



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:08 PM UTC

PDB ID : 1FE8 / pdb_00001fe8
Title : CRYSTAL STRUCTURE OF THE VON WILLEBRAND FACTOR A3 DOMAIN IN COMPLEX WITH A FAB FRAGMENT OF IGG RU5 THAT INHIBITS COLLAGEN BINDING
Authors : Bouma, B.; Huizinga, E.G.; Schiphorst, M.E.; Sixma, J.J.; Kroon, J.; Gros, P.
Deposited on : 2000-07-21
Resolution : 2.03 Å(reported)

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

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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

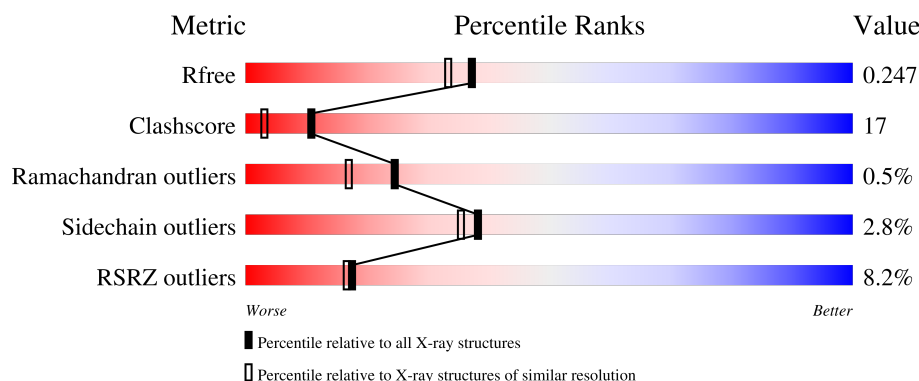
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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	196	7% 70% 23% ...
1	B	196	5% 74% 22% ..
1	C	196	13% 66% 31% ..
2	H	210	9% 65% 31% .
2	I	210	5% 66% 29% 5%

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Mol	Chain	Length	Quality of chain
2	J	210	
3	L	211	
3	M	211	
3	N	211	
4	D	2	
4	E	2	
4	F	2	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NAG	D	1	-	-	X	-
4	FUC	D	2	-	-	X	-
4	FUC	F	2	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 14872 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 VON WILLEBRAND FACTOR.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	190	Total	C	N	O	S	Se	0	0	0
			1421	899	243	273	2	4			
1	B	190	Total	C	N	O	S	Se	0	0	0
			1421	899	243	273	2	4			
1	C	191	Total	C	N	O	S	Se	0	0	0
			1426	902	244	274	2	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	916	GLY	-	cloning artifact	? P04275
A	917	SER	-	cloning artifact	? P04275
A	918	HIS	-	cloning artifact	? P04275
A	919	MSE	-	cloning artifact	? P04275
A	947	MSE	MET	modified residue	? P04275
A	998	MSE	MET	modified residue	? P04275
A	1022	MSE	MET	modified residue	? P04275
A	1097	MSE	MET	modified residue	? P04275
B	916	GLY	-	cloning artifact	? P04275
B	917	SER	-	cloning artifact	? P04275
B	918	HIS	-	cloning artifact	? P04275
B	919	MSE	-	cloning artifact	? P04275
B	947	MSE	MET	modified residue	? P04275
B	998	MSE	MET	modified residue	? P04275
B	1022	MSE	MET	modified residue	? P04275
B	1097	MSE	MET	modified residue	? P04275
C	916	GLY	-	cloning artifact	? P04275
C	917	SER	-	cloning artifact	? P04275
C	918	HIS	-	cloning artifact	? P04275
C	919	MSE	-	cloning artifact	? P04275
C	947	MSE	MET	modified residue	? P04275
C	998	MSE	MET	modified residue	? P04275
C	1022	MSE	MET	modified residue	? P04275

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1097	MSE	MET	modified residue	? P04275

- Molecule 2 is a protein called IMMUNOGLOBULIN IGG RU5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	210	Total	C	N	O	S	0	0	0
			1578	999	258	316	5			
2	I	210	Total	C	N	O	S	0	0	0
			1578	999	258	316	5			
2	J	210	Total	C	N	O	S	0	0	0
			1578	999	258	316	5			

- Molecule 3 is a protein called IMMUNOGLOBULIN IGG RU5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	0
			1629	1009	277	337	6			
3	M	211	Total	C	N	O	S	0	0	0
			1629	1009	277	337	6			
3	N	211	Total	C	N	O	S	0	0	0
			1629	1009	277	337	6			

- Molecule 4 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.



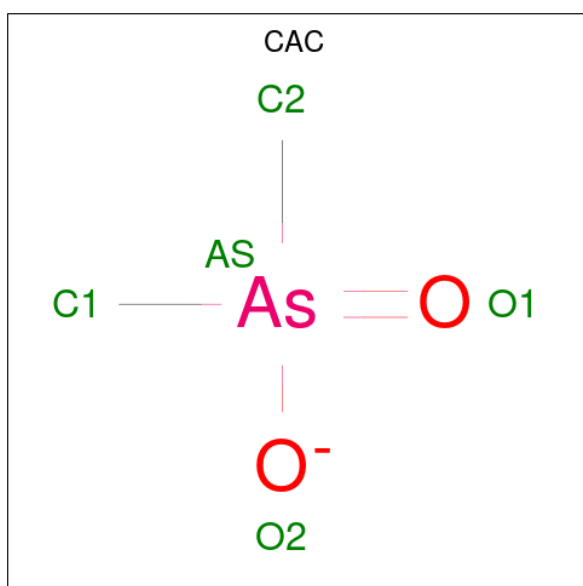
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	2	Total	C	N	O	0	0	0
			24	14	1	9			
4	E	2	Total	C	N	O	0	0	0
			24	14	1	9			
4	F	2	Total	C	N	O	0	0	0
			24	14	1	9			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	H	1	Total	C	N	O	0	0
			14	8	1	5		
5	I	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is CACODYLATE ION (CCD ID: CAC) (formula: $C_2H_6AsO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	L	1	Total	As	C	O	0	0
			5	1	2	2		

