



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:07 PM UTC

PDB ID : 4XHJ / pdb_00004xhj
Title : gHgL of Varicella-zoster virus in complex with human neutralizing antibodies.
Authors : Xing, Y.
Deposited on : 2015-01-05
Resolution : 3.16 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

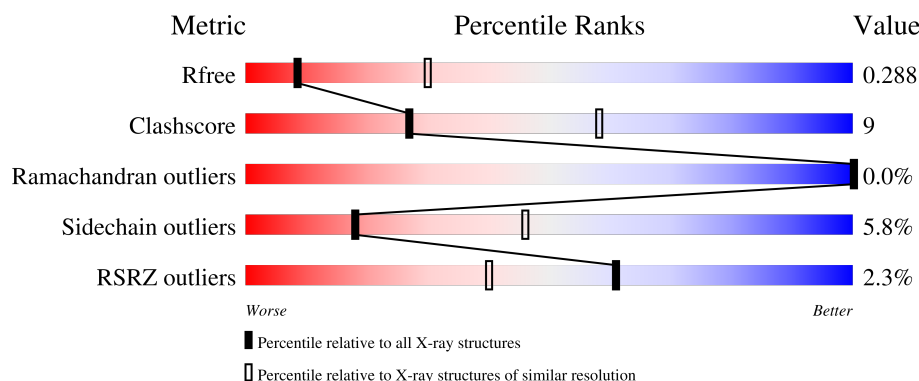
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.16 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	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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	833	% 62% 24% • 12%
1	E	833	3% 60% 23% • 15%
2	B	138	% 69% 24% • •
2	F	138	% 67% 22% 5% 6%
3	C	215	4% 79% 18% •

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Mol	Chain	Length	Quality of chain
3	G	215	3%73%17%9%
4	D	282	%55%16%27%
4	H	282	%55%18%27%
5	I	3	33%67%
6	J	5	60%20%20%
6	M	5	100%
7	K	2	50%50%
7	L	2	50%50%

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 19505 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 Envelope glycoprotein H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	729	Total	C	N	O	S	0	0	0
			5610	3602	936	1047	25			
1	E	705	Total	C	N	O	S	0	0	0
			5452	3508	906	1015	23			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	796	GLY	-	expression tag	UNP Q775J3
A	797	SER	-	expression tag	UNP Q775J3
A	798	GLU	-	expression tag	UNP Q775J3
A	799	ASN	-	expression tag	UNP Q775J3
A	800	LEU	-	expression tag	UNP Q775J3
A	801	TYR	-	expression tag	UNP Q775J3
A	802	PHE	-	expression tag	UNP Q775J3
A	803	GLN	-	expression tag	UNP Q775J3
A	804	GLY	-	expression tag	UNP Q775J3
A	805	SER	-	expression tag	UNP Q775J3
A	806	TRP	-	expression tag	UNP Q775J3
A	807	SER	-	expression tag	UNP Q775J3
A	808	HIS	-	expression tag	UNP Q775J3
A	809	PRO	-	expression tag	UNP Q775J3
A	810	GLN	-	expression tag	UNP Q775J3
A	811	PHE	-	expression tag	UNP Q775J3
A	812	GLU	-	expression tag	UNP Q775J3
A	813	LYS	-	expression tag	UNP Q775J3
A	814	GLY	-	expression tag	UNP Q775J3
A	815	GLY	-	expression tag	UNP Q775J3
A	816	GLY	-	expression tag	UNP Q775J3
A	817	SER	-	expression tag	UNP Q775J3
A	818	GLY	-	expression tag	UNP Q775J3
A	819	GLY	-	expression tag	UNP Q775J3
A	820	GLY	-	expression tag	UNP Q775J3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	821	SER	-	expression tag	UNP Q775J3
A	822	GLY	-	expression tag	UNP Q775J3
A	823	GLY	-	expression tag	UNP Q775J3
A	824	GLY	-	expression tag	UNP Q775J3
A	825	SER	-	expression tag	UNP Q775J3
A	826	TRP	-	expression tag	UNP Q775J3
A	827	SER	-	expression tag	UNP Q775J3
A	828	HIS	-	expression tag	UNP Q775J3
A	829	PRO	-	expression tag	UNP Q775J3
A	830	GLN	-	expression tag	UNP Q775J3
A	831	PHE	-	expression tag	UNP Q775J3
A	832	GLU	-	expression tag	UNP Q775J3
A	833	LYS	-	expression tag	UNP Q775J3
E	796	GLY	-	expression tag	UNP Q775J3
E	797	SER	-	expression tag	UNP Q775J3
E	798	GLU	-	expression tag	UNP Q775J3
E	799	ASN	-	expression tag	UNP Q775J3
E	800	LEU	-	expression tag	UNP Q775J3
E	801	TYR	-	expression tag	UNP Q775J3
E	802	PHE	-	expression tag	UNP Q775J3
E	803	GLN	-	expression tag	UNP Q775J3
E	804	GLY	-	expression tag	UNP Q775J3
E	805	SER	-	expression tag	UNP Q775J3
E	806	TRP	-	expression tag	UNP Q775J3
E	807	SER	-	expression tag	UNP Q775J3
E	808	HIS	-	expression tag	UNP Q775J3
E	809	PRO	-	expression tag	UNP Q775J3
E	810	GLN	-	expression tag	UNP Q775J3
E	811	PHE	-	expression tag	UNP Q775J3
E	812	GLU	-	expression tag	UNP Q775J3
E	813	LYS	-	expression tag	UNP Q775J3
E	814	GLY	-	expression tag	UNP Q775J3
E	815	GLY	-	expression tag	UNP Q775J3
E	816	GLY	-	expression tag	UNP Q775J3
E	817	SER	-	expression tag	UNP Q775J3
E	818	GLY	-	expression tag	UNP Q775J3
E	819	GLY	-	expression tag	UNP Q775J3
E	820	GLY	-	expression tag	UNP Q775J3
E	821	SER	-	expression tag	UNP Q775J3
E	822	GLY	-	expression tag	UNP Q775J3
E	823	GLY	-	expression tag	UNP Q775J3
E	824	GLY	-	expression tag	UNP Q775J3

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Chain	Residue	Modelled	Actual	Comment	Reference
E	825	SER	-	expression tag	UNP Q775J3
E	826	TRP	-	expression tag	UNP Q775J3
E	827	SER	-	expression tag	UNP Q775J3
E	828	HIS	-	expression tag	UNP Q775J3
E	829	PRO	-	expression tag	UNP Q775J3
E	830	GLN	-	expression tag	UNP Q775J3
E	831	PHE	-	expression tag	UNP Q775J3
E	832	GLU	-	expression tag	UNP Q775J3
E	833	LYS	-	expression tag	UNP Q775J3

- Molecule 2 is a protein called Envelope glycoprotein L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	132	Total	C	N	O	S	0	0	0
			1007	647	163	192	5			
2	F	130	Total	C	N	O	S	0	0	0
			997	641	161	190	5			

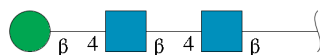
- Molecule 3 is a protein called Fab-RC light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	207	Total	C	N	O	S	0	0	0
			1569	984	265	315	5			
3	G	196	Total	C	N	O	S	0	0	0
			1481	936	247	293	5			

- Molecule 4 is a protein called Fab-RC heavy chain.

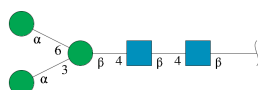
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	207	Total	C	N	O	S	0	0	0
			1569	992	265	299	13			
4	H	206	Total	C	N	O	S	0	0	0
			1561	988	264	296	13			

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 6 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
6	J	5	Total	C	N	O	0	0	0
			61	34	2	25			
6	M	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 7 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
7	L	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

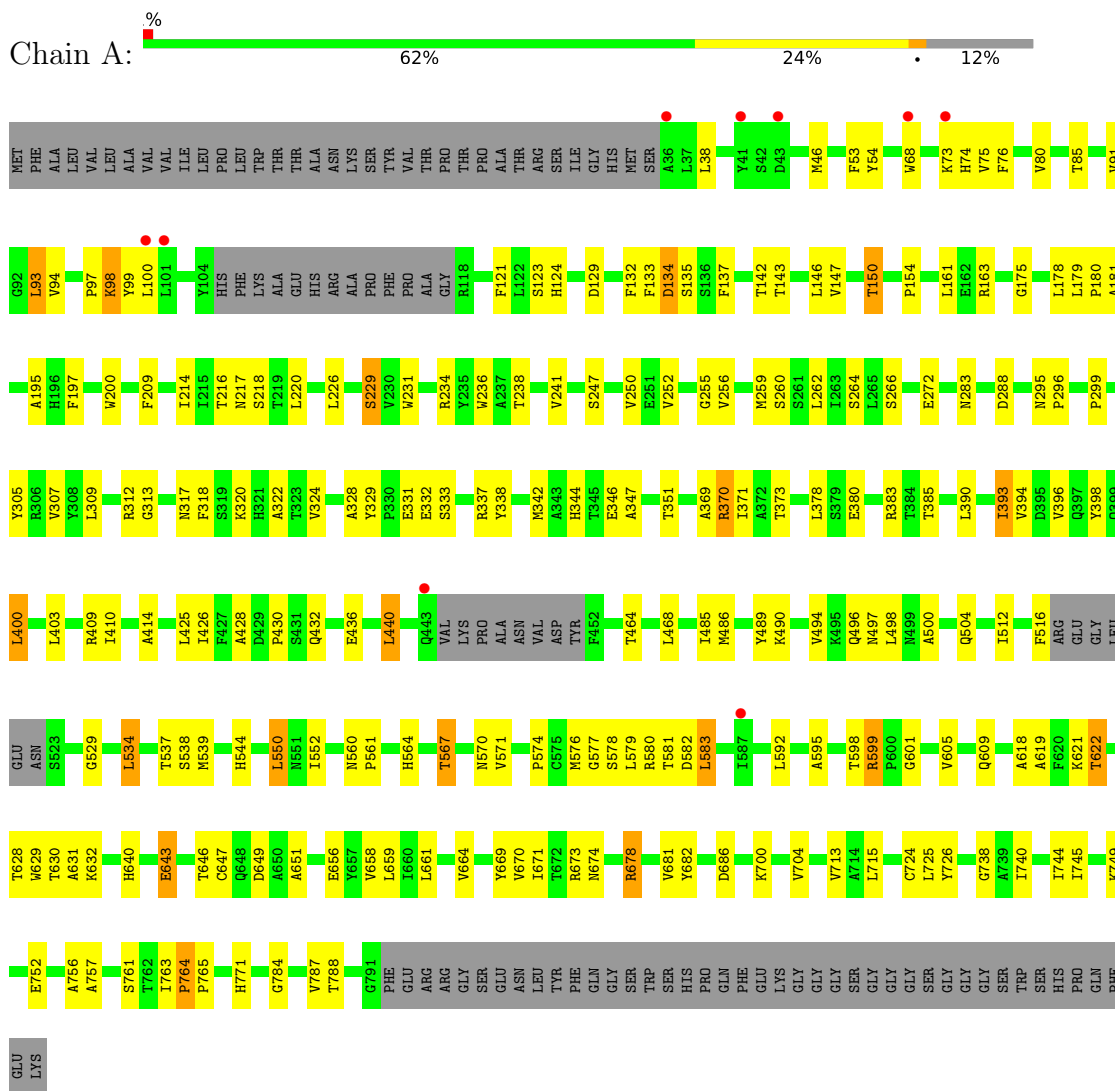


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	E	1	Total	C	N	O	0	0
			14	8	1	5		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Envelope glycoprotein H

