



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:29 AM UTC

PDB ID : 6S9J / pdb_00006s9j
Title : Crystal structure of Tfr1 mimicry in complex with GP1 from MACV
Authors : Diskin, R.; Cohen-Dvashi, H.
Deposited on : 2019-07-15
Resolution : 2.70 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 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

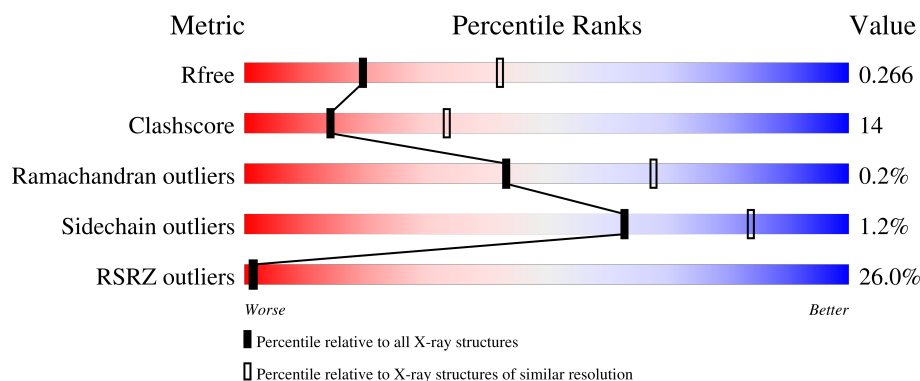
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	B	168	4% 71% 21% 8%
1	D	168	12% 71% 18% 10%
1	F	168	55% 53% 36% 11%
1	H	168	46% 60% 31% 8%
2	A	178	11% 65% 20% 15%

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Mol	Chain	Length	Quality of chain
2	C	178	
2	E	178	
2	G	178	
3	I	5	
4	J	2	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9715 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 Pre-glycoprotein polyprotein GP complex.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	155	Total	C	N	O	S	0	0	0
			1247	785	216	232	14			
1	D	152	Total	C	N	O	S	0	0	0
			1218	766	211	227	14			
1	F	149	Total	C	N	O	S	0	0	0
			1192	750	207	221	14			
1	H	154	Total	C	N	O	S	0	0	0
			1238	780	215	229	14			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	245	GLY	-	expression tag	UNP Q8AZ57
B	246	SER	-	expression tag	UNP Q8AZ57
B	247	HIS	-	expression tag	UNP Q8AZ57
B	248	HIS	-	expression tag	UNP Q8AZ57
B	249	HIS	-	expression tag	UNP Q8AZ57
B	250	HIS	-	expression tag	UNP Q8AZ57
B	251	HIS	-	expression tag	UNP Q8AZ57
B	252	HIS	-	expression tag	UNP Q8AZ57
D	245	GLY	-	expression tag	UNP Q8AZ57
D	246	SER	-	expression tag	UNP Q8AZ57
D	247	HIS	-	expression tag	UNP Q8AZ57
D	248	HIS	-	expression tag	UNP Q8AZ57
D	249	HIS	-	expression tag	UNP Q8AZ57
D	250	HIS	-	expression tag	UNP Q8AZ57
D	251	HIS	-	expression tag	UNP Q8AZ57
D	252	HIS	-	expression tag	UNP Q8AZ57
F	245	GLY	-	expression tag	UNP Q8AZ57
F	246	SER	-	expression tag	UNP Q8AZ57
F	247	HIS	-	expression tag	UNP Q8AZ57
F	248	HIS	-	expression tag	UNP Q8AZ57
F	249	HIS	-	expression tag	UNP Q8AZ57

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Chain	Residue	Modelled	Actual	Comment	Reference
F	250	HIS	-	expression tag	UNP Q8AZ57
F	251	HIS	-	expression tag	UNP Q8AZ57
F	252	HIS	-	expression tag	UNP Q8AZ57
H	245	GLY	-	expression tag	UNP Q8AZ57
H	246	SER	-	expression tag	UNP Q8AZ57
H	247	HIS	-	expression tag	UNP Q8AZ57
H	248	HIS	-	expression tag	UNP Q8AZ57
H	249	HIS	-	expression tag	UNP Q8AZ57
H	250	HIS	-	expression tag	UNP Q8AZ57
H	251	HIS	-	expression tag	UNP Q8AZ57
H	252	HIS	-	expression tag	UNP Q8AZ57

- Molecule 2 is a protein called Transferrin receptor 1, Transferrin receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	151	Total	C	N	O	S	0	0	0
			1170	737	201	229	3			
2	C	149	Total	C	N	O	S	0	0	0
			1159	731	199	226	3			
2	E	149	Total	C	N	O	S	0	0	0
			1159	731	199	226	3			
2	G	149	Total	C	N	O	S	0	0	0
			1159	731	199	226	3			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	194	CYS	GLY	conflict	UNP A0A060BIS8
A	?	-	VAL	deletion	UNP A0A060BIS8
A	283	LYS	MET	conflict	UNP A0A060BIS8
A	288	TYR	PHE	conflict	UNP A0A060BIS8
A	291	SER	VAL	conflict	UNP A0A060BIS8
A	295	GLU	ILE	conflict	UNP A0A060BIS8
A	324	GLN	PHE	conflict	UNP A0A060BIS8
A	380	CYS	ALA	conflict	UNP A0A060BIS8
A	384	GLY	-	expression tag	UNP A0A060BIS8
A	385	THR	-	expression tag	UNP A0A060BIS8
A	386	SER	-	expression tag	UNP A0A060BIS8
A	387	GLY	-	expression tag	UNP A0A060BIS8
A	388	HIS	-	expression tag	UNP A0A060BIS8
A	389	HIS	-	expression tag	UNP A0A060BIS8
A	390	HIS	-	expression tag	UNP A0A060BIS8

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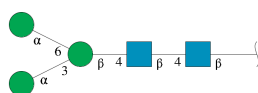
Chain	Residue	Modelled	Actual	Comment	Reference
A	391	HIS	-	expression tag	UNP A0A060BIS8
A	392	HIS	-	expression tag	UNP A0A060BIS8
A	393	HIS	-	expression tag	UNP A0A060BIS8
C	194	CYS	GLY	conflict	UNP A0A060BIS8
C	?	-	VAL	deletion	UNP A0A060BIS8
C	283	LYS	MET	conflict	UNP A0A060BIS8
C	288	TYR	PHE	conflict	UNP A0A060BIS8
C	291	SER	VAL	conflict	UNP A0A060BIS8
C	295	GLU	ILE	conflict	UNP A0A060BIS8
C	324	GLN	PHE	conflict	UNP A0A060BIS8
C	380	CYS	ALA	conflict	UNP A0A060BIS8
C	384	GLY	-	expression tag	UNP A0A060BIS8
C	385	THR	-	expression tag	UNP A0A060BIS8
C	386	SER	-	expression tag	UNP A0A060BIS8
C	387	GLY	-	expression tag	UNP A0A060BIS8
C	388	HIS	-	expression tag	UNP A0A060BIS8
C	389	HIS	-	expression tag	UNP A0A060BIS8
C	390	HIS	-	expression tag	UNP A0A060BIS8
C	391	HIS	-	expression tag	UNP A0A060BIS8
C	392	HIS	-	expression tag	UNP A0A060BIS8
C	393	HIS	-	expression tag	UNP A0A060BIS8
E	194	CYS	GLY	conflict	UNP A0A060BIS8
E	?	-	VAL	deletion	UNP A0A060BIS8
E	283	LYS	MET	conflict	UNP A0A060BIS8
E	288	TYR	PHE	conflict	UNP A0A060BIS8
E	291	SER	VAL	conflict	UNP A0A060BIS8
E	295	GLU	ILE	conflict	UNP A0A060BIS8
E	324	GLN	PHE	conflict	UNP A0A060BIS8
E	380	CYS	ALA	conflict	UNP A0A060BIS8
E	384	GLY	-	expression tag	UNP A0A060BIS8
E	385	THR	-	expression tag	UNP A0A060BIS8
E	386	SER	-	expression tag	UNP A0A060BIS8
E	387	GLY	-	expression tag	UNP A0A060BIS8
E	388	HIS	-	expression tag	UNP A0A060BIS8
E	389	HIS	-	expression tag	UNP A0A060BIS8
E	390	HIS	-	expression tag	UNP A0A060BIS8
E	391	HIS	-	expression tag	UNP A0A060BIS8
E	392	HIS	-	expression tag	UNP A0A060BIS8
E	393	HIS	-	expression tag	UNP A0A060BIS8
G	194	CYS	GLY	conflict	UNP A0A060BIS8
G	?	-	VAL	deletion	UNP A0A060BIS8
G	283	LYS	MET	conflict	UNP A0A060BIS8