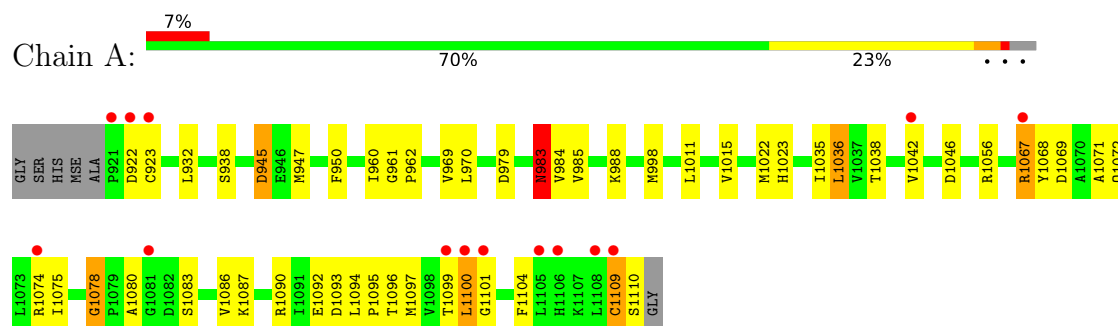
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	84	Total 84	O 84	0	0
7	H	87	Total 87	O 87	0	0
7	L	163	Total 163	O 163	0	0
7	B	79	Total 79	O 79	0	0
7	I	107	Total 107	O 107	0	0
7	M	139	Total 139	O 139	0	0
7	C	58	Total 58	O 58	0	0
7	J	48	Total 48	O 48	0	0
7	N	113	Total 113	O 113	0	0

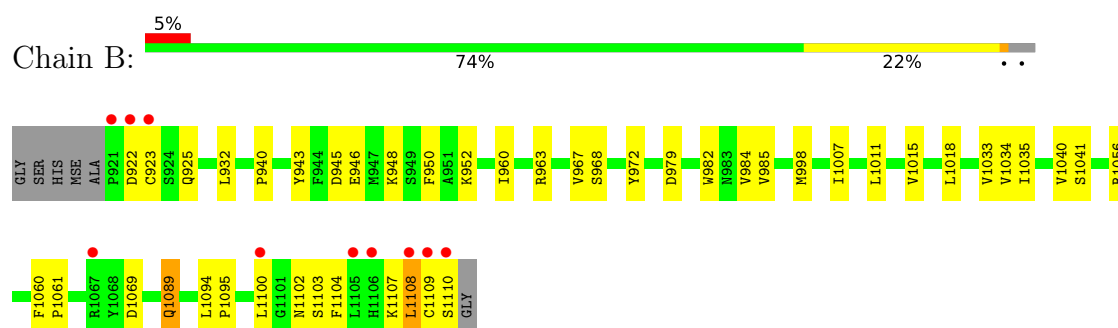
3 Residue-property plots [i](#)