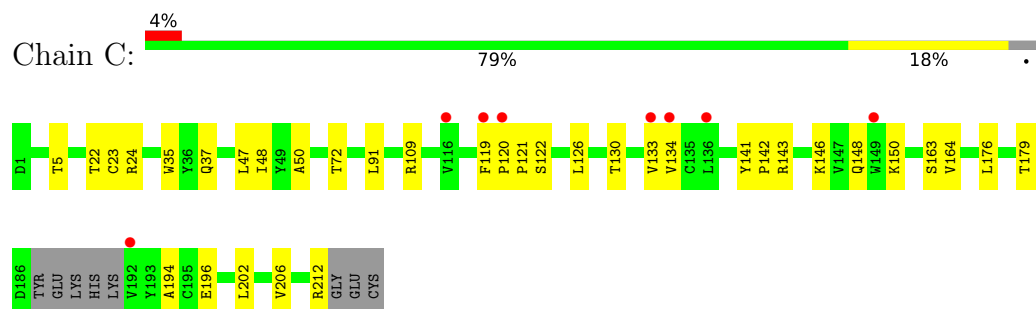


- Molecule 1: Envelope glycoprotein H

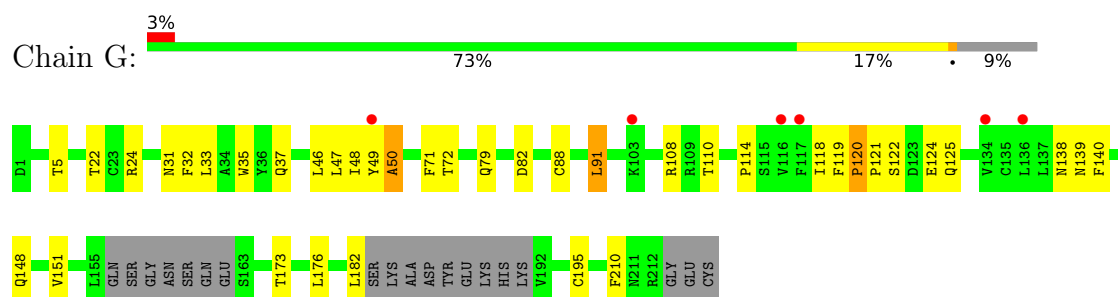




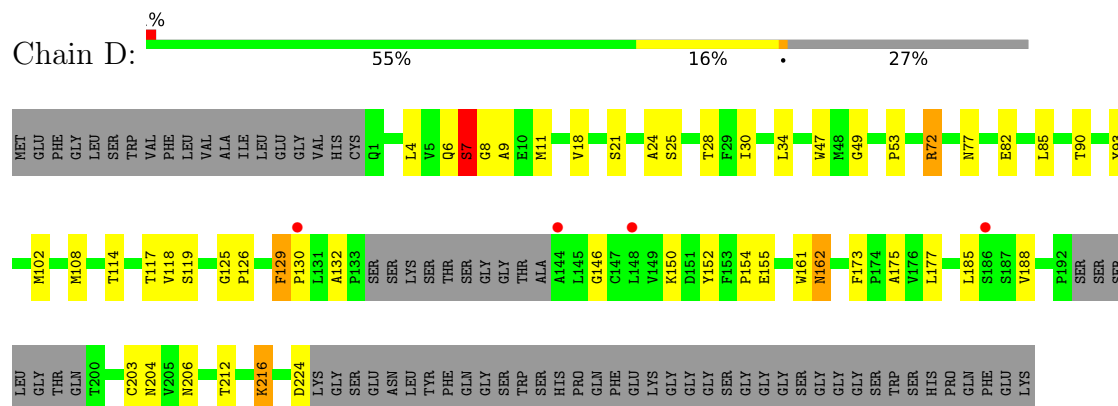
- Molecule 3: Fab-RC light chain



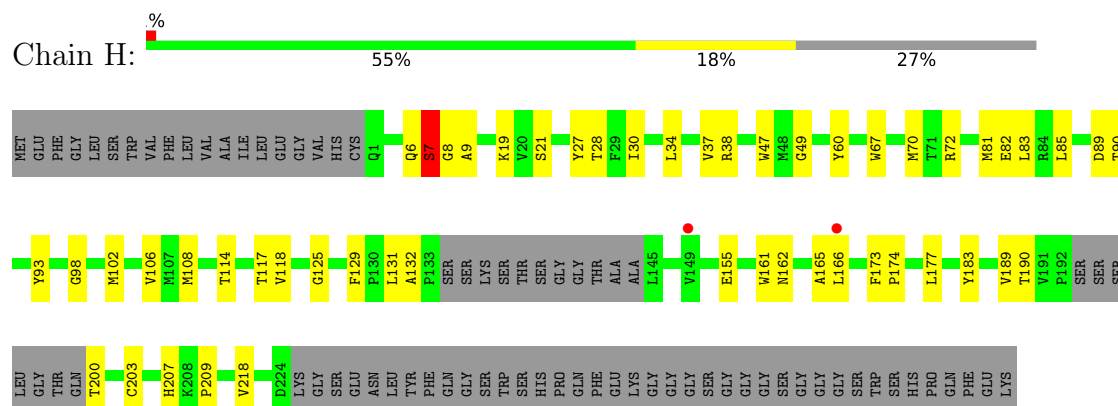
- Molecule 3: Fab-RC light chain



- Molecule 4: Fab-RC heavy chain



- Molecule 4: Fab-RC heavy chain



- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  33% 67%

NAG1
NAG2
BMA3

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

Chain J:  60% 20% 20%

NAG1
NAG2
BMA3
MAN4
MAN5

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

Chain M:  100%

NAG1
NAG2
BMA3
MAN4
MAN5

- Molecule 7: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  50% 50%

NAG1
NAG2

- Molecule 7: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  50% 50%

NAG1
NAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	147.00Å 147.00Å 198.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.10 – 3.16 41.10 – 3.16	Depositor EDS
% Data completeness (in resolution range)	99.6 (41.10-3.16) 99.7 (41.10-3.16)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 3.18Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1839)	Depositor
R, R_{free}	0.236 , 0.278 0.244 , 0.288	Depositor DCC
R_{free} test set	3636 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	101.8	Xtriage
Anisotropy	0.601	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 70.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.047 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	19505	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.35	0/5749	0.82	7/7854 (0.1%)
1	E	0.33	0/5582	0.81	9/7619 (0.1%)
2	B	0.40	1/1034 (0.1%)	0.97	6/1416 (0.4%)
2	F	0.32	0/1024	0.96	4/1402 (0.3%)
3	C	0.32	0/1600	0.81	3/2173 (0.1%)
3	G	0.31	0/1511	0.78	2/2055 (0.1%)
4	D	0.31	0/1608	0.82	6/2191 (0.3%)
4	H	0.30	0/1600	0.80	5/2180 (0.2%)
All	All	0.33	1/19708 (0.0%)	0.83	42/26890 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
4	D	0	1
4	H	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	64	VAL	C-N	-5.79	1.26	1.33

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	7	SER	N-CA-C	7.64	120.83	108.76
3	G	120	PRO	N-CA-C	7.38	119.71	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	36	LYS	CA-C-N	-7.26	113.00	120.85
2	F	36	LYS	C-N-CA	-7.26	113.00	120.85
4	H	7	SER	N-CA-C	7.26	120.23	108.76
2	B	36	LYS	CA-C-N	-7.25	113.02	120.85
2	B	36	LYS	C-N-CA	-7.25	113.02	120.85
1	E	54	TYR	CA-C-N	7.09	128.70	119.84
1	E	54	TYR	C-N-CA	7.09	128.70	119.84
2	B	131	VAL	N-CA-C	6.87	117.58	110.36
2	B	30	PRO	N-CA-CB	6.77	110.36	103.25
1	A	599	ARG	CA-C-N	6.52	126.29	119.05
1	A	599	ARG	C-N-CA	6.52	126.29	119.05
3	C	120	PRO	N-CA-C	6.40	118.51	110.70
4	D	125	GLY	CA-C-N	6.33	126.52	120.31
4	D	125	GLY	C-N-CA	6.33	126.52	120.31
1	E	764	PRO	N-CA-C	6.29	118.38	110.70
3	C	119	PHE	CA-C-N	6.15	126.72	120.38
3	C	119	PHE	C-N-CA	6.15	126.72	120.38
1	A	764	PRO	N-CA-C	6.03	118.06	110.70
4	H	125	GLY	CA-C-N	6.00	126.19	120.31
4	H	125	GLY	C-N-CA	6.00	126.19	120.31
2	F	99	ASN	CA-C-N	5.82	125.02	118.97
2	F	99	ASN	C-N-CA	5.82	125.02	118.97
1	E	599	ARG	CA-C-N	5.71	125.83	119.32
1	E	599	ARG	C-N-CA	5.71	125.83	119.32
2	B	142	ILE	CA-C-N	-5.70	114.88	120.52
2	B	142	ILE	C-N-CA	-5.70	114.88	120.52
3	G	50	ALA	N-CA-C	-5.68	101.55	110.36
4	D	129	PHE	CA-C-N	5.50	125.50	119.89
4	D	129	PHE	C-N-CA	5.50	125.50	119.89
4	H	129	PHE	CA-C-N	5.49	125.49	119.89
4	H	129	PHE	C-N-CA	5.49	125.49	119.89
1	E	133	PHE	N-CA-C	5.45	116.90	111.07
4	D	162	ASN	CB-CA-C	-5.38	110.39	116.63
1	A	133	PHE	N-CA-C	5.27	116.70	111.07
1	E	85	THR	N-CA-C	-5.26	106.45	112.92
1	A	681	VAL	N-CA-C	5.11	115.21	107.80
1	A	744	ILE	N-CA-C	5.09	115.96	109.30
1	A	85	THR	N-CA-C	-5.06	106.69	112.92
1	E	415	HIS	CA-C-N	5.05	125.32	119.47
1	E	415	HIS	C-N-CA	5.05	125.32	119.47

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	D	7	SER	Peptide
4	H	7	SER	Peptide

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	5610	0	5452	122	0
1	E	5452	0	5349	117	0
2	B	1007	0	972	22	0
2	F	997	0	969	30	0
3	C	1569	0	1527	20	0
3	G	1481	0	1440	22	0
4	D	1569	0	1517	29	0
4	H	1561	0	1510	28	0
5	I	39	0	34	1	0
6	J	61	0	52	1	0
6	M	61	0	52	0	0
7	K	28	0	25	0	0
7	L	28	0	25	0	0
8	A	28	0	26	0	0
8	E	14	0	13	0	0
All	All	19505	0	18963	354	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 9.