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Chain	Residue	Modelled	Actual	Comment	Reference
G	288	TYR	PHE	conflict	UNP A0A060BIS8
G	291	SER	VAL	conflict	UNP A0A060BIS8
G	295	GLU	ILE	conflict	UNP A0A060BIS8
G	324	GLN	PHE	conflict	UNP A0A060BIS8
G	380	CYS	ALA	conflict	UNP A0A060BIS8
G	384	GLY	-	expression tag	UNP A0A060BIS8
G	385	THR	-	expression tag	UNP A0A060BIS8
G	386	SER	-	expression tag	UNP A0A060BIS8
G	387	GLY	-	expression tag	UNP A0A060BIS8
G	388	HIS	-	expression tag	UNP A0A060BIS8
G	389	HIS	-	expression tag	UNP A0A060BIS8
G	390	HIS	-	expression tag	UNP A0A060BIS8
G	391	HIS	-	expression tag	UNP A0A060BIS8
G	392	HIS	-	expression tag	UNP A0A060BIS8
G	393	HIS	-	expression tag	UNP A0A060BIS8

- 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	I	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 4 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
4	J	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



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

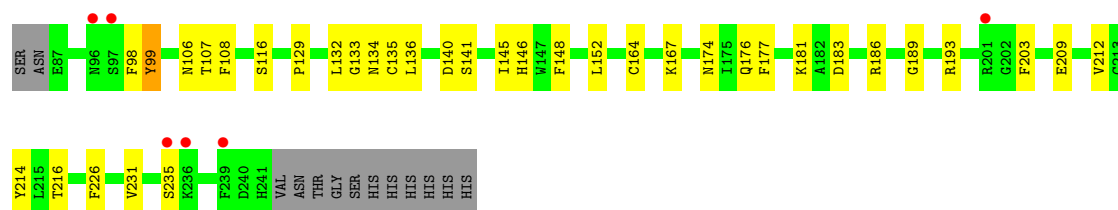
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	8	Total	O	0	0
			8	8		
6	A	8	Total	O	0	0
			8	8		
6	D	2	Total	O	0	0
			2	2		
6	C	9	Total	O	0	0
			9	9		
6	F	5	Total	O	0	0
			5	5		
6	E	6	Total	O	0	0
			6	6		
6	H	2	Total	O	0	0
			2	2		
6	G	2	Total	O	0	0
			2	2		

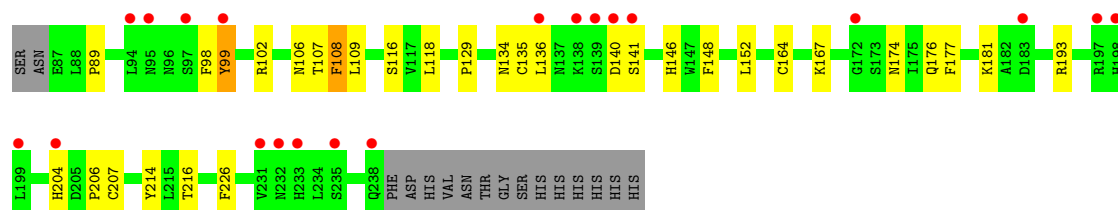
3 Residue-property plots

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.

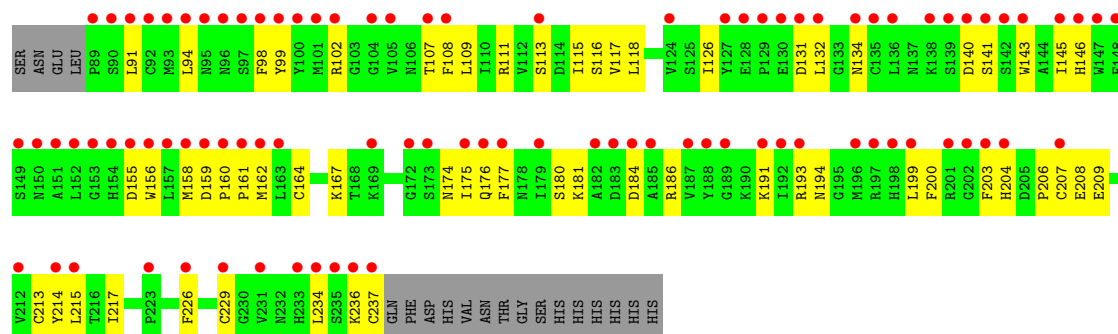
- Molecule 1: Pre-glycoprotein polyprotein GP complex



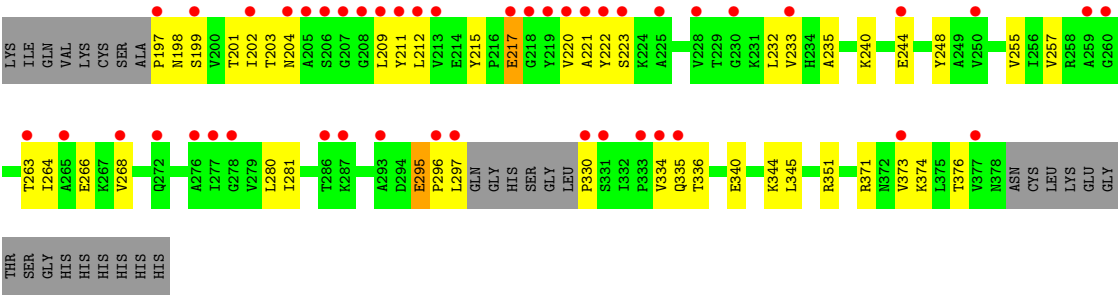
- Molecule 1: Pre-glycoprotein polyprotein GP complex



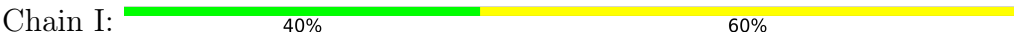
- Molecule 1: Pre-glycoprotein polyprotein GP complex



- Molecule 1: Pre-glycoprotein polyprotein GP complex



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, α , β , γ	104.63Å 104.63Å 281.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.56 – 2.70 49.56 – 2.70	Depositor EDS
% Data completeness (in resolution range)	83.1 (49.56-2.70) 83.1 (49.56-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.69Å)	Xtriage
Refinement program	PHENIX (dev_2871: ???)	Depositor
R, R_{free}	0.241 , 0.267 0.241 , 0.266	Depositor DCC
R_{free} test set	2000 reflections (4.54%)	wwPDB-VP
Wilson B-factor (Å ²)	43.4	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 57.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	9715	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.82% 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, BMA, MAN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	B	0.30	0/1279	0.54	0/1724
1	D	0.28	0/1248	0.49	0/1682
1	F	0.24	0/1222	0.54	0/1646
1	H	0.28	0/1270	0.58	1/1712 (0.1%)
2	A	0.35	0/1189	0.58	1/1611 (0.1%)
2	C	0.27	0/1178	0.56	0/1595
2	E	0.27	0/1178	0.56	0/1595
2	G	0.26	0/1178	0.73	2/1595 (0.1%)
All	All	0.28	0/9742	0.57	4/13160 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	295	GLU	CA-C-N	10.97	131.06	120.31
2	G	295	GLU	C-N-CA	10.97	131.06	120.31
1	H	183	ASP	CB-CA-C	-5.61	110.09	116.54
2	A	224	LYS	CA-CB-CG	-5.28	103.54	114.10