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

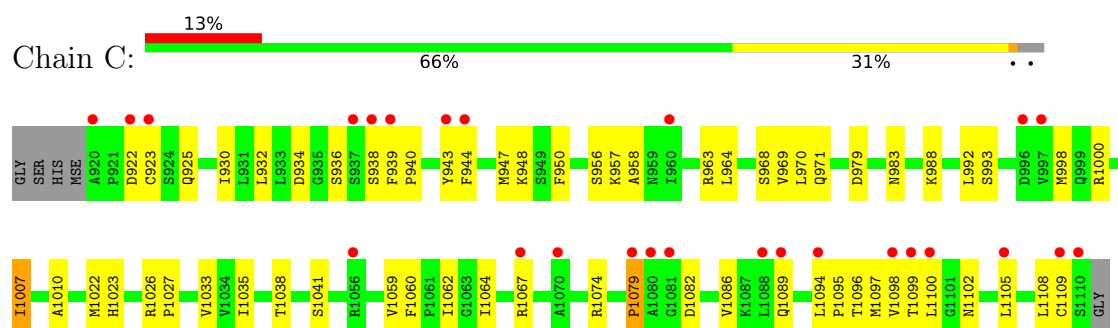
• Molecule 1: VON WILLEBRAND FACTOR



• Molecule 1: VON WILLEBRAND FACTOR

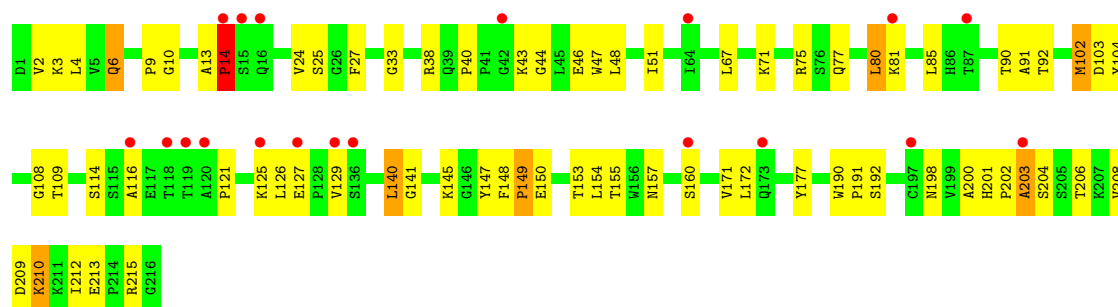


• Molecule 1: VON WILLEBRAND FACTOR

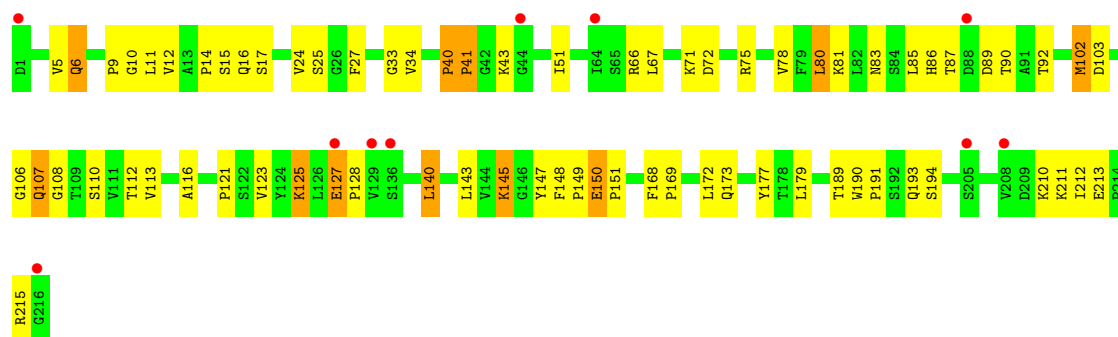


• Molecule 2: IMMUNOGLOBULIN IGG RU5

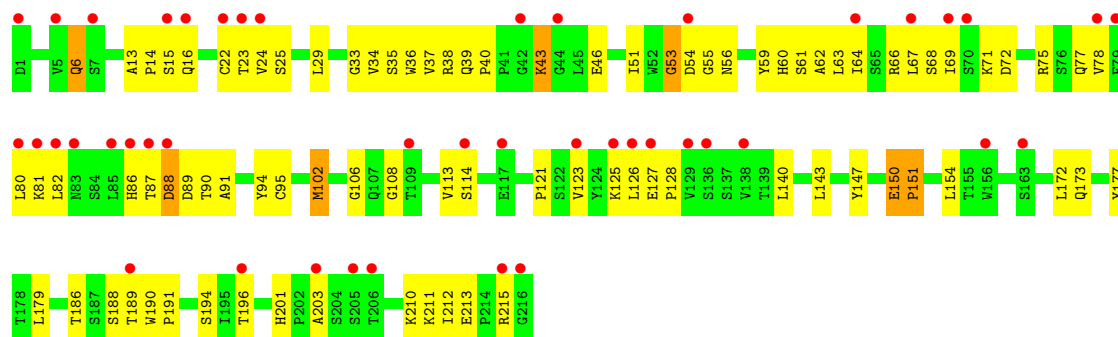




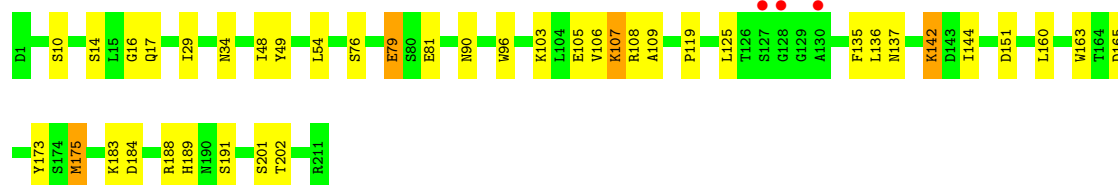
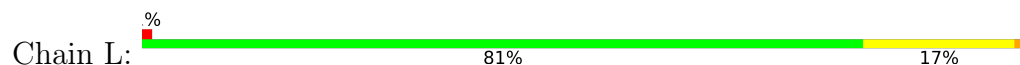
• Molecule 2: IMMUNOGLOBULIN IGG RU5



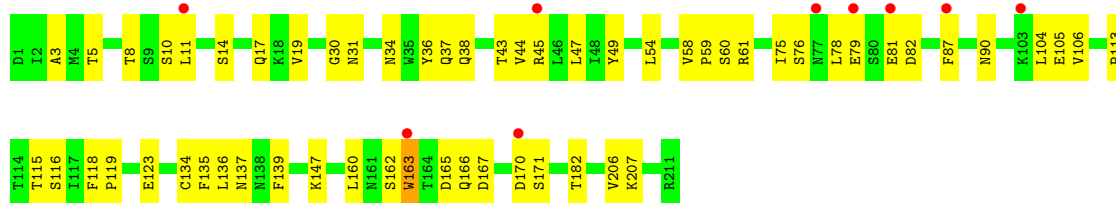
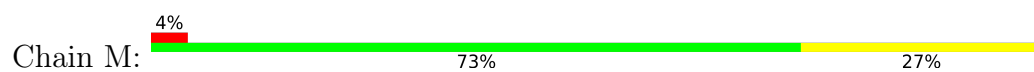
• Molecule 2: IMMUNOGLOBULIN IGG RU5