All (354) 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:98:LYS:HD2	2:B:68:SER:HB3	1.57	0.86
1:E:175:GLY:HA3	1:E:333:SER:HA	1.63	0.79
1:E:73:LYS:O	1:E:98:LYS:NZ	2.17	0.78
3:G:5:THR:HB	3:G:24:ARG:HB2	1.69	0.75
1:E:61:GLU:HB3	1:E:64:LYS:HB2	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:757:ALA:HB3	1:E:761:SER:HB3	1.68	0.74
1:E:161:LEU:HB3	1:E:324:VAL:HG12	1.73	0.70
1:E:394:VAL:HG11	1:E:464:THR:HG22	1.74	0.70
3:C:202:LEU:HD13	3:C:206:VAL:HG22	1.74	0.69
1:A:175:GLY:HA3	1:A:333:SER:HA	1.72	0.69
1:A:537:THR:HA	1:A:578:SER:HB2	1.75	0.68
1:E:98:LYS:HD2	2:F:68:SER:HB3	1.75	0.68
1:A:134:ASP:OD1	6:J:1:NAG:H82	1.94	0.68
1:A:163:ARG:NH2	1:A:331:GLU:OE2	2.27	0.67
1:E:579:LEU:HA	1:E:631:ALA:HB2	1.76	0.67
3:C:5:THR:HB	3:C:24:ARG:HB2	1.77	0.66
1:A:73:LYS:O	1:A:98:LYS:NZ	2.27	0.66
1:A:757:ALA:HB3	1:A:761:SER:HB3	1.76	0.66
2:F:144:GLU:O	2:F:157:ARG:NH2	2.29	0.66
2:B:149:TRP:CD1	2:B:150:GLY:H	2.14	0.66
1:E:163:ARG:NH2	1:E:331:GLU:OE2	2.28	0.66
1:A:342:MET:HB2	1:A:370:ARG:HD2	1.77	0.65
1:A:579:LEU:HA	1:A:631:ALA:HB2	1.78	0.65
1:E:129:ASP:OD2	1:E:135:SER:OG	2.13	0.65
2:F:149:TRP:CD1	2:F:150:GLY:H	2.14	0.65
4:D:175:ALA:HA	4:D:185:LEU:HB3	1.79	0.65
1:A:266:SER:HA	2:B:146:GLY:HA2	1.79	0.65
1:A:763:ILE:HG22	1:A:765:PRO:HD2	1.79	0.64
1:E:580:ARG:HG2	1:E:659:LEU:HD12	1.79	0.64
1:A:394:VAL:HG11	1:A:464:THR:HG22	1.77	0.64
3:C:122:SER:O	3:C:126:LEU:N	2.30	0.63
1:E:763:ILE:HG22	1:E:765:PRO:HD2	1.80	0.63
1:E:593:LEU:HD11	1:E:616:PHE:HD1	1.63	0.63
2:F:134:ASP:O	4:H:28:THR:OG1	2.18	0.62
1:A:238:THR:H	1:A:255:GLY:HA2	1.64	0.62
1:A:649:ASP:OD2	1:A:678:ARG:NH1	2.32	0.62
4:H:6:GLN:NE2	4:H:93:TYR:O	2.28	0.61
4:D:53:PRO:O	4:D:72:ARG:NH1	2.33	0.61
1:A:143:THR:O	2:B:99:ASN:ND2	2.29	0.61
1:A:580:ARG:HG2	1:A:659:LEU:HD12	1.83	0.61
1:A:75:VAL:HG23	1:A:98:LYS:HZ1	1.65	0.61
2:F:130:GLY:O	2:F:133:VAL:HG23	2.01	0.61
1:A:129:ASP:OD2	1:A:135:SER:OG	2.19	0.60
1:E:342:MET:HB2	1:E:370:ARG:HD2	1.84	0.60
1:E:534:LEU:HD22	1:E:745:ILE:HD11	1.84	0.60
1:E:724:CYS:HB2	1:E:749:LYS:HE2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:LEU:O	1:A:229:SER:OG	2.18	0.60
3:G:48:ILE:HG22	3:G:50:ALA:O	2.02	0.59
1:A:529:GLY:HA3	1:A:640:HIS:CD2	2.38	0.59
1:E:504:GLN:NE2	1:E:570:ASN:OD1	2.36	0.59
1:A:671:ILE:HG21	1:A:740:ILE:HD11	1.84	0.59
4:H:90:THR:HG23	4:H:117:THR:HA	1.85	0.59
2:B:108:THR:HG22	2:B:128:LEU:HD13	1.83	0.59
4:D:11:MET:HB2	4:D:154:PRO:HG3	1.83	0.59
1:E:628:THR:HA	1:E:632:LYS:HB2	1.83	0.59
4:H:85:LEU:HD13	4:H:118:VAL:HG22	1.82	0.59
1:A:161:LEU:HB3	1:A:324:VAL:HG12	1.85	0.58
1:A:534:LEU:HD22	1:A:745:ILE:HD11	1.86	0.58
1:E:70:ASN:HB2	2:F:90:LYS:HD2	1.85	0.57
2:F:40:ASP:HB2	2:F:43:LEU:HG	1.86	0.57
1:E:37:LEU:HD12	1:E:60:GLU:HG2	1.86	0.57
1:E:537:THR:HA	1:E:578:SER:HB2	1.85	0.57
1:E:179:LEU:O	1:E:181:ALA:N	2.37	0.57
1:A:313:GLY:HA2	1:A:317:ASN:HB2	1.87	0.57
1:E:238:THR:H	1:E:255:GLY:HA2	1.70	0.57
3:G:108:ARG:NH2	3:G:110:THR:OG1	2.37	0.57
3:C:130:THR:HG22	3:C:183:SER:HA	1.86	0.57
1:A:259:MET:HE1	1:A:373:THR:HG22	1.87	0.57
1:A:567:THR:HG22	1:A:570:ASN:HB2	1.86	0.57
1:E:682:TYR:OH	1:E:784:GLY:O	2.19	0.57
1:A:68:TRP:CD1	2:B:41:VAL:HB	2.40	0.56
1:A:178:LEU:HD11	1:A:338:TYR:HA	1.87	0.56
1:E:582:ASP:HB3	1:E:658:VAL:HG13	1.86	0.56
4:D:146:GLY:HA3	4:D:188:VAL:HA	1.87	0.56
2:F:108:THR:HG22	2:F:128:LEU:HD13	1.87	0.56
1:A:715:LEU:HD13	1:A:756:ALA:HA	1.88	0.56
3:C:134:VAL:HG12	3:C:179:THR:HG23	1.88	0.56
1:E:226:LEU:N	1:E:272:GLU:OE1	2.37	0.56
1:A:179:LEU:O	1:A:181:ALA:N	2.39	0.56
1:A:724:CYS:HB2	1:A:749:LYS:HE2	1.87	0.56
1:A:226:LEU:HD13	1:A:262:LEU:HD11	1.87	0.55
4:D:85:LEU:HD13	4:D:118:VAL:HG22	1.89	0.55
1:A:486:MET:O	1:A:490:LYS:HG3	2.07	0.55
1:A:647:CYS:C	1:A:649:ASP:H	2.14	0.55
1:E:197:PHE:HD2	1:E:390:LEU:HD21	1.71	0.55
1:E:266:SER:HA	2:F:146:GLY:HA2	1.88	0.55
1:E:259:MET:HE1	1:E:373:THR:HG22	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:133:VAL:O	3:C:180:LEU:HB3	2.08	0.54
2:B:79:LYS:NZ	2:B:134:ASP:OD2	2.39	0.54
1:E:713:VAL:HG22	1:E:787:VAL:HG11	1.89	0.54
1:A:195:ALA:HB2	1:A:209:PHE:CD1	2.43	0.54
1:E:671:ILE:HG21	1:E:740:ILE:HD11	1.88	0.54
1:A:68:TRP:HZ3	1:A:93:LEU:HD22	1.72	0.54
4:D:146:GLY:HA2	4:D:161:TRP:CH2	2.42	0.54
3:G:37:GLN:HB2	3:G:47:LEU:HD11	1.88	0.54
1:E:624:MET:O	1:E:628:THR:HG22	2.09	0.53
4:H:207:HIS:CD2	4:H:209:PRO:HD2	2.43	0.53
1:E:264:SER:HB2	2:F:147:CYS:HA	1.90	0.53
1:A:393:ILE:HD11	1:A:428:ALA:HB2	1.90	0.53
1:A:618:ALA:O	1:A:622:THR:HG22	2.08	0.53
1:A:264:SER:HB2	2:B:147:CYS:HA	1.91	0.53
1:A:504:GLN:NE2	1:A:570:ASN:OD1	2.42	0.53
1:E:529:GLY:HA3	1:E:640:HIS:CD2	2.43	0.53
1:E:543:ALA:O	1:E:546:THR:OG1	2.25	0.53
1:A:765:PRO:HA	1:A:771:HIS:NE2	2.23	0.53
4:D:4:LEU:HD23	4:D:24:ALA:HA	1.91	0.53
1:A:197:PHE:HD2	1:A:390:LEU:HD21	1.73	0.52
1:E:179:LEU:HB2	1:E:563:LYS:HD3	1.91	0.52
1:A:220:LEU:HD11	1:A:234:ARG:HB3	1.90	0.52
1:E:49:LYS:HZ3	1:E:68:TRP:HZ2	1.55	0.52
1:E:247:SER:H	1:E:414:ALA:HB2	1.75	0.52
2:B:134:ASP:O	4:D:28:THR:OG1	2.28	0.52
1:E:489:TYR:HB2	1:E:539:MET:HE2	1.92	0.52
1:A:661:LEU:HB2	1:A:669:TYR:HB2	1.91	0.52
4:D:130:PRO:HD3	4:D:216:LYS:HD3	1.91	0.52
2:F:133:VAL:HG12	4:H:27:TYR:HB2	1.92	0.52
2:F:62:PRO:HB2	2:F:64:VAL:HG12	1.92	0.52
1:A:214:ILE:O	1:A:218:SER:OG	2.27	0.52
1:A:646:THR:HG21	1:A:700:LYS:HA	1.92	0.52
1:A:68:TRP:HD1	2:B:41:VAL:HB	1.75	0.51
3:C:212:ARG:NH2	4:D:224:ASP:H	2.09	0.51
1:E:143:THR:O	2:F:99:ASN:ND2	2.37	0.51
2:B:40:ASP:HB2	2:B:43:LEU:HG	1.91	0.51
1:A:226:LEU:N	1:A:272:GLU:OE1	2.41	0.51
1:A:646:THR:OG1	1:A:678:ARG:NH2	2.41	0.51
1:E:393:ILE:HD11	1:E:428:ALA:HB2	1.91	0.51
1:E:661:LEU:HD21	1:E:743:ILE:HG21	1.93	0.51
1:A:725:LEU:HG	5:I:2:NAG:H83	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:ALA:O	1:A:504:GLN:HG3	2.12	0.50
3:C:37:GLN:HB2	3:C:47:LEU:HD11	1.93	0.50
1:A:305:TYR:OH	1:A:380:GLU:HG2	2.12	0.50
1:E:195:ALA:HB2	1:E:209:PHE:CD1	2.46	0.50
1:A:628:THR:HA	1:A:632:LYS:HB2	1.93	0.50
4:H:161:TRP:CZ3	4:H:203:CYS:HB3	2.47	0.50
1:E:269:LEU:O	1:E:365:TYR:OH	2.23	0.50
1:A:595:ALA:O	1:A:598:THR:OG1	2.24	0.50
3:G:91:LEU:HD21	4:H:106:VAL:HG22	1.93	0.50
1:E:147:VAL:O	2:F:139:SER:OG	2.30	0.49
1:E:620:PHE:HD2	1:E:621:LYS:HE2	1.77	0.49
1:A:121:PHE:HE2	1:A:383:ARG:HH12	1.60	0.49
1:A:651:ALA:HB1	1:A:656:GLU:HB2	1.94	0.49
1:A:686:ASP:OD1	1:A:686:ASP:N	2.35	0.49
2:B:144:GLU:O	2:B:157:ARG:NH2	2.43	0.49
1:A:147:VAL:O	2:B:139:SER:OG	2.29	0.49
3:C:48:ILE:HG22	3:C:50:ALA:O	2.11	0.49
1:E:340:LEU:HD13	1:E:399:GLN:HG3	1.95	0.49
4:H:155:GLU:HG2	4:H:183:TYR:CD1	2.48	0.49
1:A:74:HIS:ND1	1:A:97:PRO:HA	2.28	0.49
1:A:247:SER:H	1:A:414:ALA:HB2	1.78	0.48
1:A:180:PRO:HB2	1:A:337:ARG:HH12	1.77	0.48
1:A:260:SER:OG	1:A:272:GLU:OE2	2.30	0.48
1:E:651:ALA:HB1	1:E:656:GLU:HB2	1.95	0.48
1:E:381:MET:O	1:E:385:THR:OG1	2.28	0.48
4:H:166:LEU:HD13	4:H:189:VAL:HG21	1.96	0.48
1:A:344:HIS:CD2	1:A:403:LEU:HD21	2.48	0.48
1:A:124:HIS:HB3	1:A:236:TRP:CG	2.49	0.48
1:E:68:TRP:CD1	2:F:41:VAL:HB	2.48	0.48
4:D:9:ALA:HA	4:D:114:THR:HG23	1.96	0.48
1:E:256:VAL:HG13	1:E:383:ARG:HH11	1.78	0.48
4:D:129:PHE:HE2	4:D:150:LYS:HE2	1.79	0.48
1:E:347:ALA:O	1:E:351:THR:HG23	2.13	0.48
4:H:6:GLN:HA	4:H:21:SER:O	2.14	0.48
3:C:179:THR:HG22	3:C:181:THR:HG23	1.96	0.48
2:F:154:LEU:H	2:F:154:LEU:HG	1.35	0.48
1:A:146:LEU:HD23	1:A:150:THR:HG22	1.96	0.47
4:D:7:SER:HA	4:D:8:GLY:HA2	1.58	0.47
1:E:226:LEU:HD13	1:E:262:LEU:HD11	1.96	0.47
1:A:80:VAL:HA	1:A:91:VAL:HG12	1.96	0.47
1:E:312:ARG:NH2	2:F:140:THR:O	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:592:LEU:HD22	1:E:623:MET:HE2	1.96	0.47
1:E:669:TYR:OH	1:E:781:PHE:O	2.21	0.47
3:G:119:PHE:HB3	4:H:131:LEU:HB3	1.95	0.47
1:A:318:PHE:CD1	1:A:322:ALA:HB3	2.49	0.47
3:G:138:ASN:ND2	3:G:139:ASN:OD1	2.47	0.47
1:A:351:THR:HG22	1:A:410:ILE:HD13	1.97	0.47
4:D:6:GLN:NE2	4:D:93:TYR:O	2.37	0.47
1:E:329:TYR:CE2	1:E:370:ARG:HG2	2.50	0.47
1:E:506:LEU:HD21	1:E:574:PRO:HG3	1.95	0.47
1:A:489:TYR:HB2	1:A:539:MET:HE2	1.97	0.47
4:D:162:ASN:O	4:D:204:ASN:ND2	2.47	0.47
3:G:114:PRO:HB3	3:G:140:PHE:HB3	1.97	0.47
1:E:546:THR:HA	1:E:592:LEU:HD13	1.97	0.47
3:G:32:PHE:HB3	3:G:91:LEU:HB3	1.96	0.47
1:E:260:SER:OG	1:E:272:GLU:OE2	2.33	0.47
2:B:156:GLN:HG2	2:B:158:VAL:HG23	1.96	0.47
3:C:148:GLN:HB2	3:C:196:GLU:HB3	1.97	0.47
1:E:497:ASN:OD1	1:E:498:LEU:N	2.45	0.47
4:H:162:ASN:HB2	4:H:165:ALA:HB3	1.97	0.47