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	B	1247	0	1176	27	0
1	D	1218	0	1158	26	0
1	F	1192	0	1135	51	0
1	H	1238	0	1171	57	0
2	A	1170	0	1172	21	0
2	C	1159	0	1164	20	0
2	E	1159	0	1164	36	0
2	G	1159	0	1164	52	0
3	I	61	0	52	1	0
4	J	28	0	25	0	0
5	A	14	0	13	0	0
5	B	28	0	26	0	0
6	A	8	0	0	0	0
6	B	8	0	0	0	0
6	C	9	0	0	1	0
6	D	2	0	0	0	0
6	E	6	0	0	1	0
6	F	5	0	0	2	0
6	G	2	0	0	0	0
6	H	2	0	0	1	0
All	All	9715	0	9420	271	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (271) 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:F:155:ASP:OD2	2:G:221:ALA:HB2	1.36	1.25
1:H:181:LYS:HE3	1:H:193:ARG:NH1	1.51	1.22
2:G:220:VAL:HG11	2:G:335:GLN:NE2	1.61	1.13
1:H:181:LYS:NZ	1:H:207:CYS:O	1.82	1.11
1:F:181:LYS:HE2	1:F:193:ARG:HH12	1.01	1.11
1:H:181:LYS:NZ	1:H:207:CYS:C	2.11	1.07
1:F:155:ASP:OD2	2:G:221:ALA:CB	2.03	1.06
1:H:181:LYS:CE	1:H:193:ARG:NH1	2.24	1.00
2:G:220:VAL:HG11	2:G:335:GLN:CD	1.88	0.97
2:G:220:VAL:CG1	2:G:335:GLN:NE2	2.28	0.96
1:F:155:ASP:CG	2:G:221:ALA:HB2	1.90	0.95
1:F:181:LYS:HE2	1:F:193:ARG:NH1	1.84	0.92
1:H:181:LYS:HE3	1:H:193:ARG:HH11	1.26	0.92
2:G:220:VAL:CG1	2:G:335:GLN:HE22	1.82	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:181:LYS:O	1:H:186:ARG:NH2	2.02	0.90
1:H:181:LYS:CE	1:H:193:ARG:HH12	1.87	0.87
2:E:217:GLU:N	2:E:217:GLU:OE1	2.07	0.85
1:F:94:LEU:HD21	1:F:236:LYS:HE3	1.60	0.83
1:D:129:PRO:HB3	1:D:146:HIS:HD2	1.44	0.81
1:H:181:LYS:HZ2	1:H:207:CYS:C	1.83	0.81
2:G:297:LEU:O	2:G:297:LEU:HD12	1.80	0.81
1:H:181:LYS:CD	1:H:193:ARG:HH12	1.94	0.79
2:C:203:THR:HG22	2:C:209:LEU:HG	1.66	0.77
2:E:262:ILE:HB	2:E:266:GLU:OE1	1.86	0.76
2:E:198:ASN:ND2	2:E:220:VAL:HG21	2.00	0.75
1:H:181:LYS:HZ2	1:H:207:CYS:CB	1.99	0.74
2:G:198:ASN:ND2	2:G:220:VAL:CG2	2.51	0.73
1:B:129:PRO:HB3	1:B:146:HIS:CD2	2.24	0.73
2:G:220:VAL:HG11	2:G:335:GLN:HE22	1.42	0.73
1:H:181:LYS:NZ	1:H:207:CYS:HB2	2.04	0.72
2:E:203:THR:HG22	2:E:209:LEU:HG	1.71	0.72
1:H:118:LEU:H	1:H:176:GLN:HE22	1.36	0.72
2:G:204:ASN:HA	2:G:371:ARG:HG2	1.71	0.72
2:G:203:THR:HG22	2:G:209:LEU:HG	1.72	0.71
2:E:199:SER:HG	2:E:376:THR:HG1	1.30	0.71
1:D:129:PRO:HB3	1:D:146:HIS:CD2	2.26	0.70
2:C:233:VAL:HG13	2:C:248:TYR:CE2	2.27	0.70
2:G:220:VAL:HG13	2:G:335:GLN:HE22	1.56	0.69
2:C:272:GLN:NE2	2:C:331:SER:OG	2.25	0.69
2:A:233:VAL:HG13	2:A:248:TYR:CE2	2.28	0.69
2:G:281:ILE:HB	2:G:336:THR:HG22	1.75	0.68
1:H:180:SER:OG	1:H:212:VAL:HA	1.92	0.68
1:B:136:LEU:O	1:B:167:LYS:NZ	2.24	0.68
2:A:203:THR:HG22	2:A:209:LEU:HG	1.76	0.67
1:H:98:PHE:HD1	1:H:109:LEU:HD11	1.59	0.66
1:H:181:LYS:NZ	1:H:207:CYS:CB	2.57	0.66
1:B:129:PRO:HB3	1:B:146:HIS:HD2	1.61	0.65
2:G:295:GLU:HB3	2:G:296:PRO:HD2	1.79	0.65
2:C:201:THR:HG22	2:C:212:LEU:HA	1.78	0.65
2:E:198:ASN:HD21	2:E:220:VAL:HG21	1.61	0.65
1:F:177:PHE:HB2	1:F:215:LEU:HB3	1.79	0.65
1:F:98:PHE:HD1	1:F:109:LEU:HD11	1.61	0.64
1:H:180:SER:OG	1:H:213:CYS:N	2.29	0.64
1:F:158:MET:HE3	2:G:334:VAL:HG11	1.79	0.64
1:F:118:LEU:HB2	1:F:176:GLN:OE1	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:PHE:C	1:D:99:TYR:HD1	2.06	0.64
1:H:113:SER:HG	2:G:211:TYR:HH	1.46	0.64
2:A:258:ARG:NH1	2:A:284:ASP:OD2	2.28	0.63
1:D:116:SER:CB	1:D:176:GLN:HE22	2.11	0.62
2:A:199:SER:HG	2:A:376:THR:HG1	1.46	0.62
2:G:201:THR:HG22	2:G:212:LEU:HA	1.82	0.62
1:D:181:LYS:HE2	1:D:193:ARG:NH1	2.14	0.62
1:B:186:ARG:NH2	1:B:209:GLU:OE2	2.33	0.61
2:G:202:ILE:HD11	2:G:345:LEU:HD23	1.82	0.61
1:H:118:LEU:HB2	1:H:176:GLN:OE1	1.99	0.61
1:H:181:LYS:HE3	1:H:193:ARG:HH12	1.44	0.61
1:D:116:SER:HB2	1:D:176:GLN:HE22	1.66	0.60
2:E:202:ILE:HD11	2:E:345:LEU:HD23	1.82	0.60
2:G:217:GLU:N	2:G:217:GLU:OE1	2.34	0.60
2:E:201:THR:HG22	2:E:212:LEU:HA	1.84	0.60
2:A:202:ILE:HD11	2:A:345:LEU:HD23	1.84	0.60
2:E:268:VAL:HG11	2:E:334:VAL:HG21	1.84	0.59
2:G:233:VAL:HG13	2:G:248:TYR:CE2	2.37	0.59
2:A:201:THR:HG22	2:A:212:LEU:HA	1.83	0.59
2:G:232:LEU:HD22	2:G:373:VAL:HG21	1.84	0.59
1:H:111:ARG:NH2	1:H:218:ASN:OD1	2.33	0.59
2:E:261:LYS:N	6:E:401:HOH:O	2.33	0.59
2:C:283:LYS:HD3	2:C:287:LYS:HE3	1.85	0.58
1:B:181:LYS:HE2	1:B:193:ARG:HH12	1.68	0.58
1:B:107:THR:C	1:B:108:PHE:HD1	2.12	0.58
2:G:295:GLU:OE1	2:G:296:PRO:HD2	2.03	0.57
1:D:107:THR:C	1:D:108:PHE:HD1	2.12	0.57
2:C:243:PHE:HB3	2:C:274:PHE:CE2	2.39	0.57
1:H:123:ASP:OD2	2:G:344:LYS:NZ	2.38	0.57
1:F:98:PHE:CE1	1:F:111:ARG:HD3	2.40	0.57
1:F:155:ASP:OD2	2:G:221:ALA:HB3	2.00	0.57
2:A:243:PHE:HB3	2:A:274:PHE:CE2	2.41	0.56
1:B:189:GLY:C	1:B:193:ARG:HE	2.14	0.56
1:B:98:PHE:C	1:B:99:TYR:HD1	2.14	0.56
2:A:285:ARG:HH21	2:A:339:ARG:HH12	1.53	0.56
1:H:174:ASN:OD1	1:H:216:THR:HG23	2.06	0.56
1:H:181:LYS:CE	1:H:193:ARG:HH11	2.05	0.56
1:H:118:LEU:N	1:H:176:GLN:HE22	2.02	0.56
1:F:117:VAL:H	1:F:176:GLN:NE2	2.03	0.55
1:H:130:GLU:O	1:H:130:GLU:HG3	2.06	0.55
1:D:164:CYS:HA	1:D:174:ASN:O	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:181:LYS:CG	1:H:193:ARG:HH12	2.19	0.55
2:G:340:GLU:HG2	2:G:344:LYS:HE2	1.89	0.55
2:G:220:VAL:HG11	2:G:335:GLN:OE1	2.04	0.55
1:B:181:LYS:CE	1:B:193:ARG:HH12	2.20	0.55
2:A:209:LEU:C	2:A:211:TYR:HD1	2.15	0.54
2:A:243:PHE:HB3	2:A:274:PHE:CD2	2.43	0.54
1:F:134:ASN:HA	1:F:167:LYS:NZ	2.22	0.54
2:E:232:LEU:HD22	2:E:373:VAL:HG21	1.89	0.54
2:C:227:GLU:OE2	2:C:374:LYS:HE3	2.08	0.54
2:G:197:PRO:HG2	2:G:215:TYR:CE1	2.42	0.54
2:G:198:ASN:ND2	2:G:220:VAL:HG22	2.22	0.54
2:E:227:GLU:OE2	2:E:374:LYS:HE3	2.07	0.54
2:C:202:ILE:HD11	2:C:345:LEU:HD23	1.91	0.53
2:E:222:TYR:CD2	2:E:222:TYR:O	2.61	0.53
1:H:129:PRO:HB3	1:H:146:HIS:CD2	2.43	0.53
1:H:116:SER:HB3	1:H:214:TYR:CE2	2.44	0.53
1:F:118:LEU:N	1:F:176:GLN:HE22	2.07	0.53
2:C:209:LEU:C	2:C:211:TYR:HD1	2.17	0.53
1:F:140:ASP:OD1	1:F:141:SER:N	2.42	0.52
2:E:233:VAL:HG13	2:E:248:TYR:CE2	2.44	0.52
1:B:140:ASP:OD1	1:B:141:SER:N	2.42	0.52
1:B:164:CYS:HA	1:B:174:ASN:O	2.09	0.52
2:E:222:TYR:O	2:E:222:TYR:HD2	1.93	0.52
2:G:295:GLU:CB	2:G:296:PRO:HD2	2.39	0.52
2:E:237:PHE:O	2:E:267:LYS:NZ	2.39	0.52
1:H:117:VAL:H	1:H:176:GLN:NE2	2.08	0.52
1:H:181:LYS:HG2	1:H:193:ARG:HH12	1.74	0.52
1:F:132:LEU:HD13	1:F:145:ILE:HG22	1.92	0.51
1:F:143:TRP:O	1:F:146:HIS:HB2	2.10	0.51
1:D:116:SER:HB3	1:D:214:TYR:CE2	2.46	0.51
2:E:215:TYR:O	2:E:217:GLU:OE1	2.28	0.51
2:E:220:VAL:HG11	2:E:335:GLN:OE1	2.09	0.51
2:E:199:SER:HB3	2:E:215:TYR:HE1	1.75	0.51
1:H:140:ASP:OD1	1:H:141:SER:N	2.44	0.51
1:D:98:PHE:HD1	1:D:109:LEU:HD11	1.76	0.50
1:D:106:ASN:HB3	1:D:108:PHE:HE1	1.77	0.50
1:F:107:THR:C	1:F:108:PHE:HD1	2.20	0.50
1:H:107:THR:C	1:H:108:PHE:HD1	2.20	0.50
2:A:263:THR:HG23	2:A:266:GLU:H	1.76	0.50