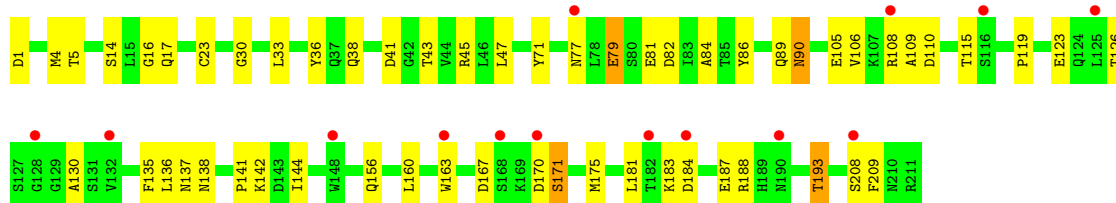
• Molecule 3: IMMUNOGLOBULIN IGG RU5



• Molecule 3: IMMUNOGLOBULIN IGG RU5



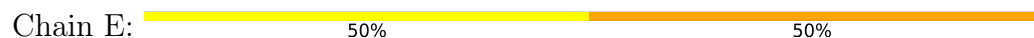
- Molecule 3: IMMUNOGLOBULIN IGG RU5



- Molecule 4: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.80Å 183.60Å 131.80Å 90.00° 116.20° 90.00°	Depositor
Resolution (Å)	29.90 – 2.03 29.90 – 2.03	Depositor EDS
% Data completeness (in resolution range)	85.9 (29.90-2.03) 85.8 (29.90-2.03)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.03Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.227 , 0.264 (Not available) , 0.247	Depositor DCC
R_{free} test set	7206 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.011 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14872	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% 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: FUC, CAC, 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	0.42	0/1443	0.95	5/1956 (0.3%)
1	B	0.41	0/1443	0.95	1/1956 (0.1%)
1	C	0.37	0/1448	0.93	2/1964 (0.1%)
2	H	0.40	0/1618	0.90	6/2215 (0.3%)
2	I	0.38	0/1618	0.94	10/2215 (0.5%)
2	J	0.34	0/1618	0.89	7/2215 (0.3%)
3	L	0.43	0/1664	0.93	6/2260 (0.3%)
3	M	0.40	0/1664	0.87	4/2260 (0.2%)
3	N	0.36	0/1664	0.85	5/2260 (0.2%)
All	All	0.39	0/14180	0.91	46/19301 (0.2%)

There are no bond length outliers.

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	150	GLU	N-CA-C	-8.28	98.01	110.01
2	I	102	MET	N-CA-C	-7.69	98.86	109.71
3	L	142	LYS	N-CA-C	7.41	121.42	112.38
1	A	938	SER	N-CA-C	7.36	119.30	111.28
2	H	103	ASP	N-CA-C	7.28	119.21	111.28
2	H	33	GLY	N-CA-C	-7.26	101.42	112.89
2	I	103	ASP	N-CA-C	7.25	118.83	111.07
3	L	165	ASP	N-CA-C	-7.05	101.43	110.53
2	J	88	ASP	N-CA-C	-6.61	105.32	113.19
1	A	1078	GLY	CA-C-N	6.53	125.77	118.97
1	A	1078	GLY	C-N-CA	6.53	125.77	118.97
2	J	55	GLY	N-CA-C	-6.36	106.94	115.21
2	I	33	GLY	N-CA-C	-6.32	102.91	112.89
2	I	106	GLY	N-CA-C	-6.15	103.18	112.89
3	L	137	ASN	N-CA-C	6.11	119.66	109.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	M	137	ASN	N-CA-C	6.05	119.57	109.95
2	I	87	THR	N-CA-C	-6.02	105.04	112.38
2	J	106	GLY	N-CA-C	-6.01	103.39	112.89
1	C	964	LEU	N-CA-C	5.98	114.49	108.75
3	M	5	THR	N-CA-C	5.97	118.15	108.41
2	J	102	MET	N-CA-C	-5.96	101.30	109.71
2	I	40	PRO	CA-C-N	5.89	127.21	119.84
2	I	40	PRO	C-N-CA	5.89	127.21	119.84
3	L	202	THR	N-CA-C	-5.76	106.08	113.23
1	C	1007	ILE	N-CA-C	5.74	116.47	110.62
3	L	76	SER	N-CA-C	5.67	117.46	111.28
2	I	145	LYS	N-CA-C	5.66	118.78	109.72
2	J	53	GLY	N-CA-C	-5.58	108.54	114.67
1	A	1109	CYS	N-CA-C	5.57	115.90	108.38
3	N	137	ASN	N-CA-C	5.53	119.08	110.17
2	I	150	GLU	N-CA-C	-5.53	102.24	110.20
3	M	206	VAL	N-CA-C	5.41	116.38	108.53
3	L	201	SER	N-CA-C	5.38	118.32	109.76
2	H	92	THR	N-CA-C	-5.33	100.62	109.46
3	N	138	ASN	N-CA-C	5.33	118.59	111.24
3	N	171	SER	N-CA-C	-5.30	106.76	113.23
1	A	983	ASN	N-CA-C	5.21	121.65	113.72
3	N	193	THR	N-CA-C	5.16	117.84	109.06
3	N	142	LYS	N-CA-C	5.09	117.56	111.71
2	H	150	GLU	N-CA-C	-5.08	102.89	110.20
1	B	1034	VAL	N-CA-C	-5.05	100.62	107.99
2	H	13	ALA	CA-C-N	5.04	126.14	119.84
2	H	13	ALA	C-N-CA	5.04	126.14	119.84
2	I	168	PHE	N-CA-C	5.03	117.24	110.29
3	M	165	ASP	N-CA-C	-5.03	103.30	110.59
2	J	33	GLY	N-CA-C	-5.01	104.97	112.89

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1421	0	1433	42	0
1	B	1421	0	1433	31	0
1	C	1426	0	1437	50	0
2	H	1578	0	1540	68	0
2	I	1578	0	1540	54	0
2	J	1578	0	1540	93	0
3	L	1629	0	1554	36	0
3	M	1629	0	1554	45	0
3	N	1629	0	1554	41	0
4	D	24	0	22	8	0
4	E	24	0	22	6	0
4	F	24	0	22	8	0
5	H	14	0	13	2	0
5	I	14	0	13	1	0
6	L	5	0	0	0	0
7	A	84	0	0	3	0
7	B	79	0	0	1	0
7	C	58	0	0	2	0
7	H	87	0	0	2	0
7	I	107	0	0	5	0
7	J	48	0	0	2	0
7	L	163	0	0	8	0
7	M	139	0	0	9	0
7	N	113	0	0	2	0
All	All	14872	0	13677	458	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 17.