1:A:425:LEU:HB2	1:A:428:ALA:HB3	1.97	0.46
3:C:150:LYS:HG2	3:C:194:ALA:HB3	1.97	0.46
1:E:305:TYR:OH	1:E:380:GLU:HG2	2.15	0.46
1:E:351:THR:HG22	1:E:410:ILE:HD13	1.95	0.46
1:A:582:ASP:HB3	1:A:658:VAL:CG1	2.45	0.46
4:D:6:GLN:HA	4:D:21:SER:O	2.14	0.46
1:A:238:THR:O	1:A:426:ILE:HG12	2.16	0.46
1:A:497:ASN:OD1	1:A:498:LEU:N	2.45	0.46
1:A:544:HIS:NE2	1:A:574:PRO:HD2	2.30	0.46
1:E:132:PHE:CE1	2:F:106:GLY:HA3	2.51	0.46
4:H:7:SER:HA	4:H:8:GLY:HA2	1.59	0.46
4:H:60:TYR:HE1	4:H:70:MET:HG3	1.80	0.46
3:G:79:GLN:HB2	3:G:82:ASP:OD1	2.16	0.46
1:A:142:THR:OG1	2:B:154:LEU:HD13	2.16	0.45
1:A:347:ALA:O	1:A:351:THR:HG23	2.16	0.45
1:A:216:THR:OG1	1:A:217:ASN:N	2.49	0.45
4:D:130:PRO:HG3	4:D:216:LYS:HB3	1.98	0.45
1:A:371:ILE:HD11	1:A:403:LEU:HB3	1.98	0.45
4:H:37:VAL:HG22	4:H:47:TRP:HA	1.98	0.45
1:E:382:GLY:HA3	1:E:388:PHE:CZ	2.51	0.45
3:G:120:PRO:HB3	3:G:210:PHE:CE1	2.50	0.45
1:E:75:VAL:HG23	1:E:98:LYS:NZ	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:400:LEU:HD12	1:E:400:LEU:HA	1.80	0.45
3:C:109:ARG:H	3:C:109:ARG:HD2	1.81	0.45
4:D:90:THR:HG23	4:D:117:THR:HA	1.98	0.45
1:A:646:THR:HG1	1:A:678:ARG:HH21	1.62	0.45
1:E:75:VAL:HG23	1:E:98:LYS:HZ1	1.81	0.45
1:A:643:GLU:HA	1:A:646:THR:HG23	1.98	0.45
1:E:96:PHE:CG	2:F:124:LEU:HD11	2.52	0.45
1:E:100:LEU:HD12	1:E:100:LEU:HA	1.81	0.45
1:E:742:ASP:HB3	1:E:744:ILE:HD13	1.99	0.45
1:A:396:VAL:O	1:A:400:LEU:HB2	2.17	0.44
1:E:396:VAL:O	1:E:400:LEU:HB2	2.17	0.44
1:A:582:ASP:HB3	1:A:658:VAL:HG13	1.97	0.44
2:F:133:VAL:HG12	4:H:27:TYR:CB	2.48	0.44
1:A:409:ARG:NH2	1:A:628:THR:O	2.51	0.44
4:H:38:ARG:HH12	4:H:89:ASP:HA	1.82	0.44
1:E:70:ASN:OD1	1:E:71:ASP:N	2.51	0.44
1:E:199:THR:OG1	1:E:201:ASP:OD2	2.31	0.44
1:E:54:TYR:HA	1:E:55:PRO:HD2	1.84	0.44
1:A:351:THR:HG21	1:A:629:TRP:CZ3	2.53	0.44
1:A:605:VAL:O	1:A:609:GLN:HG2	2.17	0.44
2:F:156:GLN:HG2	2:F:158:VAL:HG23	1.99	0.44
3:G:46:LEU:HD21	3:G:49:TYR:HB3	2.00	0.44
3:G:118:ILE:HD12	3:G:195:CYS:HB2	2.00	0.43
3:C:141:TYR:CG	3:C:142:PRO:HA	2.53	0.43
1:E:67:HIS:O	1:E:67:HIS:ND1	2.51	0.43
1:E:783:ASN:OD1	1:E:783:ASN:N	2.49	0.43
1:A:256:VAL:HG13	1:A:383:ARG:HH11	1.82	0.43
1:A:577:GLY:HA2	1:A:630:THR:HG21	2.00	0.43
2:B:154:LEU:H	2:B:154:LEU:HG	1.32	0.43
1:A:550:LEU:HD12	1:A:550:LEU:HA	1.85	0.43
1:E:290:THR:HB	3:G:49:TYR:OH	2.19	0.43
2:F:134:ASP:HA	4:H:27:TYR:HA	2.00	0.43
4:H:9:ALA:HA	4:H:114:THR:HG23	2.00	0.43
1:E:647:CYS:C	1:E:649:ASP:H	2.26	0.43
2:F:132:TRP:HB3	2:F:135:SER:OG	2.18	0.43
3:C:121:PRO:HG2	4:D:132:ALA:H	1.84	0.43
1:E:194:LYS:NZ	1:E:429:ASP:OD2	2.49	0.43
1:E:587:ILE:HD12	1:E:587:ILE:H	1.84	0.43
1:A:75:VAL:HG13	2:B:73:GLY:HA2	2.00	0.43
1:A:329:TYR:CE2	1:A:370:ARG:HG2	2.54	0.43
1:A:592:LEU:HD23	1:A:619:ALA:HB1	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:328:ALA:O	1:E:332:GLU:HG3	2.19	0.43
2:B:62:PRO:HB2	2:B:64:VAL:HG12	2.01	0.42
1:E:482:ARG:HH12	1:E:528:ASP:CG	2.26	0.42
1:E:491:GLY:O	1:E:495:LYS:HG2	2.18	0.42
3:C:23:CYS:HB2	3:C:35:TRP:CH2	2.54	0.42
1:E:146:LEU:HD23	1:E:150:THR:HG22	2.00	0.42
1:A:560:ASN:HB3	1:A:564:HIS:ND1	2.35	0.42
1:A:581:THR:OG1	1:A:658:VAL:HA	2.19	0.42
1:A:713:VAL:HG22	1:A:787:VAL:HG11	2.00	0.42
1:A:726:TYR:HB2	1:A:752:GLU:CD	2.44	0.42
1:E:567:THR:OG1	1:E:569:PRO:HD2	2.19	0.42
4:H:98:GLY:HA2	4:H:106:VAL:O	2.19	0.42
4:D:18:VAL:O	4:D:82:GLU:HA	2.19	0.42
4:D:161:TRP:CZ3	4:D:203:CYS:HB3	2.54	0.42
1:E:80:VAL:HA	1:E:91:VAL:HG12	2.01	0.42
1:E:142:THR:OG1	2:F:154:LEU:HD13	2.20	0.42
1:A:180:PRO:O	1:A:337:ARG:NH1	2.46	0.42
1:A:351:THR:HG21	1:A:629:TRP:HZ3	1.85	0.42
4:D:11:MET:HE3	4:D:119:SER:HB3	2.01	0.42
4:D:24:ALA:HB3	4:D:77:ASN:HB3	2.02	0.42
2:F:36:LYS:NZ	2:F:47:GLU:OE1	2.42	0.42
1:A:68:TRP:CZ3	1:A:93:LEU:HD22	2.54	0.42
1:A:154:PRO:HD3	1:A:312:ARG:NH2	2.35	0.42
1:E:309:LEU:HD13	1:E:314:LEU:HD21	2.01	0.42
1:E:470:ARG:HD2	1:E:470:ARG:HA	1.90	0.42
1:A:231:TRP:HD1	1:A:283:ASN:HB3	1.85	0.42
4:D:126:PRO:HD2	4:D:212:THR:HG21	2.02	0.42
1:E:60:GLU:HG3	1:E:61:GLU:N	2.35	0.42
1:E:332:GLU:OE2	1:E:370:ARG:NH2	2.53	0.42
1:E:344:HIS:CD2	1:E:403:LEU:HD21	2.55	0.42
1:E:425:LEU:HB2	1:E:428:ALA:HB3	2.01	0.42
1:E:569:PRO:HA	1:E:572:TYR:HB2	2.00	0.42
1:E:659:LEU:HB3	1:E:671:ILE:HG13	2.01	0.42
1:A:550:LEU:HB3	1:A:599:ARG:NH1	2.34	0.42
4:D:173:PHE:HE1	4:D:188:VAL:HG22	1.85	0.42
1:E:593:LEU:HD21	1:E:616:PHE:HA	2.02	0.42
1:E:200:TRP:CE2	1:E:440:LEU:HD22	2.55	0.41
3:G:121:PRO:HG2	4:H:132:ALA:HB3	2.02	0.41
2:B:78:THR:O	2:B:135:SER:HB3	2.20	0.41
1:E:68:TRP:CZ3	1:E:93:LEU:HD13	2.55	0.41
1:E:580:ARG:NH2	1:E:582:ASP:OD2	2.37	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:PHE:CE1	2:B:73:GLY:HA3	2.54	0.41
1:A:241:VAL:HB	1:A:252:VAL:HB	2.03	0.41
1:A:432:GLN:O	1:A:436:GLU:HG3	2.20	0.41
3:C:150:LYS:HD3	3:C:150:LYS:HA	1.86	0.41
3:C:163:SER:OG	4:D:173:PHE:HB3	2.21	0.41
1:E:295:ASN:HA	1:E:296:PRO:HD2	1.84	0.41
3:G:139:ASN:HA	3:G:173:THR:HB	2.01	0.41
4:H:47:TRP:CZ2	4:H:49:GLY:HA2	2.55	0.41
2:B:76:VAL:HG22	2:B:85:VAL:HG22	2.02	0.41
1:E:500:ALA:O	1:E:504:GLN:HG3	2.20	0.41
1:E:544:HIS:NE2	1:E:574:PRO:HD2	2.36	0.41
1:A:295:ASN:HA	1:A:296:PRO:HD2	1.87	0.41
1:A:328:ALA:O	1:A:332:GLU:HG3	2.20	0.41
1:A:673:ARG:NH2	1:A:738:GLY:O	2.54	0.41
1:E:596:ILE:N	1:E:597:PRO:HD2	2.35	0.41
4:H:67:TRP:O	4:H:83:LEU:HD12	2.20	0.41
1:A:132:PHE:HB2	1:A:137:PHE:HA	2.03	0.41
1:A:369:ALA:O	1:A:373:THR:HG23	2.20	0.41
1:A:560:ASN:HA	1:A:561:PRO:HD2	1.91	0.41
1:E:351:THR:HG21	1:E:629:TRP:CZ3	2.56	0.41
1:A:75:VAL:HG23	1:A:98:LYS:NZ	2.33	0.41
1:A:299:PRO:HD2	1:A:307:VAL:O	2.21	0.41
1:A:393:ILE:HD12	1:A:393:ILE:HA	1.87	0.41
1:E:74:HIS:ND1	1:E:97:PRO:HA	2.35	0.41
3:G:35:TRP:CZ3	3:G:88:CYS:HB3	2.56	0.41
4:D:47:TRP:CZ2	4:D:49:GLY:HA2	2.55	0.41
1:E:37:LEU:HD22	1:E:53:PHE:HB2	2.03	0.41
1:E:625:ILE:O	1:E:629:TRP:HB2	2.21	0.41
1:A:38:LEU:HD13	2:B:123:ALA:HB1	2.02	0.41
1:A:538:SER:HB2	1:A:745:ILE:HG13	2.03	0.41
1:E:231:TRP:HD1	1:E:283:ASN:HB3	1.86	0.41
2:F:64:VAL:HG21	2:F:70:ALA:HB2	2.03	0.41
3:G:182:LEU:HD23	3:G:182:LEU:HA	1.92	0.41
4:H:19:LYS:HE2	4:H:82:GLU:OE2	2.21	0.41
1:A:200:TRP:CE2	1:A:440:LEU:HD22	2.56	0.41
1:A:318:PHE:HB3	1:A:320:LYS:O	2.21	0.41
1:A:583:LEU:HD13	1:A:583:LEU:HA	1.74	0.41
1:A:763:ILE:HA	1:A:764:PRO:HD2	1.97	0.41
1:E:216:THR:OG1	1:E:217:ASN:N	2.53	0.41
3:G:120:PRO:HD2	3:G:121:PRO:HD3	2.02	0.41
4:H:177:LEU:H	4:H:177:LEU:HD12	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:143:ARG:NH2	3:C:164:VAL:HG21	2.37	0.40
3:C:146:LYS:HD2	3:C:146:LYS:HA	1.87	0.40
4:D:152:TYR:OH	4:D:155:GLU:OE1	2.36	0.40
1:E:130:VAL:HG21	2:F:113:GLU:OE2	2.21	0.40
1:E:478:LEU:O	1:E:482:ARG:HG3	2.22	0.40
2:F:153:ARG:HH11	2:F:153:ARG:HB3	1.86	0.40
4:H:81:MET:HE3	4:H:81:MET:HB3	2.01	0.40
1:A:430:PRO:HB3	1:A:468:LEU:HD12	2.04	0.40
1:E:39:ARG:O	1:E:93:LEU:N	2.52	0.40
1:E:121:PHE:HE2	1:E:383:ARG:HH12	1.69	0.40
1:E:152:PHE:CZ	1:E:229:SER:HB2	2.56	0.40
1:E:418:THR:HB	2:F:160:GLN:NE2	2.36	0.40
4:D:24:ALA:O	4:D:25:SER:OG	2.38	0.40
3:G:33:LEU:HD22	3:G:71:PHE:CG	2.56	0.40
4:H:173:PHE:HA	4:H:174:PRO:HD3	1.98	0.40
1:A:601:GLY:O	1:A:605:VAL:HG23	2.21	0.40
1:A:682:TYR:OH	1:A:784:GLY:O	2.38	0.40
1:E:179:LEU:HD12	1:E:179:LEU:HA	1.86	0.40
2:F:136:THR:HA	2:F:137:PRO:HD3	1.88	0.40
3:G:122:SER:HB3	3:G:125:GLN:HB3	2.03	0.40
3:G:124:GLU:OE1	3:G:124:GLU:N	2.52	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	721/833 (87%)	687 (95%)	34 (5%)	0	100	100
1	E	691/833 (83%)	660 (96%)	31 (4%)	0	100	100
2	B	130/138 (94%)	119 (92%)	10 (8%)	1 (1%)	16	46
2	F	128/138 (93%)	115 (90%)	13 (10%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	203/215 (94%)	196 (97%)	7 (3%)	0	100	100
3	G	190/215 (88%)	182 (96%)	8 (4%)	0	100	100
4	D	201/282 (71%)	192 (96%)	9 (4%)	0	100	100
4	H	200/282 (71%)	192 (96%)	8 (4%)	0	100	100
All	All	2464/2936 (84%)	2343 (95%)	120 (5%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	30	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	604/717 (84%)	560 (93%)	44 (7%)	13	38
1	E	591/717 (82%)	554 (94%)	37 (6%)	16	44
2	B	111/121 (92%)	103 (93%)	8 (7%)	13	39
2	F	111/121 (92%)	104 (94%)	7 (6%)	16	44
3	C	176/187 (94%)	172 (98%)	4 (2%)	44	66
3	G	165/187 (88%)	158 (96%)	7 (4%)	26	55
4	D	173/233 (74%)	165 (95%)	8 (5%)	24	53
4	H	172/233 (74%)	164 (95%)	8 (5%)	23	52
All	All	2103/2516 (84%)	1980 (94%)	123 (6%)	18	46