1:F:164:CYS:HA	1:F:174:ASN:O	2.12	0.50
2:E:199:SER:HB3	2:E:215:TYR:CE1	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:135:CYS:HB2	1:H:136:LEU:HD12	1.93	0.50
2:G:201:THR:OG1	2:G:374:LYS:HB3	2.11	0.50
2:A:272:GLN:NE2	2:A:331:SER:OG	2.45	0.50
1:H:181:LYS:HZ1	1:H:207:CYS:C	1.98	0.50
1:B:106:ASN:HB3	1:B:108:PHE:HE1	1.77	0.49
1:D:140:ASP:OD1	1:D:141:SER:N	2.45	0.49
1:B:135:CYS:HB2	1:B:136:LEU:HD12	1.94	0.49
2:E:340:GLU:HG2	2:E:344:LYS:HE2	1.93	0.49
2:G:217:GLU:CD	2:G:217:GLU:H	2.21	0.49
1:H:118:LEU:HD11	1:H:174:ASN:O	2.12	0.49
1:B:183:ASP:OD2	3:I:2:NAG:H3	2.12	0.49
2:E:209:LEU:C	2:E:211:TYR:HD1	2.21	0.49
1:H:181:LYS:HG2	1:H:193:ARG:NH1	2.28	0.49
1:F:140:ASP:O	1:F:143:TRP:HD1	1.95	0.49
2:C:352:ASP:OD1	2:C:364:LYS:NZ	2.46	0.49
1:H:120:LYS:HG2	1:H:163:LEU:HD21	1.95	0.49
1:F:126:ILE:HG22	1:F:156:TRP:HZ3	1.77	0.48
1:H:154:HIS:CD2	1:H:162:MET:HE1	2.48	0.48
1:D:134:ASN:HA	1:D:167:LYS:HE3	1.95	0.48
1:H:115:ILE:HD13	2:G:202:ILE:HG22	1.96	0.48
2:C:263:THR:HG23	2:C:266:GLU:H	1.79	0.48
1:D:118:LEU:HD12	1:D:176:GLN:HB2	1.96	0.47
2:C:367:LEU:HD22	2:C:373:VAL:HG23	1.96	0.47
2:A:339:ARG:NH1	2:A:343:GLU:OE2	2.46	0.47
1:B:116:SER:HB3	1:B:214:TYR:CE2	2.49	0.47
1:F:186:ARG:NH2	1:F:209:GLU:OE2	2.47	0.47
1:H:117:VAL:O	2:G:344:LYS:HD3	2.15	0.47
1:F:155:ASP:O	1:F:159:ASP:HB2	2.15	0.47
1:F:118:LEU:H	1:F:176:GLN:HE22	1.61	0.47
1:D:135:CYS:HB2	1:D:136:LEU:HD12	1.97	0.47
2:C:240:LYS:O	2:C:244:GLU:HG3	2.15	0.47
2:A:227:GLU:OE2	2:A:374:LYS:HE3	2.15	0.47
1:D:204:HIS:O	1:D:206:PRO:HD3	2.15	0.47
1:F:126:ILE:HG23	1:F:131:ASP:HB2	1.97	0.46
2:E:344:LYS:O	2:E:347:GLN:HG2	2.15	0.46
2:G:199:SER:OG	2:G:376:THR:OG1	2.29	0.46
2:G:235:ALA:HB3	2:G:257:VAL:HG12	1.97	0.46
1:D:181:LYS:HE3	1:D:207:CYS:HB2	1.97	0.46
1:B:99:TYR:HE2	1:B:203:PHE:HD1	1.61	0.46
2:A:340:GLU:HG2	2:A:344:LYS:HE2	1.96	0.46
1:D:99:TYR:CD1	1:D:99:TYR:N	2.83	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:116:SER:HB3	1:F:214:TYR:CE2	2.50	0.46
2:G:297:LEU:HD12	2:G:297:LEU:C	2.41	0.46
1:H:226:PHE:HE2	2:G:209:LEU:HB3	1.81	0.46
1:F:181:LYS:HB2	1:F:208:GLU:O	2.16	0.46
1:B:212:VAL:HG12	1:B:214:TYR:CE2	2.51	0.46
1:F:194:ASN:ND2	6:F:302:HOH:O	2.35	0.46
1:F:226:PHE:CE2	2:E:209:LEU:HB3	2.51	0.46
1:H:164:CYS:HA	1:H:174:ASN:O	2.16	0.46
1:H:181:LYS:HZ2	1:H:207:CYS:HB3	1.78	0.46
2:G:209:LEU:C	2:G:211:TYR:HD1	2.24	0.46
1:F:134:ASN:HA	1:F:167:LYS:HZ1	1.79	0.45
1:D:99:TYR:HD1	1:D:99:TYR:N	2.13	0.45
1:H:116:SER:HB3	1:H:214:TYR:CZ	2.51	0.45
1:F:115:ILE:HD13	2:E:202:ILE:HG22	1.98	0.45
1:F:113:SER:OG	2:E:211:TYR:OH	2.26	0.45
2:E:262:ILE:HD12	2:E:266:GLU:HB3	1.97	0.45
1:H:98:PHE:C	1:H:99:TYR:HD1	2.25	0.45
1:F:99:TYR:HE2	1:F:203:PHE:HD1	1.65	0.45
1:H:126:ILE:HB	1:H:162:MET:SD	2.57	0.45
1:B:174:ASN:OD1	1:B:216:THR:HG23	2.17	0.45
1:F:158:MET:HE1	2:G:264:ILE:HG22	1.99	0.45
1:H:169:LYS:O	6:H:301:HOH:O	2.21	0.45
1:B:231:VAL:O	1:B:235:SER:CB	2.65	0.45
1:H:226:PHE:CE2	2:G:209:LEU:HB3	2.52	0.45
2:G:222:TYR:O	2:G:222:TYR:CD2	2.70	0.45
2:E:225:ALA:HA	2:E:377:VAL:O	2.17	0.45
1:H:89:PRO:HA	1:H:102:ARG:O	2.17	0.44
1:H:204:HIS:O	1:H:206:PRO:HD3	2.17	0.44
2:G:223:SER:OG	2:G:330:PRO:O	2.22	0.44
2:A:350:GLU:O	2:A:351:ARG:HB2	2.17	0.44
2:G:240:LYS:O	2:G:244:GLU:HG3	2.18	0.44
1:H:140:ASP:O	1:H:143:TRP:HD1	2.01	0.43
2:G:198:ASN:ND2	2:G:220:VAL:HG21	2.31	0.43
2:E:263:THR:O	2:E:266:GLU:HB2	2.18	0.43
1:H:155:ASP:HB3	1:H:159:ASP:CG	2.43	0.43
1:D:89:PRO:HA	1:D:102:ARG:O	2.19	0.43
1:D:118:LEU:HB2	1:D:176:GLN:OE1	2.18	0.43
2:E:216:PRO:HA	2:E:217:GLU:OE1	2.19	0.43
1:F:229:CYS:O	6:F:301:HOH:O	2.21	0.43
2:E:198:ASN:O	2:E:216:PRO:HD2	2.19	0.43
2:E:217:GLU:N	2:E:217:GLU:CD	2.74	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:226:PHE:CE2	2:C:209:LEU:HB3	2.53	0.43
1:F:98:PHE:C	1:F:99:TYR:HD1	2.26	0.43
1:F:126:ILE:O	1:F:161:PRO:HB3	2.18	0.43
1:H:181:LYS:CE	1:H:207:CYS:HB2	2.48	0.43
1:B:148:PHE:CE2	1:B:152:LEU:HD11	2.54	0.43
1:D:226:PHE:HE2	2:C:209:LEU:HB3	1.84	0.43
1:F:180:SER:HA	1:F:213:CYS:HB3	2.01	0.43
2:E:203:THR:CG2	2:E:209:LEU:HG	2.45	0.43
2:C:344:LYS:O	2:C:347:GLN:HG2	2.19	0.43
1:F:162:MET:HB2	1:F:176:GLN:O	2.19	0.43
1:H:186:ARG:NH2	1:H:209:GLU:HG2	2.34	0.43
1:H:229:CYS:HA	1:H:234:LEU:HD13	2.01	0.43
1:F:91:LEU:HB3	1:F:200:PHE:CE1	2.55	0.42
2:G:233:VAL:O	2:G:255:VAL:HA	2.19	0.42
2:G:263:THR:HG23	2:G:266:GLU:H	1.84	0.42
2:A:198:ASN:ND2	2:A:220:VAL:HG11	2.34	0.42
1:B:99:TYR:HE2	1:B:203:PHE:CD1	2.36	0.42
1:B:108:PHE:HE2	1:B:145:ILE:HG12	1.83	0.42
1:D:174:ASN:OD1	1:D:216:THR:HG23	2.19	0.42
2:C:243:PHE:HB3	2:C:274:PHE:CD2	2.55	0.42
1:F:91:LEU:HD13	1:F:199:LEU:HB3	2.01	0.42
1:B:133:GLY:O	1:B:167:LYS:HE2	2.20	0.42
1:B:226:PHE:N	1:B:226:PHE:CD1	2.87	0.42
1:F:99:TYR:HE2	1:F:203:PHE:CD1	2.38	0.41
1:F:184:ASP:OD1	2:G:297:LEU:HA	2.20	0.41
1:B:231:VAL:O	1:B:235:SER:HB3	2.20	0.41
2:C:339:ARG:NH1	6:C:403:HOH:O	2.53	0.41
1:F:102:ARG:HH22	1:F:234:LEU:HD23	1.84	0.41
2:E:295:GLU:CD	2:E:296:PRO:HD2	2.45	0.41
2:A:203:THR:CG2	2:A:209:LEU:HG	2.47	0.41
1:D:226:PHE:N	1:D:226:PHE:CD1	2.88	0.41
2:A:367:LEU:HD22	2:A:373:VAL:HG23	2.03	0.41
2:G:268:VAL:HG11	2:G:334:VAL:HG21	2.02	0.41
1:F:181:LYS:HG3	1:F:207:CYS:HB3	2.03	0.41
1:B:116:SER:HB2	1:B:176:GLN:HE22	1.86	0.41
2:A:233:VAL:O	2:A:255:VAL:HA	2.21	0.41
1:D:148:PHE:CE2	1:D:152:LEU:HD11	2.56	0.41
1:H:118:LEU:HA	2:G:344:LYS:NZ	2.36	0.41
1:H:165:ARG:HG3	1:H:169:LYS:HD3	2.02	0.41
2:E:367:LEU:HD22	2:E:373:VAL:HG23	2.03	0.41
2:A:259:ALA:HA	2:A:267:LYS:HE3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:242:ASP:O	2:C:246:LEU:HG	2.21	0.40
1:F:160:PRO:HA	1:F:161:PRO:HD3	1.99	0.40
1:F:159:ASP:OD2	1:F:191:LYS:NZ	2.40	0.40
1:F:175:ILE:HB	1:F:217:ILE:HB	2.03	0.40
1:F:204:HIS:O	1:F:206:PRO:HD3	2.22	0.40
1:F:226:PHE:N	1:F:226:PHE:CD1	2.90	0.40
2:G:203:THR:CG2	2:G:209:LEU:HG	2.48	0.40
2:G:280:LEU:HA	2:G:335:GLN:O	2.21	0.40
1:B:132:LEU:C	1:B:134:ASN:H	2.30	0.40
2:C:295:GLU:CD	2:C:296:PRO:HD2	2.47	0.40
2:E:264:ILE:HG23	2:E:281:ILE:CD1	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	153/168 (91%)	146 (95%)	7 (5%)	0	100	100
1	D	150/168 (89%)	145 (97%)	5 (3%)	0	100	100
1	F	147/168 (88%)	140 (95%)	7 (5%)	0	100	100
1	H	152/168 (90%)	148 (97%)	4 (3%)	0	100	100
2	A	147/178 (83%)	141 (96%)	5 (3%)	1 (1%)	18	41
2	C	145/178 (82%)	137 (94%)	8 (6%)	0	100	100
2	E	145/178 (82%)	136 (94%)	9 (6%)	0	100	100
2	G	145/178 (82%)	138 (95%)	6 (4%)	1 (1%)	18	41
All	All	1184/1384 (86%)	1131 (96%)	51 (4%)	2 (0%)	43	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	G	351	ARG
2	A	351	ARG