All (458) 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:C:944:PHE:HA	1:C:947:MSE:HE3	1.31	1.06
2:J:90:THR:HG22	2:J:113:VAL:H	1.22	1.04
2:H:121:PRO:HB3	2:H:147:TYR:HB3	1.48	0.96
2:J:150:GLU:HG3	2:J:151:PRO:HA	1.52	0.92
1:B:1108:LEU:HD23	1:B:1108:LEU:H	1.36	0.91
2:I:40:PRO:HG2	2:I:43:LYS:HB2	1.58	0.85
2:I:121:PRO:HB3	2:I:147:TYR:HB3	1.59	0.83
1:B:1035:ILE:HG23	1:B:1061:PRO:HA	1.62	0.82
1:A:988:LYS:HB2	4:D:2:FUC:H4	1.60	0.82
2:H:38:ARG:HG2	2:H:48:LEU:HD21	1.61	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:6:GLN:H	2:I:107:GLN:HE22	1.29	0.81
2:I:125:LYS:HZ3	2:I:213:GLU:H	1.29	0.79
2:H:127:GLU:HG2	2:H:213:GLU:O	1.82	0.79
1:C:1062:ILE:HD13	1:C:1097:MSE:HE1	1.64	0.78
2:J:90:THR:HG22	2:J:113:VAL:N	1.99	0.78
2:J:43:LYS:HE3	2:J:43:LYS:HA	1.65	0.77
2:J:121:PRO:HB3	2:J:147:TYR:HB3	1.67	0.76
4:E:1:NAG:H61	4:E:2:FUC:H5	1.68	0.76
1:A:1097:MSE:HE2	1:A:1104:PHE:CD1	2.21	0.76
1:C:993:SER:HB2	2:J:54:ASP:HB2	1.68	0.75
3:L:142:LYS:HB2	7:L:1058:HOH:O	1.85	0.75
3:M:38:GLN:HG3	7:M:312:HOH:O	1.87	0.75
4:D:1:NAG:H62	4:D:2:FUC:H3	1.68	0.74
2:I:127:GLU:HG2	2:I:213:GLU:O	1.87	0.73
4:F:1:NAG:H61	4:F:2:FUC:H5	1.69	0.73
2:J:127:GLU:HG2	2:J:213:GLU:O	1.88	0.73
2:J:6:GLN:H	2:J:6:GLN:HE21	1.37	0.73
5:H:553:NAG:C1	4:D:1:NAG:H4	2.19	0.73
2:J:40:PRO:HB2	2:J:43:LYS:HB2	1.69	0.72
1:C:944:PHE:CA	1:C:947:MSE:HE3	2.16	0.72
1:C:944:PHE:HA	1:C:947:MSE:CE	2.16	0.72
2:J:90:THR:CG2	2:J:113:VAL:H	2.01	0.71
3:N:183:LYS:O	3:N:187:GLU:HG2	1.91	0.70
2:J:6:GLN:NE2	2:J:108:GLY:H	1.89	0.70
3:M:79:GLU:HG2	3:M:81:GLU:H	1.57	0.69
2:J:86:HIS:C	2:J:88:ASP:H	1.99	0.69
3:M:167:ASP:OD2	3:M:170:ASP:HB2	1.93	0.69
2:H:51:ILE:HD13	2:H:71:LYS:HG2	1.73	0.69
4:E:1:NAG:H61	4:E:2:FUC:H3	1.73	0.69
2:I:125:LYS:NZ	2:I:213:GLU:H	1.91	0.68
1:A:1094:LEU:HB3	1:A:1095:PRO:HD3	1.76	0.68
2:H:125:LYS:HE3	2:H:210:LYS:HD2	1.75	0.67
3:M:54:LEU:HD11	3:M:60:SER:HA	1.75	0.67
3:N:108:ARG:HG2	3:N:108:ARG:HH11	1.59	0.67
2:I:6:GLN:NE2	2:I:107:GLN:NE2	2.42	0.67
3:N:184:ASP:HB3	3:N:188:ARG:NH1	2.09	0.67
7:I:584:HOH:O	4:E:1:NAG:H4	1.94	0.66
2:H:6:GLN:H	2:H:6:GLN:HE21	1.43	0.65
2:I:72:ASP:OD1	2:I:75:ARG:HD3	1.95	0.65
3:L:108:ARG:HG2	3:L:109:ALA:N	2.11	0.65
2:J:87:THR:O	2:J:90:THR:HG23	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:553:NAG:H82	4:D:1:NAG:O4	1.97	0.65
2:H:80:LEU:C	2:H:80:LEU:HD13	2.22	0.65
3:L:14:SER:HB2	3:L:17:GLN:HE21	1.62	0.64
2:H:47:TRP:HD1	2:H:102:MET:HE1	1.63	0.64
2:I:145:LYS:HE2	7:I:582:HOH:O	1.97	0.64
1:A:1056:ARG:HG2	1:A:1056:ARG:HH11	1.63	0.63
3:M:78:LEU:HD11	3:M:104:LEU:HD21	1.80	0.63
3:L:105:GLU:HG3	3:L:173:TYR:OH	1.98	0.63
2:H:102:MET:HE3	3:L:96:TRP:CZ3	2.33	0.63
3:N:105:GLU:HG2	3:N:106:VAL:N	2.14	0.63
1:B:932:LEU:HD23	1:B:1035:ILE:HD13	1.81	0.63
1:B:1040:VAL:HG12	1:B:1069:ASP:HB2	1.79	0.63
2:I:6:GLN:NE2	2:I:108:GLY:H	1.97	0.63
2:H:6:GLN:HB2	2:H:109:THR:HG23	1.80	0.63
2:J:150:GLU:HG3	2:J:151:PRO:CA	2.29	0.63
1:B:923:CYS:HB2	1:B:1110:SER:O	1.99	0.62
2:J:23:THR:HG22	2:J:77:GLN:HG2	1.80	0.62