All (123) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	MET
1	A	53	PHE
1	A	54	TYR

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Mol	Chain	Res	Type
1	A	93	LEU
1	A	94	VAL
1	A	98	LYS
1	A	99	TYR
1	A	100	LEU
1	A	123	SER
1	A	134	ASP
1	A	150	THR
1	A	229	SER
1	A	250	VAL
1	A	288	ASP
1	A	309	LEU
1	A	346	GLU
1	A	370	ARG
1	A	378	LEU
1	A	385	THR
1	A	393	ILE
1	A	398	TYR
1	A	400	LEU
1	A	440	LEU
1	A	485	ILE
1	A	494	VAL
1	A	496	GLN
1	A	512	ILE
1	A	516	PHE
1	A	534	LEU
1	A	550	LEU
1	A	552	ILE
1	A	567	THR
1	A	571	VAL
1	A	576	MET
1	A	583	LEU
1	A	621	LYS
1	A	622	THR
1	A	643	GLU
1	A	664	VAL
1	A	670	VAL
1	A	674	ASN
1	A	678	ARG
1	A	704	VAL
1	A	788	THR
2	B	44	ILE

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Mol	Chain	Res	Type
2	B	46	THR
2	B	104	LEU
2	B	114	GLU
2	B	116	LYS
2	B	152	ASP
2	B	154	LEU
2	B	157	ARG
3	C	22	THR
3	C	72	THR
3	C	91	LEU
3	C	176	LEU
4	D	30	ILE
4	D	34	LEU
4	D	72	ARG
4	D	102	MET
4	D	108	MET
4	D	177	LEU
4	D	206	ASN
4	D	216	LYS
1	E	41	TYR
1	E	50	LEU
1	E	67	HIS
1	E	93	LEU
1	E	94	VAL
1	E	98	LYS
1	E	100	LEU
1	E	123	SER
1	E	150	THR
1	E	199	THR
1	E	250	VAL
1	E	288	ASP
1	E	309	LEU
1	E	334	LEU
1	E	346	GLU
1	E	358	ASP
1	E	370	ARG
1	E	378	LEU
1	E	385	THR
1	E	391	ASP
1	E	398	TYR
1	E	400	LEU
1	E	440	LEU

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Mol	Chain	Res	Type
1	E	485	ILE
1	E	496	GLN
1	E	512	ILE
1	E	534	LEU
1	E	550	LEU
1	E	552	ILE
1	E	583	LEU
1	E	621	LYS
1	E	630	THR
1	E	632	LYS
1	E	643	GLU
1	E	670	VAL
1	E	678	ARG
1	E	711	GLU
2	F	44	ILE
2	F	47	GLU
2	F	104	LEU
2	F	132	TRP
2	F	153	ARG
2	F	154	LEU
2	F	157	ARG
3	G	22	THR
3	G	31	ASN
3	G	72	THR
3	G	91	LEU
3	G	148	GLN
3	G	151	VAL
3	G	176	LEU
4	H	30	ILE
4	H	34	LEU
4	H	72	ARG
4	H	102	MET
4	H	108	MET
4	H	190	THR
4	H	200	THR
4	H	218	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	HIS
1	A	206	HIS

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Mol	Chain	Res	Type
1	A	397	GLN
3	C	55	GLN
3	C	90	GLN
3	C	92	ASN
3	C	159	ASN
3	C	161	GLN
1	E	144	GLN
1	E	145	HIS
1	E	206	HIS
2	F	156	GLN
2	F	160	GLN
3	G	90	GLN

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](#)

17 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	I	1	1,5	14,14,15	0.59	0	17,19,21	0.89	1 (5%)
5	NAG	I	2	5	14,14,15	0.44	0	17,19,21	1.08	0
5	BMA	I	3	5	11,11,12	0.71	0	15,15,17	0.66	0
6	NAG	J	1	6,2	14,14,15	0.95	1 (7%)	17,19,21	1.16	2 (11%)
6	NAG	J	2	6	14,14,15	0.57	0	17,19,21	0.71	0
6	BMA	J	3	6	11,11,12	0.99	1 (9%)	15,15,17	1.27	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	MAN	J	4	6	11,11,12	0.69	0	15,15,17	0.75	0
6	MAN	J	5	6	11,11,12	0.64	0	15,15,17	0.46	0
7	NAG	K	1	7,1	14,14,15	0.66	0	17,19,21	1.01	0
7	NAG	K	2	7	14,14,15	0.61	0	17,19,21	0.90	1 (5%)
7	NAG	L	1	7,1	14,14,15	0.47	0	17,19,21	0.83	1 (5%)
7	NAG	L	2	7	14,14,15	0.61	0	17,19,21	0.71	0
6	NAG	M	1	6,2	14,14,15	0.52	0	17,19,21	0.95	1 (5%)
6	NAG	M	2	6	14,14,15	0.59	0	17,19,21	0.94	1 (5%)
6	BMA	M	3	6	11,11,12	1.18	0	15,15,17	0.94	1 (6%)
6	MAN	M	4	6	11,11,12	0.59	0	15,15,17	0.91	1 (6%)
6	MAN	M	5	6	11,11,12	0.51	0	15,15,17	0.93	1 (6%)