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	B	140/152 (92%)	138 (99%)	2 (1%)	59	82
1	D	137/152 (90%)	134 (98%)	3 (2%)	45	74
1	F	134/152 (88%)	133 (99%)	1 (1%)	76	90
1	H	139/152 (91%)	136 (98%)	3 (2%)	45	74
2	A	129/152 (85%)	129 (100%)	0	100	100
2	C	128/152 (84%)	128 (100%)	0	100	100
2	E	128/152 (84%)	125 (98%)	3 (2%)	44	73
2	G	128/152 (84%)	127 (99%)	1 (1%)	73	88
All	All	1063/1216 (87%)	1050 (99%)	13 (1%)	63	84

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	99	TYR
1	B	177	PHE
1	D	99	TYR
1	D	108	PHE
1	D	177	PHE
1	F	237	CYS
2	E	217	GLU
2	E	220	VAL
2	E	222	TYR
1	H	179	ILE
1	H	183	ASP
1	H	187	VAL
2	G	217	GLU

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

Mol	Chain	Res	Type
1	B	146	HIS
1	B	150	ASN
1	B	176	GLN
2	A	272	GLN
2	A	372	ASN
1	D	146	HIS
1	D	176	GLN
2	C	272	GLN
1	F	176	GLN
1	H	146	HIS
1	H	176	GLN
1	H	233	HIS
2	G	335	GLN
2	G	372	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 ⓘ

7 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	I	1	1,3	14,14,15	0.40	0	17,19,21	0.55	0
3	NAG	I	2	3	14,14,15	0.46	0	17,19,21	0.83	0
3	BMA	I	3	3	11,11,12	0.99	0	15,15,17	0.92	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MAN	I	4	3	11,11,12	1.40	1 (9%)	15,15,17	1.59	4 (26%)
3	MAN	I	5	3	11,11,12	0.77	0	15,15,17	0.79	0
4	NAG	J	1	1,4	14,14,15	0.49	0	17,19,21	0.62	1 (5%)
4	NAG	J	2	4	14,14,15	0.33	0	17,19,21	0.43	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	I	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1
3	BMA	I	3	3	-	2/2/19/22	0/1/1/1
3	MAN	I	4	3	-	0/2/19/22	0/1/1/1
3	MAN	I	5	3	-	2/2/19/22	0/1/1/1
4	NAG	J	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	4	MAN	C1-C2	4.16	1.62	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	4	MAN	C1-O5-C5	3.05	116.28	112.19
3	I	4	MAN	C1-C2-C3	2.97	113.97	109.64
3	I	4	MAN	O2-C2-C3	-2.26	105.47	110.15
3	I	4	MAN	C3-C4-C5	-2.16	106.31	110.23
4	J	1	NAG	C1-O5-C5	2.04	114.93	112.19
3	I	3	BMA	C3-C4-C5	2.01	113.89	110.23