2:I:194:SER:HB2	2:I:211:LYS:HE2	1.81	0.62
3:L:183:LYS:HE3	7:L:1112:HOH:O	1.98	0.62
2:J:87:THR:H	2:J:113:VAL:HG11	1.65	0.62
1:B:945:ASP:HB2	7:B:488:HOH:O	2.00	0.62
3:M:10:SER:HB2	7:M:276:HOH:O	2.00	0.62
2:H:102:MET:HE3	3:L:96:TRP:HZ3	1.63	0.61
2:H:172:LEU:HB2	2:H:177:TYR:CE1	2.34	0.61
3:L:184:ASP:HB3	3:L:188:ARG:HH12	1.65	0.61
1:C:934:ASP:O	1:C:947:MSE:HE1	2.00	0.61
2:H:47:TRP:CD1	2:H:102:MET:HE1	2.36	0.61
4:D:1:NAG:H62	4:D:2:FUC:H5	1.82	0.61
2:J:38:ARG:HD2	2:J:46:GLU:OE1	2.01	0.61
1:A:979:ASP:OD2	1:A:998:MSE:HE3	1.99	0.61
1:C:1062:ILE:CD1	1:C:1097:MSE:HE1	2.30	0.60
2:J:140:LEU:HD12	2:J:212:ILE:HG21	1.83	0.60
2:J:39:GLN:C	2:J:91:ALA:HB1	2.26	0.60
2:J:86:HIS:O	2:J:88:ASP:N	2.34	0.60
3:L:79:GLU:OE2	1:C:1023:HIS:HE1	1.83	0.60
2:H:6:GLN:HE22	2:H:108:GLY:H	1.49	0.60
2:I:5:VAL:HA	2:I:107:GLN:HE21	1.67	0.59
1:C:1007:ILE:HG23	1:C:1035:ILE:HD11	1.83	0.59
2:J:67:LEU:HD11	2:J:80:LEU:HD21	1.84	0.59
2:J:87:THR:HA	2:J:113:VAL:HB	1.84	0.59
2:H:44:GLY:HA3	7:L:1065:HOH:O	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:67:LEU:HD11	2:J:80:LEU:CD2	2.32	0.59
2:I:6:GLN:H	2:I:6:GLN:HE21	1.49	0.59
3:M:11:LEU:HD21	7:M:344:HOH:O	2.02	0.59
2:J:6:GLN:HE22	2:J:108:GLY:H	1.48	0.59
1:B:1104:PHE:O	1:B:1107:LYS:HB2	2.03	0.59
1:B:979:ASP:OD2	1:B:998:MSE:HE3	2.03	0.58
2:I:6:GLN:NE2	2:I:107:GLN:HE22	2.00	0.58
1:C:988:LYS:O	1:C:992:LEU:HD13	2.03	0.58
3:N:16:GLY:O	3:N:77:ASN:HA	2.03	0.58
2:H:210:LYS:HB2	2:H:210:LYS:NZ	2.19	0.58
3:M:54:LEU:HD22	3:M:58:VAL:HB	1.85	0.58
2:H:125:LYS:HD3	2:H:127:GLU:OE1	2.04	0.58
2:H:154:LEU:HD23	2:H:155:THR:N	2.19	0.58
2:I:173:GLN:HG2	7:I:638:HOH:O	2.04	0.58
1:C:956:SER:HA	1:C:988:LYS:HD3	1.85	0.58
1:C:1074:ARG:HG3	1:C:1082:ASP:OD2	2.04	0.58
2:H:116:ALA:HB3	2:H:148:PHE:CE2	2.39	0.58
2:J:80:LEU:C	2:J:80:LEU:HD13	2.29	0.57
2:J:186:THR:HB	2:J:189:THR:HG23	1.85	0.57
2:H:208:VAL:HG12	2:H:209:ASP:N	2.19	0.57
2:H:6:GLN:NE2	2:H:108:GLY:H	2.01	0.57
4:E:1:NAG:H61	4:E:2:FUC:C5	2.34	0.57
1:A:1086:VAL:HG21	1:A:1097:MSE:CE	2.35	0.57
2:H:43:LYS:HB3	7:H:596:HOH:O	2.05	0.57
3:L:105:GLU:HG2	3:L:106:VAL:N	2.20	0.57
1:A:1022:MSE:HG2	7:N:286:HOH:O	2.05	0.57
1:C:943:TYR:O	1:C:947:MSE:HG3	2.05	0.57
2:J:14:PRO:C	2:J:16:GLN:H	2.12	0.57
2:I:6:GLN:HE21	2:I:107:GLN:HE22	1.51	0.57
2:J:38:ARG:HD3	2:J:40:PRO:HG3	1.87	0.56
3:L:16:GLY:N	1:C:1022:MSE:HE1	2.21	0.56
2:I:67:LEU:HD11	2:I:80:LEU:CD2	2.36	0.56
3:M:43:THR:HA	7:M:277:HOH:O	2.05	0.56
2:J:186:THR:HG22	2:J:188:SER:H	1.70	0.56
4:D:1:NAG:H62	4:D:2:FUC:C3	2.35	0.56
3:L:151:ASP:HA	3:L:191:SER:OG	2.06	0.56
4:E:1:NAG:H61	4:E:2:FUC:C3	2.36	0.56
2:I:5:VAL:HA	2:I:107:GLN:NE2	2.20	0.56
2:I:140:LEU:HG	2:I:212:ILE:HG21	1.88	0.56
1:C:939:PHE:HB3	1:C:940:PRO:CD	2.35	0.55
2:J:16:GLN:HG2	7:J:567:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:157:ASN:HB3	2:H:160:SER:HB3	1.88	0.55
2:J:201:HIS:CE1	2:J:203:ALA:HB3	2.41	0.55
2:I:116:ALA:HB3	2:I:148:PHE:CE2	2.42	0.55
3:M:31:ASN:HD21	3:N:156:GLN:NE2	2.06	0.54
1:C:1095:PRO:O	1:C:1099:THR:HG23	2.08	0.54