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	I	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	I	2	5	-	0/6/23/26	0/1/1/1
5	BMA	I	3	5	-	0/2/19/22	0/1/1/1
6	NAG	J	1	6,2	-	2/6/23/26	0/1/1/1
6	NAG	J	2	6	-	0/6/23/26	0/1/1/1
6	BMA	J	3	6	-	1/2/19/22	0/1/1/1
6	MAN	J	4	6	-	0/2/19/22	0/1/1/1
6	MAN	J	5	6	-	0/2/19/22	0/1/1/1
7	NAG	K	1	7,1	-	0/6/23/26	0/1/1/1
7	NAG	K	2	7	-	2/6/23/26	0/1/1/1
7	NAG	L	1	7,1	-	4/6/23/26	0/1/1/1
7	NAG	L	2	7	-	2/6/23/26	0/1/1/1
6	NAG	M	1	6,2	-	2/6/23/26	0/1/1/1
6	NAG	M	2	6	-	0/6/23/26	0/1/1/1
6	BMA	M	3	6	-	2/2/19/22	0/1/1/1
6	MAN	M	4	6	-	0/2/19/22	0/1/1/1
6	MAN	M	5	6	-	1/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	J	1	NAG	O5-C1	-3.19	1.38	1.43
6	J	3	BMA	O5-C1	-2.29	1.39	1.43

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	M	5	MAN	C1-O5-C5	2.97	116.16	112.19
6	M	2	NAG	C4-C3-C2	2.86	115.21	111.02
6	J	1	NAG	O5-C1-C2	-2.70	107.11	111.29
6	J	3	BMA	C2-C3-C4	2.68	115.57	110.86
6	J	1	NAG	O4-C4-C5	-2.36	103.52	109.32
6	M	1	NAG	O5-C1-C2	-2.29	107.74	111.29
7	K	2	NAG	C1-O5-C5	2.29	115.25	112.19
5	I	1	NAG	C4-C3-C2	2.24	114.30	111.02
7	L	1	NAG	C1-O5-C5	2.20	115.14	112.19
6	M	4	MAN	C1-O5-C5	2.13	115.03	112.19
6	J	3	BMA	C1-C2-C3	2.07	112.66	109.64
6	M	3	BMA	C3-C4-C5	2.05	113.95	110.23

There are no chirality outliers.

All (18) torsion outliers are listed below:

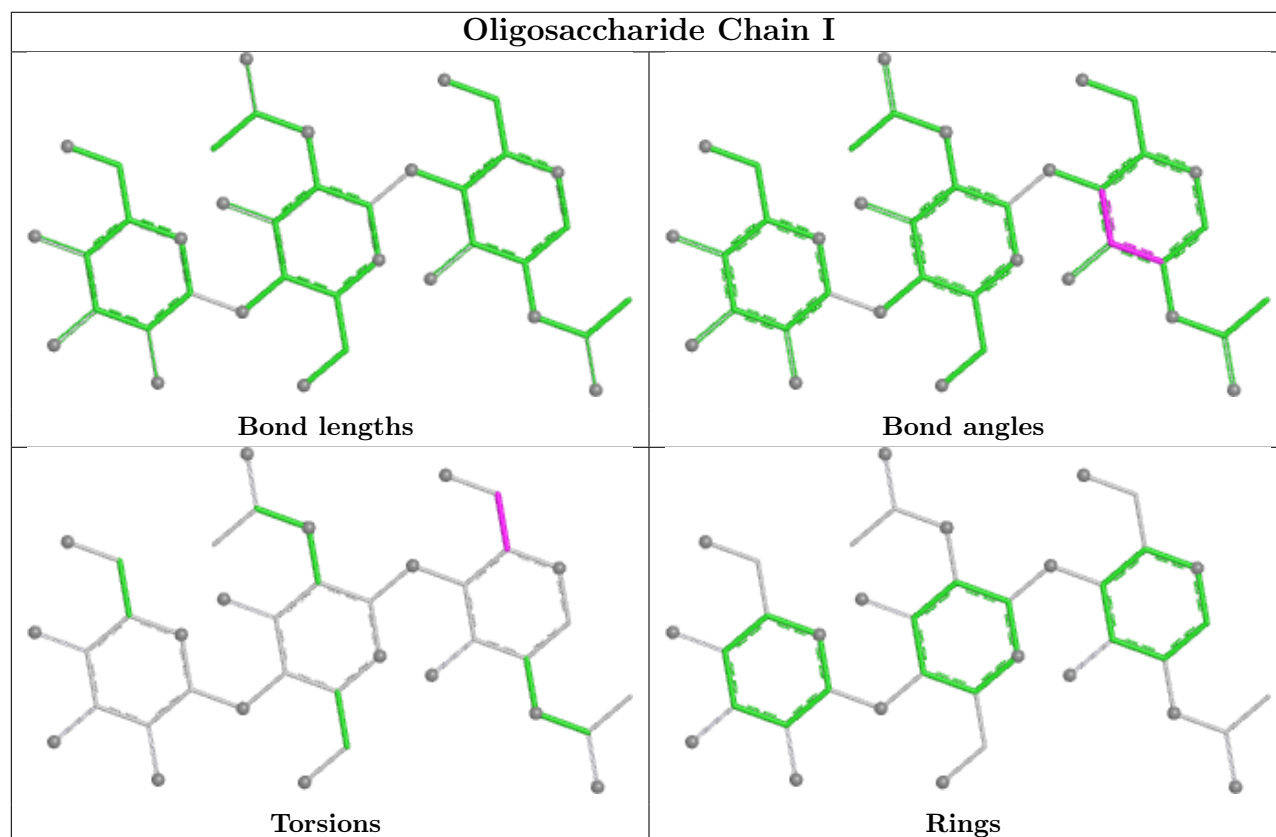
Mol	Chain	Res	Type	Atoms
7	K	2	NAG	C8-C7-N2-C2
7	K	2	NAG	O7-C7-N2-C2
6	M	3	BMA	C4-C5-C6-O6
7	L	2	NAG	C8-C7-N2-C2
7	L	2	NAG	O7-C7-N2-C2
6	M	3	BMA	O5-C5-C6-O6
5	I	1	NAG	O5-C5-C6-O6
7	L	1	NAG	C8-C7-N2-C2
5	I	1	NAG	C4-C5-C6-O6
6	J	3	BMA	O5-C5-C6-O6
6	J	1	NAG	C4-C5-C6-O6
7	L	1	NAG	C4-C5-C6-O6
7	L	1	NAG	O7-C7-N2-C2
6	M	5	MAN	O5-C5-C6-O6
7	L	1	NAG	O5-C5-C6-O6
6	M	1	NAG	C8-C7-N2-C2
6	J	1	NAG	O5-C5-C6-O6
6	M	1	NAG	O7-C7-N2-C2

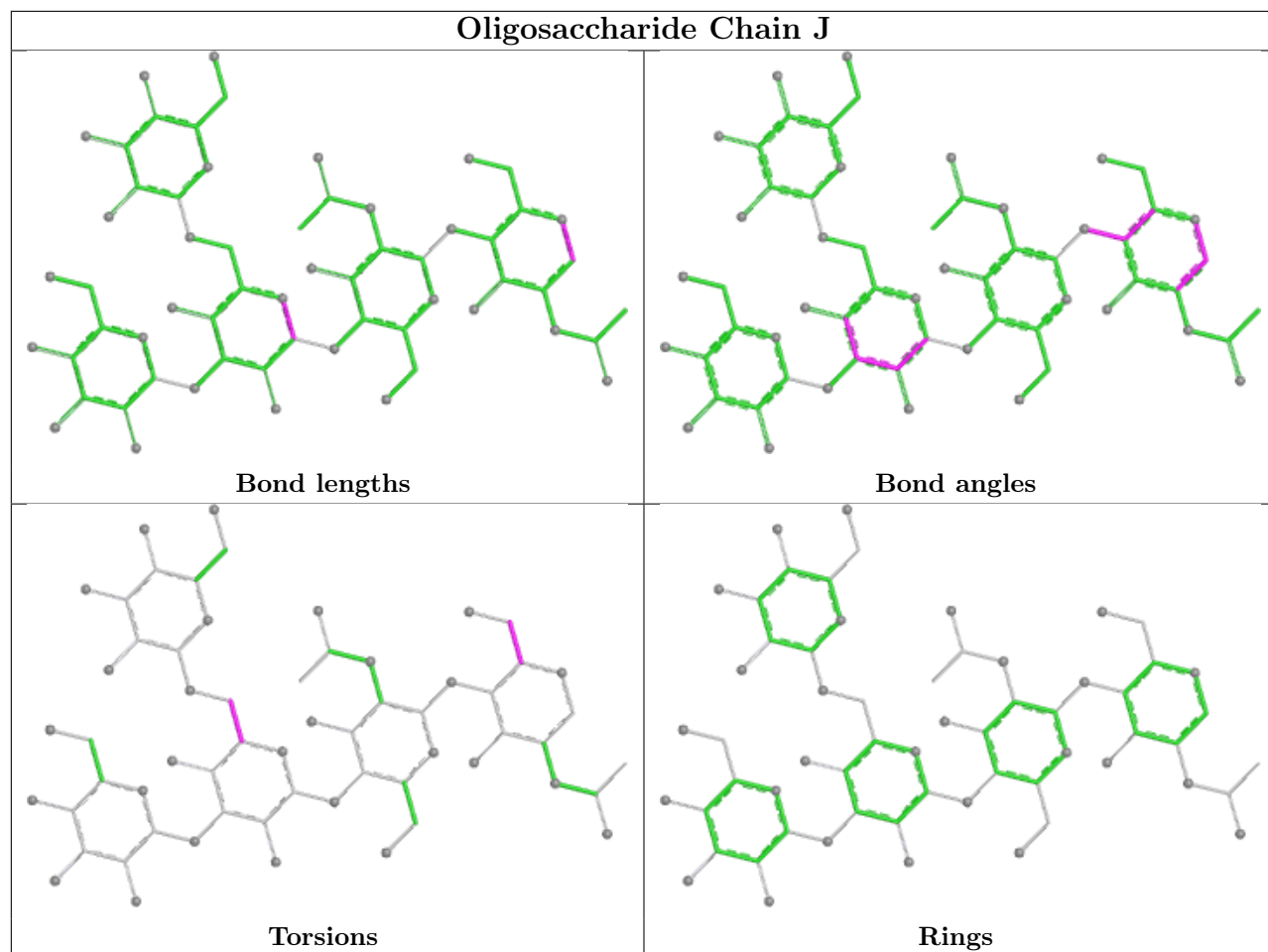
There are no ring outliers.