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	I	5	MAN	O5-C5-C6-O6
3	I	5	MAN	C4-C5-C6-O6

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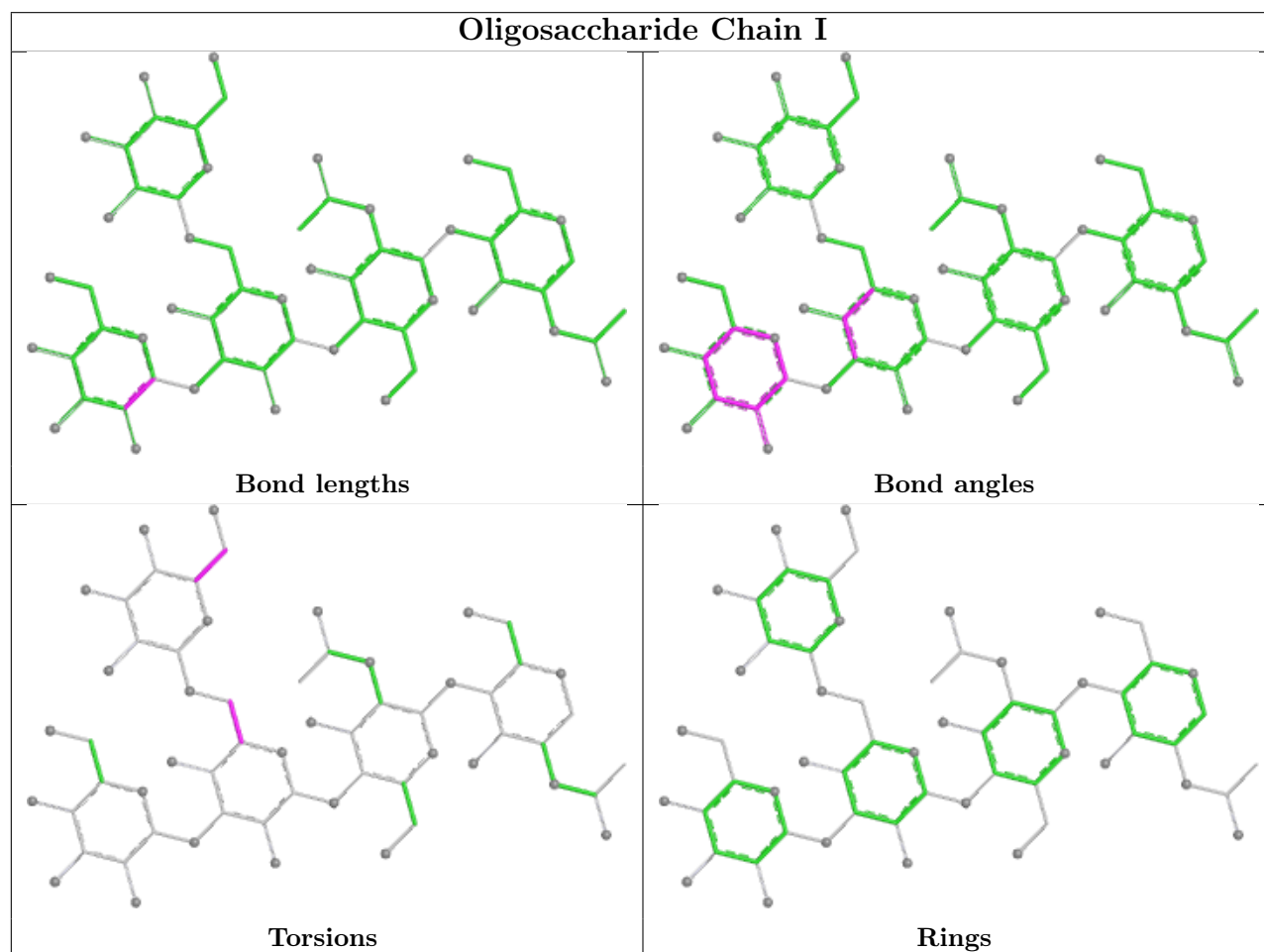
Mol	Chain	Res	Type	Atoms
3	I	3	BMA	C4-C5-C6-O6
3	I	3	BMA	O5-C5-C6-O6

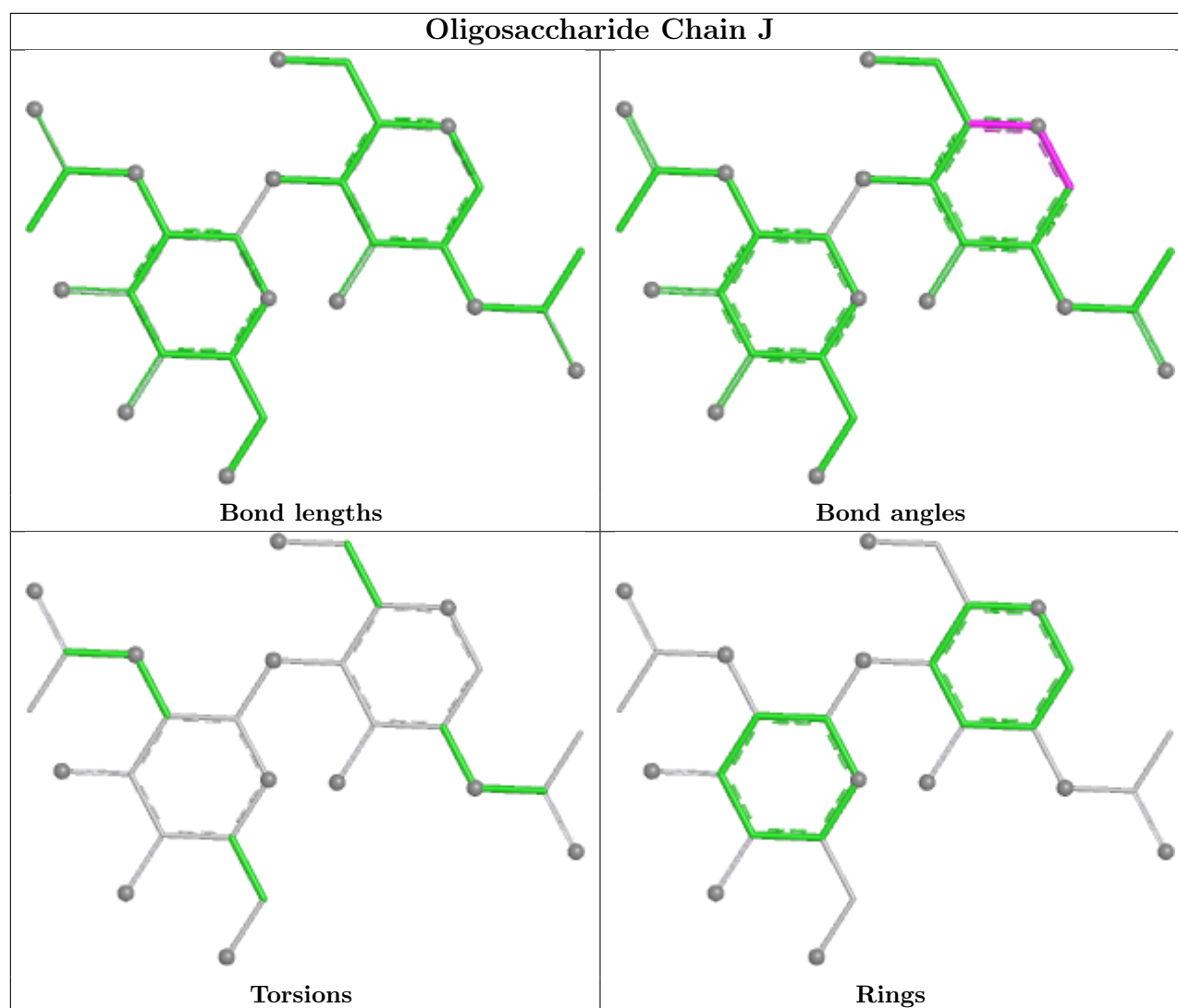
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	2	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	$\# Z > 2$
5	NAG	B	307	1	14,14,15	0.49	0	17,19,21	0.70	1 (5%)
5	NAG	B	306	1	14,14,15	0.46	0	17,19,21	0.67	1 (5%)
5	NAG	A	401	2	14,14,15	0.56	0	17,19,21	0.66	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
5	NAG	B	307	1	-	2/6/23/26	0/1/1/1
5	NAG	B	306	1	-	2/6/23/26	0/1/1/1
5	NAG	A	401	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	307	NAG	C1-O5-C5	2.55	115.61	112.19
5	B	306	NAG	C1-O5-C5	2.34	115.32	112.19
5	A	401	NAG	C1-O5-C5	2.25	115.20	112.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	306	NAG	O5-C5-C6-O6
5	A	401	NAG	O5-C5-C6-O6
5	B	306	NAG	C4-C5-C6-O6
5	A	401	NAG	C4-C5-C6-O6
5	B	307	NAG	O5-C5-C6-O6
5	B	307	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	B	155/168 (92%)	0.43	6 (3%) 43 39	19, 47, 98, 149	0
1	D	152/168 (90%)	0.75	20 (13%) 7 6	22, 57, 127, 214	0
1	F	149/168 (88%)	2.32	92 (61%) 0 0	55, 136, 174, 202	0
1	H	154/168 (91%)	2.16	78 (50%) 0 0	67, 129, 174, 223	0
2	A	151/178 (84%)	0.44	20 (13%) 7 6	13, 34, 114, 153	0
2	C	149/178 (83%)	0.54	19 (12%) 7 6	19, 44, 139, 182	0
2	E	149/178 (83%)	0.95	33 (22%) 2 2	22, 61, 159, 201	0
2	G	149/178 (83%)	1.43	46 (30%) 1 1	46, 84, 153, 184	0
All	All	1208/1384 (87%)	1.12	314 (25%) 1 1	13, 68, 165, 223	0