3:L:81:GLU:HG2	7:L:1051:HOH:O	2.06	0.54
2:I:51:ILE:HD13	2:I:71:LYS:HG2	1.90	0.54
3:N:126:THR:HG22	3:N:126:THR:O	2.07	0.53
1:B:1011:LEU:O	1:B:1015:VAL:HG23	2.08	0.53
3:M:19:VAL:HG22	3:M:75:ILE:HB	1.89	0.53
3:M:59:PRO:C	3:M:61:ARG:H	2.15	0.53
3:N:144:ILE:HG22	3:N:163:TRP:CZ3	2.43	0.53
2:I:6:GLN:H	2:I:6:GLN:NE2	2.06	0.53
1:C:1038:THR:HG22	1:C:1064:ILE:HB	1.91	0.53
2:J:6:GLN:HE21	2:J:6:GLN:N	2.03	0.53
1:A:1090:ARG:HG2	1:A:1093:ASP:OD2	2.08	0.53
3:N:108:ARG:HG2	3:N:109:ALA:N	2.22	0.53
1:B:1108:LEU:H	1:B:1108:LEU:CD2	2.16	0.53
2:I:6:GLN:HE22	2:I:108:GLY:H	1.56	0.53
1:C:1098:VAL:HG13	1:C:1105:LEU:HD21	1.90	0.53
2:J:59:TYR:HE2	2:J:69:ILE:HG13	1.74	0.53
2:I:86:HIS:O	2:I:113:VAL:HG11	2.09	0.53
3:M:105:GLU:HG2	3:M:106:VAL:N	2.23	0.53
3:N:14:SER:O	3:N:17:GLN:HB2	2.09	0.53
2:J:40:PRO:HA	2:J:91:ALA:HB2	1.90	0.52
1:C:922:ASP:OD2	1:C:925:GLN:HG2	2.08	0.52
1:A:1104:PHE:HB3	7:A:750:HOH:O	2.09	0.52
3:L:107:LYS:HB2	3:L:107:LYS:NZ	2.24	0.52
2:I:80:LEU:HD13	2:I:80:LEU:C	2.33	0.52
1:B:972:TYR:CE2	1:B:1035:ILE:HD11	2.44	0.52
2:H:171:VAL:HG11	3:L:160:LEU:HG	1.92	0.52
2:I:125:LYS:HD3	3:M:123:GLU:OE1	2.09	0.52
3:M:136:LEU:HD12	3:M:136:LEU:N	2.24	0.52
2:H:80:LEU:HD22	2:H:81:LYS:N	2.23	0.52
2:H:215:ARG:CZ	3:L:119:PRO:HG2	2.40	0.52
2:I:125:LYS:HZ2	2:I:127:GLU:CG	2.22	0.52
2:H:51:ILE:CD1	2:H:71:LYS:HG2	2.40	0.52
3:M:167:ASP:HB3	3:M:170:ASP:HB3	1.90	0.52
2:J:87:THR:C	2:J:89:ASP:H	2.15	0.52
2:I:90:THR:HG23	2:I:112:THR:HA	1.92	0.51
4:D:1:NAG:C6	4:D:2:FUC:H5	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:14:SER:CB	3:L:17:GLN:HE21	2.23	0.51
1:A:1056:ARG:HG2	1:A:1056:ARG:NH1	2.24	0.51
1:A:1042:VAL:HA	2:J:189:THR:HG21	1.93	0.51
1:A:1078:GLY:HA2	7:A:459:HOH:O	2.10	0.51
3:L:48:ILE:HD13	3:L:54:LEU:HD23	1.92	0.51
2:J:38:ARG:HG2	2:J:40:PRO:HD3	1.93	0.51
3:L:188:ARG:HD3	7:L:1142:HOH:O	2.10	0.51
1:B:1007:ILE:HD12	1:B:1041:SER:HB2	1.93	0.51
2:J:51:ILE:HG23	2:J:51:ILE:O	2.10	0.51
3:N:181:LEU:HD23	3:N:181:LEU:N	2.26	0.51
7:M:305:HOH:O	2:J:173:GLN:HG3	2.09	0.51
2:J:6:GLN:H	2:J:6:GLN:NE2	2.07	0.51
2:I:102:MET:HB2	3:M:36:TYR:CE2	2.45	0.51
3:M:166:GLN:HG2	3:M:171:SER:HA	1.93	0.50
7:J:580:HOH:O	4:F:2:FUC:H4	2.11	0.50
2:I:127:GLU:HB2	2:I:215:ARG:HG3	1.92	0.50
2:H:24:VAL:HG11	2:H:27:PHE:CE1	2.47	0.50
3:M:14:SER:HB2	3:M:17:GLN:NE2	2.26	0.50
2:J:154:LEU:C	2:J:154:LEU:HD23	2.36	0.50
1:A:1011:LEU:O	1:A:1015:VAL:HG23	2.12	0.50
3:M:163:TRP:N	3:M:163:TRP:CD1	2.79	0.50
1:C:957:LYS:HD3	1:C:1098:VAL:HG12	1.93	0.50
2:J:190:TRP:CG	2:J:191:PRO:HA	2.47	0.50
2:H:38:ARG:HD2	2:H:46:GLU:OE1	2.12	0.50
2:J:90:THR:O	2:J:91:ALA:HB2	2.11	0.50
2:J:90:THR:HG22	2:J:113:VAL:HB	1.94	0.49
2:J:172:LEU:HB2	2:J:177:TYR:CE1	2.47	0.49
2:J:56:ASN:HA	4:F:1:NAG:C8	2.41	0.49
2:J:87:THR:C	2:J:89:ASP:N	2.70	0.49
2:J:128:PRO:O	2:J:215:ARG:HD2	2.12	0.49
1:B:1108:LEU:HD23	1:B:1108:LEU:N	2.18	0.49
2:I:12:VAL:HG11	2:I:85:LEU:CD1	2.43	0.49
2:I:12:VAL:O	2:I:113:VAL:HA	2.11	0.49
3:N:108:ARG:HG2	3:N:108:ARG:NH1	2.26	0.49
3:N:163:TRP:N	3:N:163:TRP:CD1	2.81	0.49