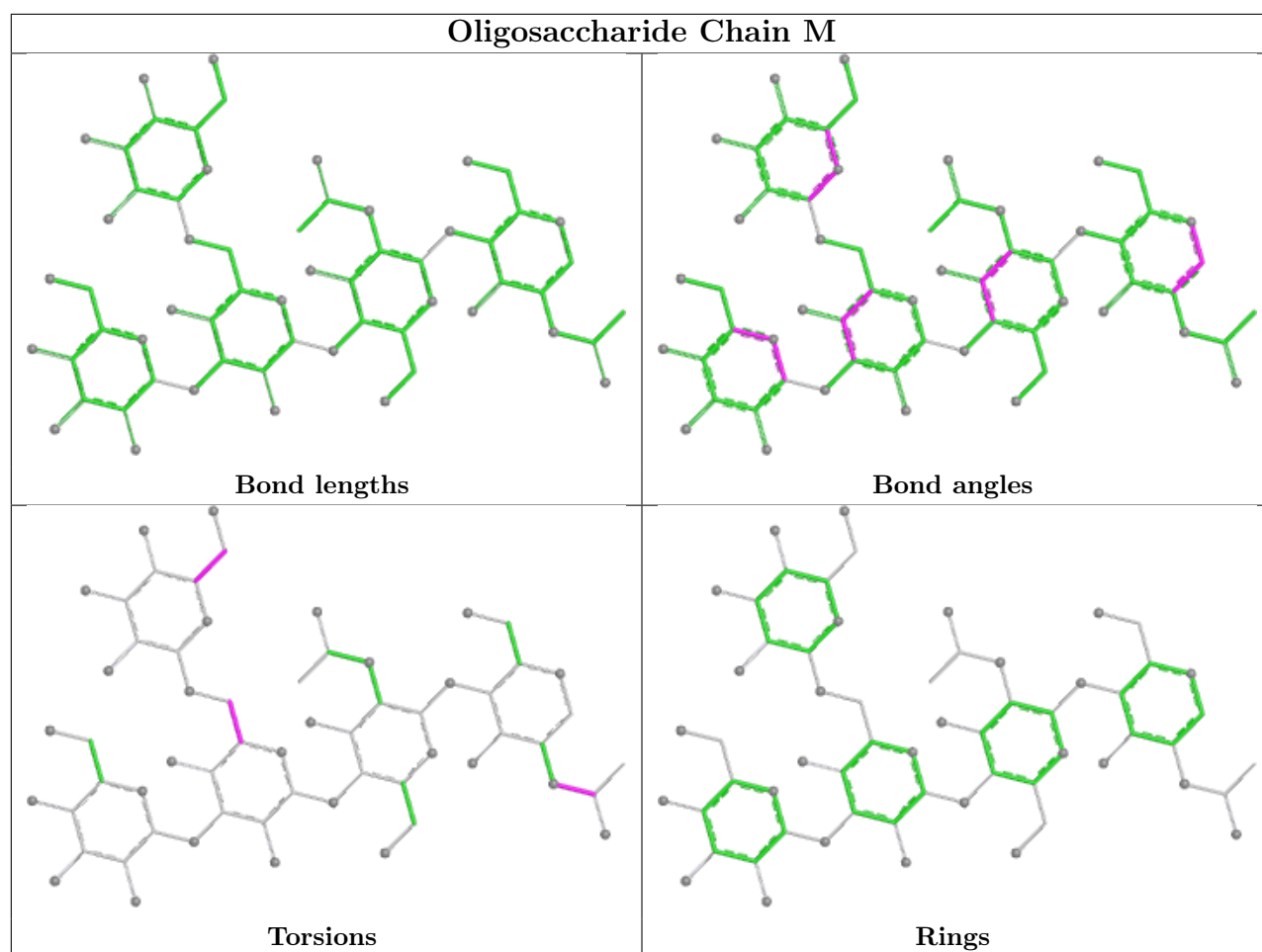
2 monomers are involved in 2 short contacts:

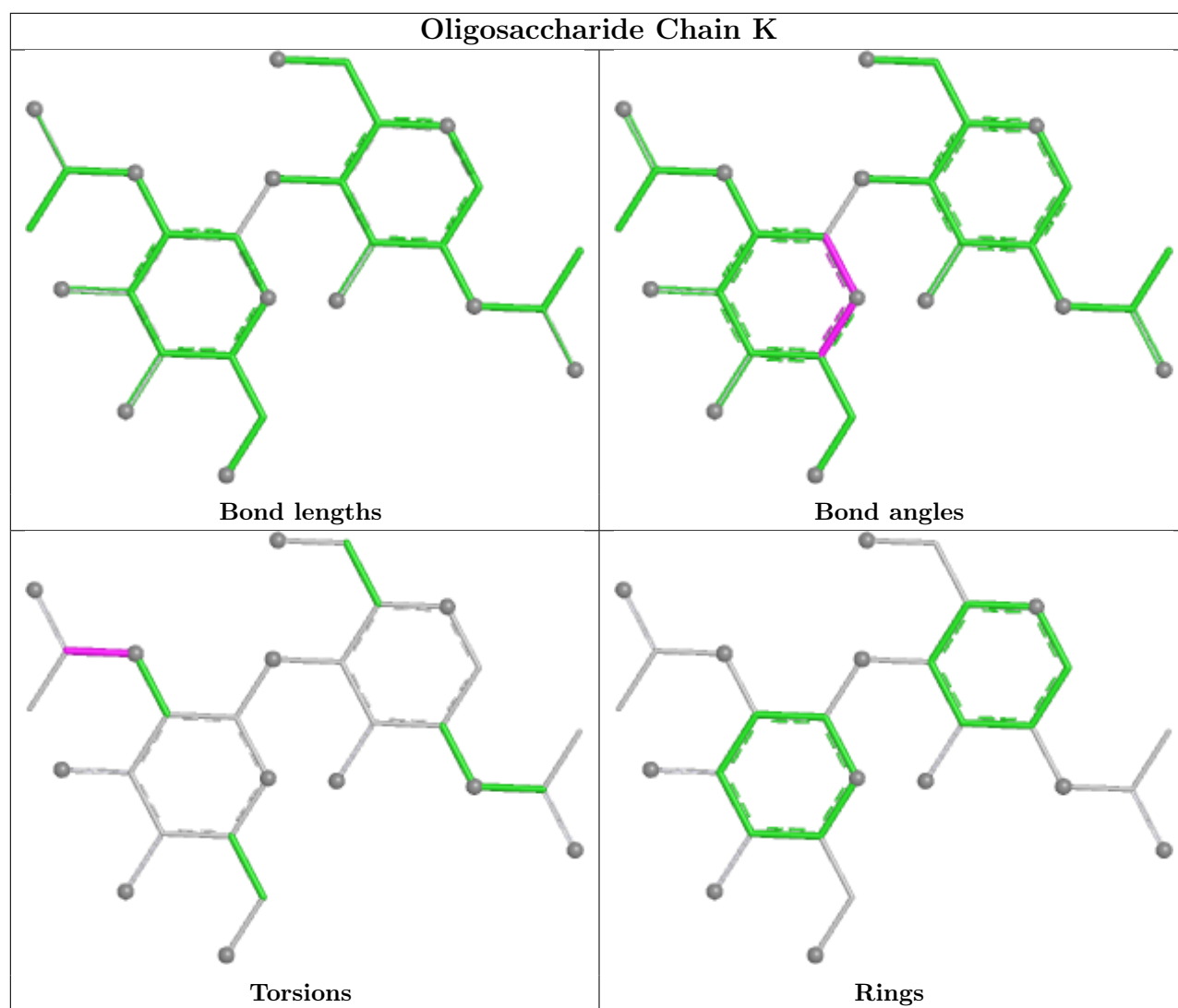
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	I	2	NAG	1	0
6	J	1	NAG	1	0

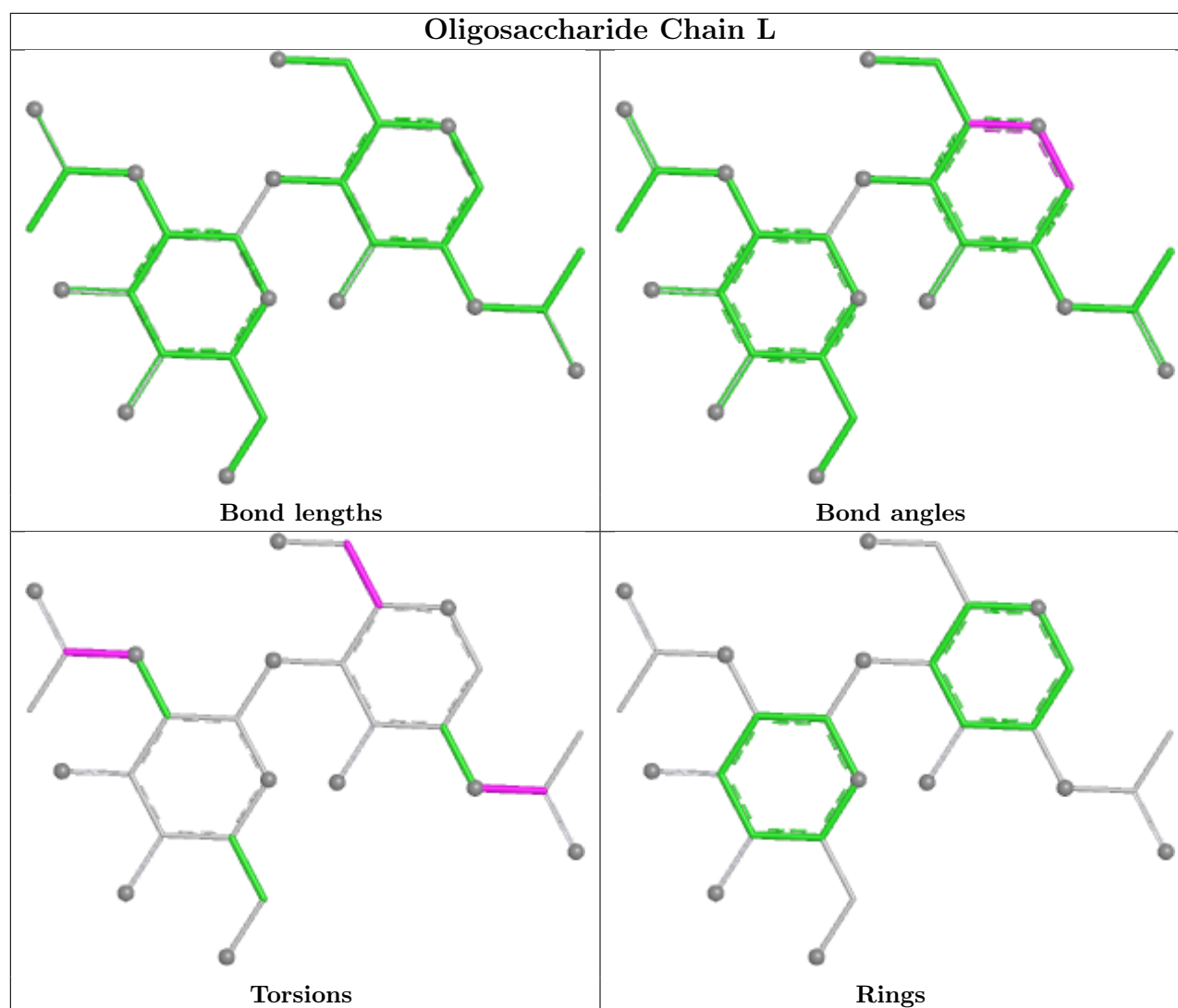
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

3 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	$\# Z > 2$
8	NAG	A	902	1	14,14,15	0.48	0	17,19,21	1.04	1 (5%)
8	NAG	A	901	1	14,14,15	0.46	0	17,19,21	0.78	0
8	NAG	E	905	1	14,14,15	0.31	0	17,19,21	0.75	1 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	A	902	1	-	2/6/23/26	0/1/1/1
8	NAG	A	901	1	-	0/6/23/26	0/1/1/1
8	NAG	E	905	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	902	NAG	C1-O5-C5	3.12	116.36	112.19
8	E	905	NAG	C1-O5-C5	2.49	115.53	112.19

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	902	NAG	C8-C7-N2-C2
8	A	902	NAG	O7-C7-N2-C2