All (314) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	330	PRO	6.3
2	E	219	TYR	5.8
2	A	297	LEU	5.4
1	F	156	TRP	5.3
2	C	330	PRO	5.3
2	E	331	SER	5.3
1	H	105	VAL	5.3
1	H	97	SER	5.2
1	F	157	LEU	5.2
1	F	97	SER	5.1
1	F	237	CYS	4.9
1	F	147	TRP	4.9
1	H	147	TRP	4.9
1	H	98	PHE	4.8
2	C	293	ALA	4.8
2	C	207	GLY	4.8

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Mol	Chain	Res	Type	RSRZ
2	A	222	TYR	4.7
1	H	151	ALA	4.7
2	G	205	ALA	4.6
1	H	88	LEU	4.6
2	C	206	SER	4.6
1	H	148	PHE	4.6
1	H	92	CYS	4.6
2	A	207	GLY	4.6
2	C	297	LEU	4.5
2	C	296	PRO	4.5
1	F	108	PHE	4.4
2	A	293	ALA	4.4
1	F	185	ALA	4.3
2	E	197	PRO	4.2
1	H	94	LEU	4.2
1	F	160	PRO	4.2
2	G	334	VAL	4.2
2	E	207	GLY	4.0
2	G	221	ALA	4.0
1	F	104	GLY	4.0
1	F	143	TRP	4.0
1	D	231	VAL	4.0
1	F	89	PRO	3.9
2	A	206	SER	3.9
2	G	260	GLY	3.9
1	H	241	HIS	3.8
1	H	139	SER	3.8
1	H	154	HIS	3.8
1	F	162	MET	3.8
2	A	223	SER	3.8
1	F	188	TYR	3.8
2	G	206	SER	3.8
1	H	220	CYS	3.7
1	F	149	SER	3.7
1	F	153	GLY	3.7
1	H	199	LEU	3.7
1	H	109	LEU	3.7
1	F	148	PHE	3.7
2	G	217	GLU	3.7
1	F	134	ASN	3.6
2	E	222	TYR	3.6
1	F	154	HIS	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	239	PHE	3.6
1	H	192	ILE	3.6
2	E	330	PRO	3.6
2	E	247	ASP	3.6
1	F	129	PRO	3.6
1	H	91	LEU	3.6
2	G	286	THR	3.5
1	D	97	SER	3.5
1	F	94	LEU	3.5
2	G	207	GLY	3.5
2	G	333	PRO	3.5
1	F	124	VAL	3.5
2	G	197	PRO	3.4
1	H	189	GLY	3.4
1	H	136	LEU	3.4
1	F	192	ILE	3.4
1	D	232	ASN	3.4
1	H	112	VAL	3.4
2	E	220	VAL	3.4
1	D	235	SER	3.3
1	H	215	LEU	3.3
1	B	97	SER	3.3
2	G	296	PRO	3.3
1	H	216	THR	3.3
1	F	136	LEU	3.3
2	G	219	TYR	3.3
2	A	205	ALA	3.3
2	A	224	LYS	3.3
1	H	177	PHE	3.3
1	F	158	MET	3.3
2	E	296	PRO	3.2
2	G	297	LEU	3.2
1	F	151	ALA	3.2
1	F	98	PHE	3.2
2	G	220	VAL	3.2
2	G	222	TYR	3.2
2	E	206	SER	3.2
1	H	193	ARG	3.2
1	H	224	SER	3.2
1	F	140	ASP	3.1
1	F	135	CYS	3.1
1	F	177	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	F	197	ARG	3.1
1	F	189	GLY	3.1
2	A	260	GLY	3.1
1	F	145	ILE	3.1
1	F	229	CYS	3.1
1	D	204	HIS	3.0
2	G	330	PRO	3.0
2	G	202	ILE	3.0
2	C	221	ALA	3.0
1	F	99	TYR	3.0
1	H	140	ASP	3.0
1	F	138	LYS	3.0
1	F	201	ARG	3.0
1	H	202	GLY	3.0
2	A	196	ALA	3.0
2	G	276	ALA	3.0
1	F	163	LEU	3.0
1	F	198	HIS	3.0
1	H	217	ILE	3.0
2	E	287	LYS	3.0
1	H	100	TYR	3.0
2	E	294	ASP	3.0
2	E	218	GLY	3.0
2	E	297	LEU	2.9
2	E	295	GLU	2.9
1	B	235	SER	2.9
1	D	197	ARG	2.9
1	F	131	ASP	2.9
2	G	230	GLY	2.9
1	F	212	VAL	2.9
1	D	198	HIS	2.9
1	F	127	TYR	2.9
1	F	202	GLY	2.9
2	G	278	GLY	2.9
1	F	182	ALA	2.9
1	F	91	LEU	2.9
1	F	132	LEU	2.9
1	F	152	LEU	2.9
1	H	174	ASN	2.9
1	F	187	VAL	2.9
2	G	213	VAL	2.9
2	C	291	SER	2.9

