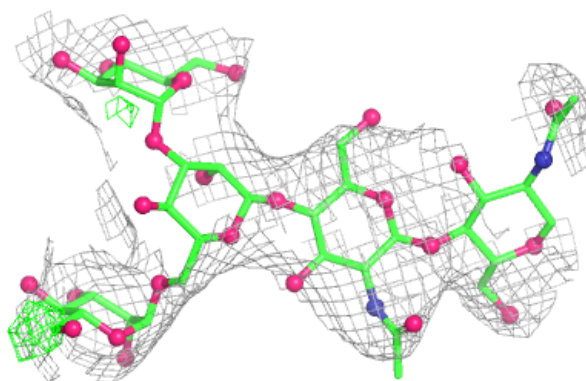
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	H	1	14/15	-	-	68,75,81,85	0
3	NAG	H	2	14/15	-	-	65,78,87,99	0
3	BMA	H	3	11/12	-	-	104,113,127,138	0
3	MAN	H	4	11/12	-	-	105,120,130,132	0
3	MAN	H	5	11/12	-	-	113,123,130,133	0
4	NAG	I	1	14/15	-	-	46,49,53,55	0
4	NAG	I	2	14/15	-	-	53,64,74,87	0
4	BMA	I	3	11/12	-	-	93,99,114,118	0
4	MAN	I	4	11/12	-	-	104,123,129,138	0
4	MAN	I	5	11/12	-	-	121,129,137,141	0
4	MAN	I	6	11/12	-	-	108,123,134,135	0
4	NAG	J	1	14/15	-	-	66,68,74,77	0
4	NAG	J	2	14/15	-	-	56,70,77,80	0
4	BMA	J	3	11/12	-	-	78,83,90,105	0
4	MAN	J	4	11/12	-	-	112,120,129,139	0
4	MAN	J	5	11/12	-	-	93,115,122,123	0
4	MAN	J	6	11/12	-	-	80,86,92,94	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.

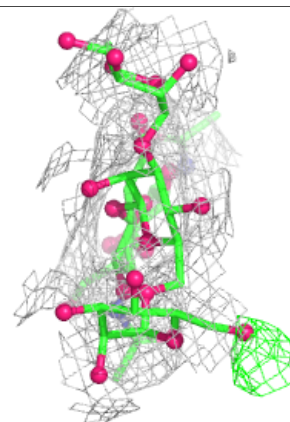
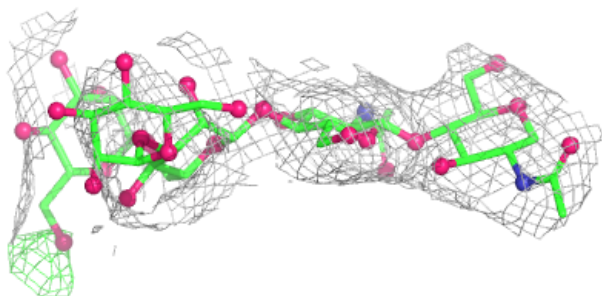
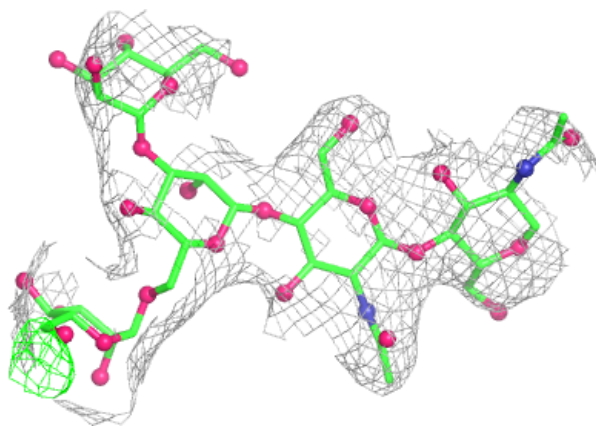


Electron density around Chain G:

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

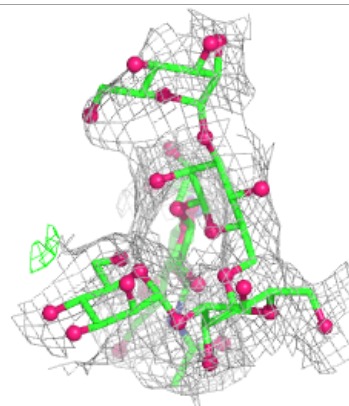
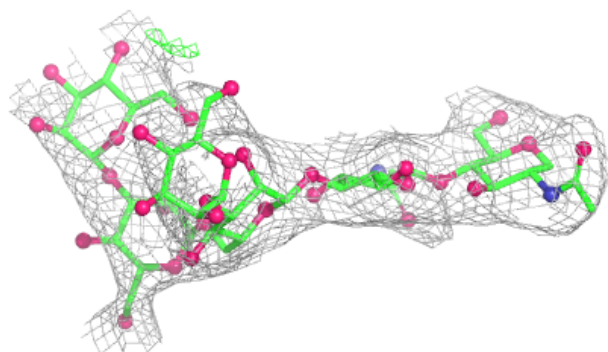
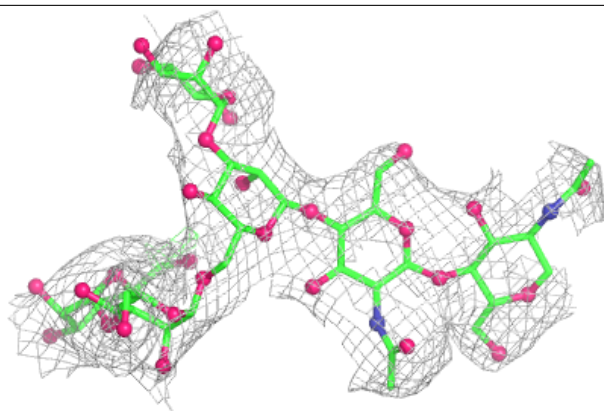
**Electron density around Chain H:**

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

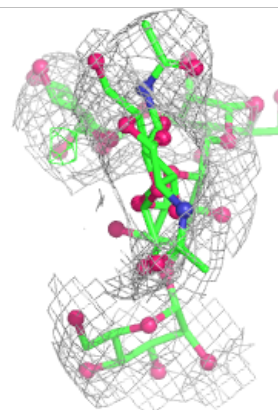
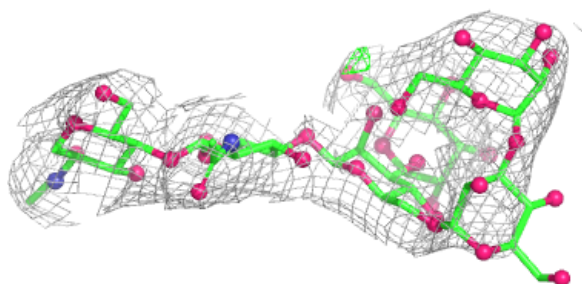
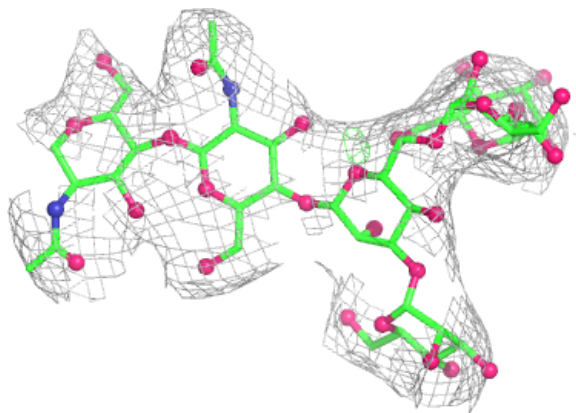


Electron density around Chain I:

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

**Electron density around Chain J:**

$2mF_o-DF_c$ (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
5	NAG	A	501	14/15	0.72	0.17	122,141,148,149	0
5	NAG	D	501	14/15	0.77	0.18	93,102,115,117	0
5	NAG	E	501	14/15	0.77	0.14	101,108,114,117	0
5	NAG	C	501	14/15	0.84	0.12	140,153,164,166	0
5	NAG	B	501	14/15	0.93	0.10	96,100,112,119	0

6.5 Other polymers [i](#)

There are no such residues in this entry.