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	729/833 (87%)	-0.28	9 (1%) 76 57	25, 58, 145, 178	0
1	E	705/833 (84%)	0.21	27 (3%) 44 26	55, 112, 159, 213	0
2	B	132/138 (95%)	-0.09	1 (0%) 82 66	53, 94, 138, 198	0
2	F	130/138 (94%)	-0.10	1 (0%) 82 66	67, 93, 118, 140	0
3	C	207/215 (96%)	0.06	8 (3%) 43 25	52, 99, 201, 212	0
3	G	196/215 (91%)	0.05	6 (3%) 51 31	58, 94, 211, 221	0
4	D	207/282 (73%)	0.00	4 (1%) 66 45	52, 116, 184, 201	0
4	H	206/282 (73%)	0.08	2 (0%) 79 62	68, 126, 196, 205	0
All	All	2512/2936 (85%)	-0.02	58 (2%) 61 40	25, 96, 182, 221	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	68	TRP	6.9
1	A	41	TYR	4.6
1	E	49	LYS	4.1
3	C	120	PRO	4.1
1	E	104	TYR	4.0
3	C	119	PHE	3.8
1	A	100	LEU	3.6
1	A	36	ALA	3.4
1	A	68	TRP	3.4
3	G	136	LEU	3.4
3	C	149	TRP	3.0
2	F	45	ILE	2.9
1	E	715	LEU	2.9
1	A	43	ASP	2.9
1	E	41	TYR	2.9
1	E	462	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
3	C	192	VAL	2.8
1	E	468	LEU	2.6
3	C	136	LEU	2.6
1	E	764	PRO	2.5
3	G	116	VAL	2.5
1	E	781	PHE	2.5
4	H	149	VAL	2.5
4	D	186	SER	2.4
4	H	166	LEU	2.4
3	C	133	VAL	2.4
1	E	174	PHE	2.4
1	E	179	LEU	2.4
3	G	103	LYS	2.4
1	E	705	SER	2.4
4	D	130	PRO	2.4
4	D	144	ALA	2.3
1	E	631	ALA	2.3
4	D	148	LEU	2.3
1	E	758	MET	2.3
1	E	59	ASP	2.2
1	E	714	ALA	2.2
3	G	49	TYR	2.2
1	E	358	ASP	2.2
1	E	452	PHE	2.2
1	E	637	LEU	2.2
1	E	639	THR	2.2
1	A	73	LYS	2.2
2	B	48	PRO	2.2
1	E	653	ARG	2.2
1	E	124	HIS	2.1
3	C	134	VAL	2.1
3	G	134	VAL	2.1
1	E	712	THR	2.1
1	A	587	ILE	2.1
1	E	704	VAL	2.1
3	C	116	VAL	2.1
3	G	117	PHE	2.1
1	E	743	ILE	2.0
1	E	56	THR	2.0
1	E	493	LEU	2.0
1	A	443	GLN	2.0
1	A	101	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

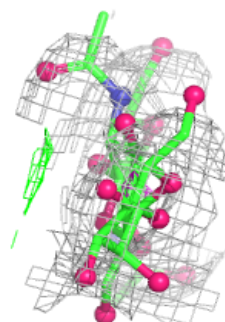
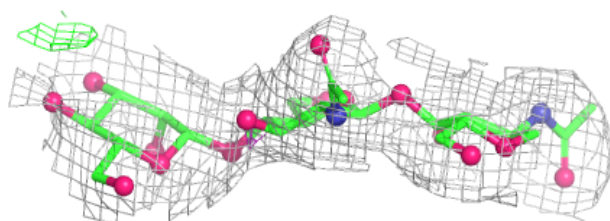
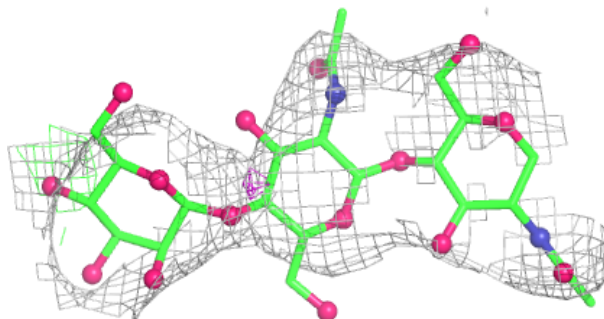
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	MAN	J	4	11/12	0.42	0.10	140,150,161,166	0
5	BMA	I	3	11/12	0.50	0.10	136,150,161,163	0
6	MAN	M	4	11/12	0.50	0.07	120,138,144,146	0
6	MAN	J	5	11/12	0.53	0.11	138,153,163,169	0
7	NAG	L	2	14/15	0.64	0.11	132,150,160,160	0
6	NAG	J	1	14/15	0.67	0.14	181,206,217,218	0
7	NAG	L	1	14/15	0.68	0.16	139,150,156,157	0
7	NAG	K	1	14/15	0.72	0.11	125,141,149,152	0
6	BMA	M	3	11/12	0.76	0.07	118,127,138,151	0
6	MAN	M	5	11/12	0.79	0.08	101,131,141,149	0
6	NAG	J	2	14/15	0.80	0.09	119,134,149,160	0
7	NAG	K	2	14/15	0.80	0.08	133,145,153,157	0
5	NAG	I	2	14/15	0.82	0.12	104,130,144,148	0
6	NAG	M	2	14/15	0.84	0.11	86,112,133,133	0
6	NAG	M	1	14/15	0.84	0.06	95,104,111,111	0
6	BMA	J	3	11/12	0.87	0.06	131,144,153,158	0
5	NAG	I	1	14/15	0.93	0.06	78,95,114,115	0

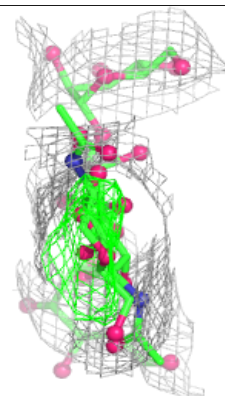
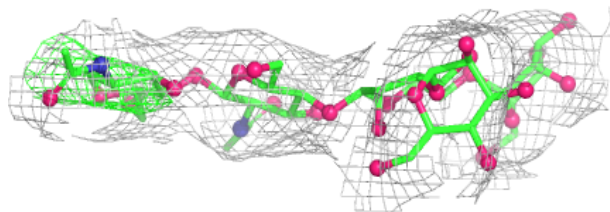
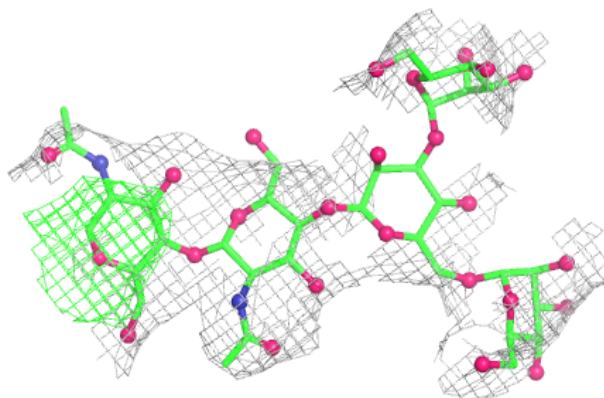
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

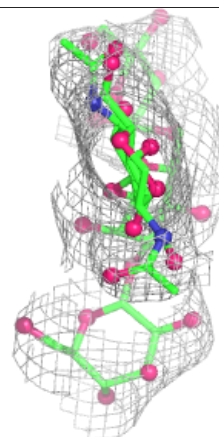
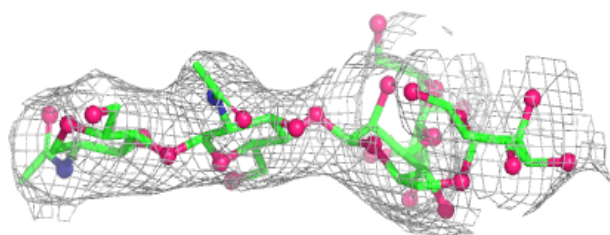
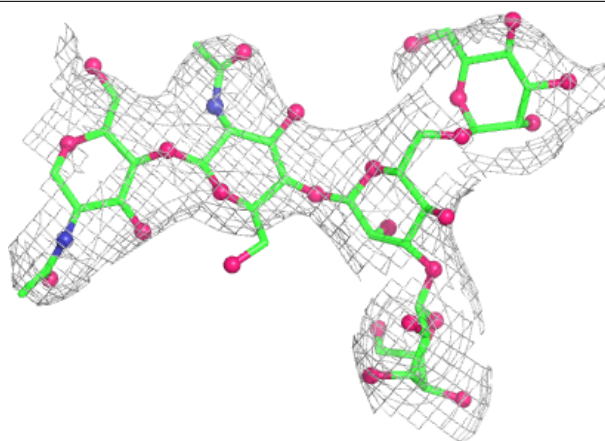
**Electron density around Chain J:**

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



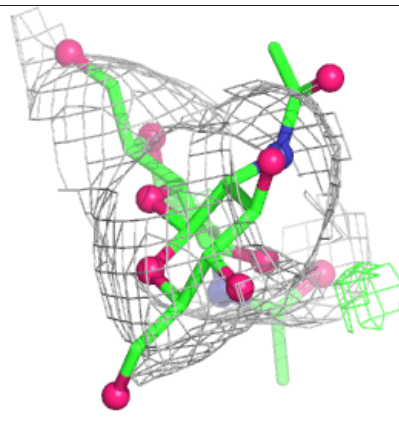
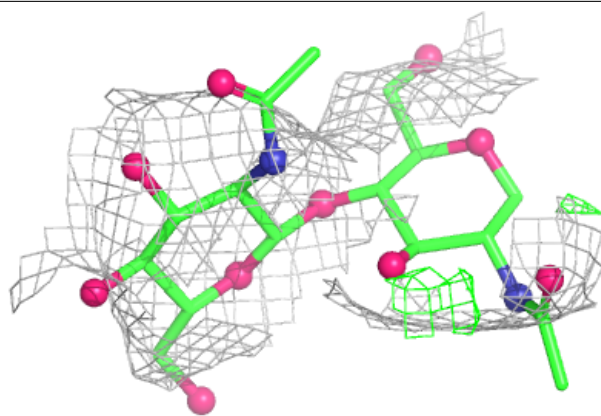
Electron density around Chain M:

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

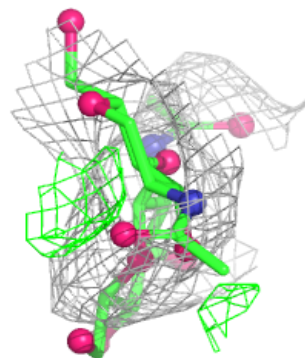
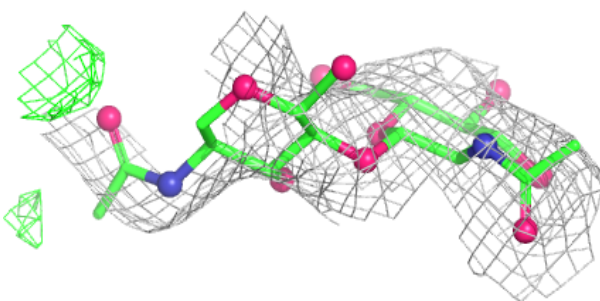
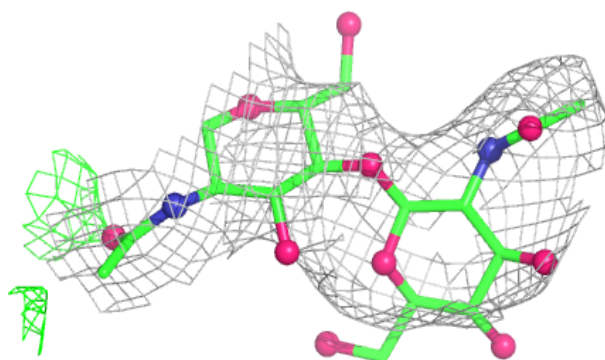


Electron density around Chain K:

$2mF_o-DF_c$ (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)



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
8	NAG	E	905	14/15	0.60	0.18	197,219,232,234	0
8	NAG	A	901	14/15	0.68	0.09	125,133,142,148	0
8	NAG	A	902	14/15	0.77	0.11	20,63,79,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.