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Mol	Chain	Res	Type	RSRZ
2	G	223	SER	2.9
1	H	110	ILE	2.9
1	F	100	TYR	2.9
1	H	214	TYR	2.9
2	E	226	THR	2.9
1	H	163	LEU	2.9
1	F	150	ASN	2.8
1	H	101	MET	2.8
2	G	272	GLN	2.8
1	F	173	SER	2.8
1	H	179	ILE	2.8
2	A	331	SER	2.8
1	F	161	PRO	2.8
1	H	173	SER	2.8
1	H	240	ASP	2.8
2	C	220	VAL	2.8
2	G	209	LEU	2.8
1	F	107	THR	2.8
2	A	217	GLU	2.8
2	A	221	ALA	2.8
2	E	248	TYR	2.8
1	H	229	CYS	2.7
2	E	377	VAL	2.7
2	G	263	THR	2.7
2	E	223	SER	2.7
1	F	226	PHE	2.7
1	H	107	THR	2.7
2	C	286	THR	2.7
2	E	205	ALA	2.7
2	E	249	ALA	2.7
1	H	108	PHE	2.7
1	D	99	TYR	2.7
2	C	287	LYS	2.7
2	G	211	TYR	2.7
1	F	141	SER	2.7
1	H	218	ASN	2.7
1	H	239	PHE	2.7
2	A	294	ASP	2.7
1	F	179	ILE	2.7
2	C	205	ALA	2.7
1	H	89	PRO	2.7
1	F	102	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	H	104	GLY	2.6
1	F	95	ASN	2.6
2	E	221	ALA	2.6
2	G	265	ALA	2.6
1	B	236	LYS	2.6
1	H	95	ASN	2.6
2	G	293	ALA	2.6
1	H	145	ILE	2.6
1	H	228	TYR	2.6
1	D	138	LYS	2.6
2	C	295	GLU	2.6
1	H	117	VAL	2.6
1	F	199	LEU	2.6
2	C	211	TYR	2.6
1	D	233	HIS	2.6
1	F	130	GLU	2.6
1	D	139	SER	2.5
1	F	90	SER	2.5
1	F	235	SER	2.5
1	H	152	LEU	2.5
1	F	193	ARG	2.5
1	H	138	LYS	2.5
2	A	292	ARG	2.5
2	G	244	GLU	2.5
2	G	218	GLY	2.5
1	H	231	VAL	2.5
1	F	203	PHE	2.5
1	B	96	ASN	2.5
1	H	219	GLN	2.5
1	H	116	SER	2.5
1	F	215	LEU	2.5
1	F	105	VAL	2.5
1	H	226	PHE	2.5
1	F	139	SER	2.5
1	H	149	SER	2.5
1	H	143	TRP	2.5
1	H	175	ILE	2.5
1	H	233	HIS	2.5
1	F	231	VAL	2.5
2	G	373	VAL	2.5
1	D	95	ASN	2.5
1	F	169	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	F	207	CYS	2.5
1	F	93	MET	2.5
1	F	234	LEU	2.4
1	F	184	ASP	2.4
2	A	287	LYS	2.4
2	G	377	VAL	2.4
1	H	234	LEU	2.4
1	F	146	HIS	2.4
1	D	140	ASP	2.4
1	F	159	ASP	2.4
1	F	183	ASP	2.4
1	D	172	GLY	2.4
1	D	136	LEU	2.4
1	F	142	SER	2.4
2	E	291	SER	2.4
1	F	92	CYS	2.4
2	E	272	GLN	2.4
2	G	250	VAL	2.4
2	E	285	ARG	2.4
2	G	199	SER	2.4
1	D	238	GLN	2.4
2	C	260	GLY	2.3
1	F	128	GLU	2.3
1	H	200	PHE	2.3
1	H	144	ALA	2.3
1	F	101	MET	2.3
1	F	196	MET	2.3
1	H	230	GLY	2.3
1	H	237	CYS	2.3
2	G	268	VAL	2.3
2	C	222	TYR	2.3
1	H	184	ASP	2.3
2	E	252	GLY	2.3
1	F	191	LYS	2.2
1	F	214	TYR	2.2
1	H	223	PRO	2.2
1	F	172	GLY	2.2
2	G	233	VAL	2.2
2	A	378	ASN	2.2
1	H	221	GLY	2.2
2	C	331	SER	2.2
2	E	260	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	201	ARG	2.2
2	E	292	ARG	2.2
1	H	194	ASN	2.2
1	F	176	GLN	2.2
1	F	236	LYS	2.2
2	G	287	LYS	2.2
2	E	293	ALA	2.2
2	G	225	ALA	2.2
2	C	217	GLU	2.2
2	A	285	ARG	2.2
2	A	296	PRO	2.2
2	C	224	LYS	2.2
1	H	172	GLY	2.2
1	H	183	ASP	2.2
1	F	113	SER	2.2
1	F	204	HIS	2.2
2	G	228	VAL	2.2
1	F	96	ASN	2.1
2	G	204	ASN	2.1
2	G	212	LEU	2.1
1	H	198	HIS	2.1
2	G	331	SER	2.1
2	G	259	ALA	2.1
1	D	94	LEU	2.1
1	H	99	TYR	2.1
1	H	157	LEU	2.1
1	D	183	ASP	2.1
1	H	115	ILE	2.1
2	E	250	VAL	2.1
1	H	196	MET	2.1
2	G	335	GLN	2.1
1	F	155	ASP	2.1
1	D	141	SER	2.1
2	E	225	ALA	2.1
1	H	176	GLN	2.1
1	F	175	ILE	2.1
2	G	277	ILE	2.1
1	H	222	ASP	2.0
1	H	185	ALA	2.0
1	F	223	PRO	2.0
1	F	233	HIS	2.0
1	H	137	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
2	E	332	ILE	2.0
2	G	208	GLY	2.0
1	D	199	LEU	2.0
2	E	378	ASN	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](#)

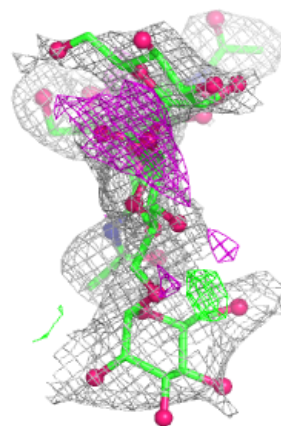
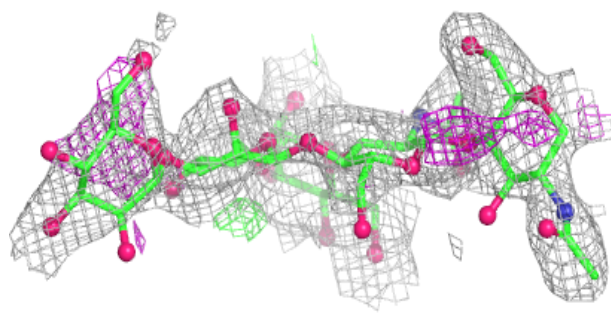
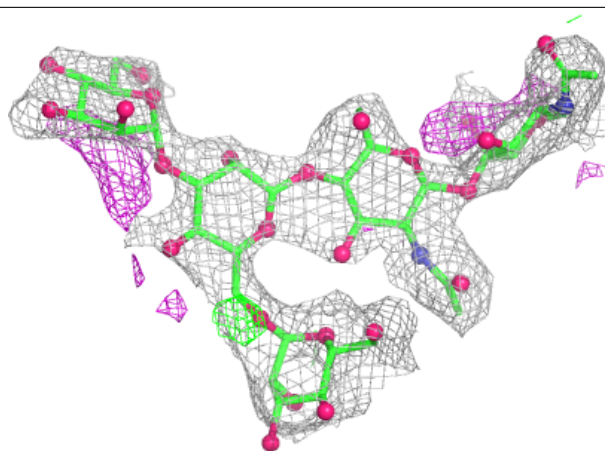
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MAN	I	5	11/12	0.33	0.22	92,103,106,107	0
3	BMA	I	3	11/12	0.64	0.17	86,93,97,98	0
4	NAG	J	2	14/15	0.70	0.13	73,79,87,88	0
3	MAN	I	4	11/12	0.71	0.15	93,98,102,106	0
3	NAG	I	2	14/15	0.72	0.16	66,86,92,95	0
3	NAG	I	1	14/15	0.79	0.15	42,60,74,80	0
4	NAG	J	1	14/15	0.92	0.08	23,47,55,62	0

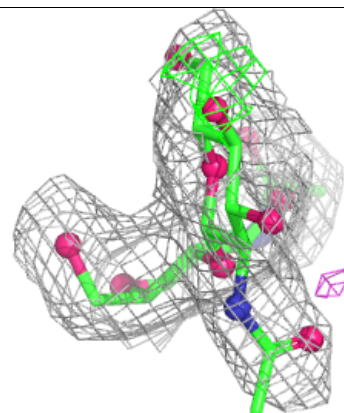
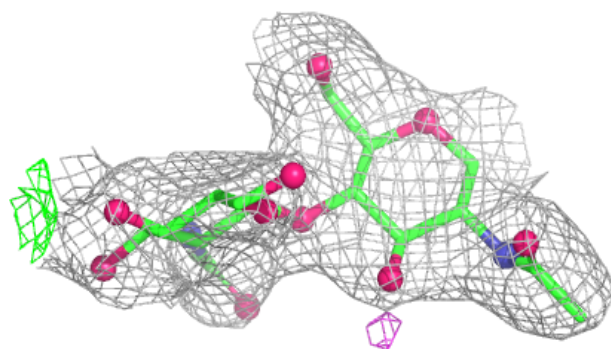
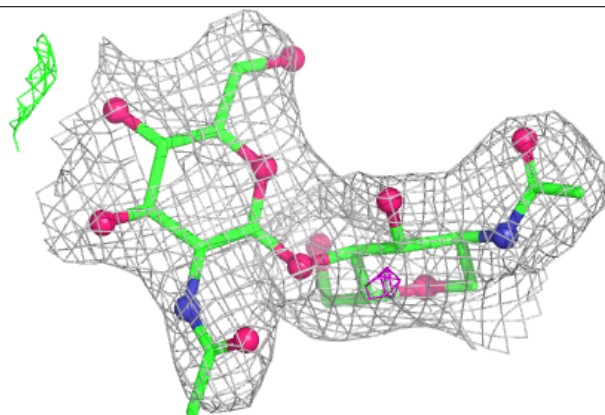
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	B	306	14/15	0.73	0.14	58,68,81,89	0
5	NAG	A	401	14/15	0.76	0.12	56,63,80,86	0
5	NAG	B	307	14/15	0.78	0.14	63,68,84,85	0

6.5 Other polymers [i](#)

There are no such residues in this entry.