2:H:126:LEU:HB2	2:H:141:GLY:CA	2.43	0.49
3:M:43:THR:C	7:M:312:HOH:O	2.55	0.49
1:B:943:TYR:O	1:B:946:GLU:HB2	2.13	0.49
1:A:923:CYS:N	1:A:1109:CYS:HB2	2.28	0.49
2:J:72:ASP:HB3	2:J:75:ARG:HB2	1.95	0.49
2:I:14:PRO:C	2:I:16:GLN:H	2.21	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:950:PHE:HB2	1:C:1094:LEU:HG	1.94	0.49
1:C:1086:VAL:HG11	1:C:1097:MSE:SE	2.62	0.49
1:A:947:MSE:HG2	1:A:1036:LEU:HD23	1.95	0.48
3:L:125:LEU:O	3:L:183:LYS:HD2	2.13	0.48
2:J:39:GLN:O	2:J:91:ALA:HB1	2.13	0.48
1:A:932:LEU:HD23	1:A:1035:ILE:HD12	1.95	0.48
1:B:922:ASP:HB3	1:B:925:GLN:CG	2.43	0.48
1:B:1094:LEU:HB3	1:B:1095:PRO:CD	2.43	0.48
4:F:1:NAG:H61	4:F:2:FUC:C5	2.40	0.48
2:J:123:VAL:HG23	2:J:123:VAL:O	2.13	0.48
3:L:14:SER:HB2	3:L:17:GLN:NE2	2.28	0.48
1:C:969:VAL:C	1:C:970:LEU:HD12	2.37	0.48
2:H:153:THR:OG1	2:H:200:ALA:HB3	2.13	0.48
2:I:81:LYS:HE3	2:I:83:ASN:OD1	2.13	0.48
3:M:59:PRO:C	3:M:61:ARG:N	2.71	0.48
2:J:37:VAL:HG23	2:J:94:TYR:HB2	1.94	0.48
1:A:984:VAL:O	1:A:985:VAL:C	2.56	0.48
2:I:16:GLN:HG2	2:I:17:SER:H	1.79	0.48
2:I:128:PRO:O	2:I:215:ARG:HD2	2.13	0.48
3:M:31:ASN:HD21	3:N:156:GLN:HE22	1.62	0.48
2:J:43:LYS:HA	2:J:43:LYS:CE	2.38	0.48
2:I:66:ARG:HH22	2:I:89:ASP:CG	2.22	0.47
2:I:107:GLN:HG3	7:I:560:HOH:O	2.14	0.47
2:H:2:VAL:HG11	2:H:104:TYR:CG	2.48	0.47
2:H:40:PRO:HB2	2:H:43:LYS:HE2	1.95	0.47
3:L:34:ASN:OD1	3:L:49:TYR:HA	2.15	0.47
2:J:14:PRO:O	2:J:16:GLN:N	2.47	0.47
3:N:119:PRO:HG3	3:N:209:PHE:CD2	2.49	0.47
4:F:1:NAG:C6	4:F:2:FUC:H5	2.42	0.47
1:C:958:ALA:O	1:C:988:LYS:NZ	2.48	0.47
1:C:1105:LEU:N	1:C:1105:LEU:HD22	2.29	0.47
1:B:968:SER:HB2	1:B:982:TRP:CE3	2.49	0.47
1:A:969:VAL:C	1:A:970:LEU:HD12	2.40	0.47
3:L:135:PHE:C	3:L:136:LEU:HD12	2.40	0.47
2:I:123:VAL:O	2:I:210:LYS:HE3	2.15	0.47
3:M:182:THR:HA	7:M:350:HOH:O	2.15	0.47
2:J:29:LEU:O	2:J:53:GLY:HA2	2.14	0.47
3:N:110:ASP:OD1	3:N:141:PRO:HD3	2.15	0.47
1:A:1038:THR:O	1:A:1067:ARG:HB2	2.15	0.47
1:C:930:ILE:HG12	1:C:968:SER:OG	2.14	0.47
1:C:979:ASP:OD2	1:C:998:MSE:HE3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:194:SER:HB2	2:J:211:LYS:HE3	1.97	0.47
2:J:196:THR:HG23	2:J:210:LYS:O	2.15	0.47
2:J:102:MET:HB2	3:N:36:TYR:CE2	2.50	0.46
1:C:1074:ARG:HH11	1:C:1074:ARG:HG2	1.79	0.46
3:N:170:ASP:O	3:N:171:SER:HB2	2.14	0.46
1:C:1060:PHE:CE1	1:C:1108:LEU:HG	2.49	0.46
2:J:140:LEU:N	2:J:140:LEU:HD22	2.30	0.46
1:A:1086:VAL:HG21	1:A:1097:MSE:HE3	1.96	0.46
2:H:80:LEU:HD22	2:H:81:LYS:H	1.81	0.46
2:H:90:THR:O	2:H:91:ALA:HB2	2.15	0.46
3:M:36:TYR:HB2	3:M:87:PHE:CZ	2.50	0.46
2:J:53:GLY:HA2	2:J:71:LYS:NZ	2.31	0.46
2:J:66:ARG:HH22	2:J:89:ASP:CG	2.22	0.46
2:J:196:THR:HG23	2:J:210:LYS:C	2.40	0.46
2:H:24:VAL:HG12	2:H:25:SER:N	2.31	0.46
2:J:56:ASN:ND2	4:F:2:FUC:H61	2.30	0.46
2:H:191:PRO:O	2:H:192:SER:C	2.58	0.46
2:J:128:PRO:HD3	2:J:140:LEU:HD13	1.98	0.46
2:H:149:PRO:HD2	2:H:203:ALA:CB	2.46	0.46
3:N:47:LEU:HD11	3:N:86:TYR:HE2	1.81	0.46
3:N:79:GLU:HB2	3:N:82:ASP:OD2	2.16	0.46
2:H:14:PRO:HD3	2:H:114:SER:C	2.41	0.46
2:J:6:GLN:NE2	2:J:6:GLN:N	2.63	0.46
2:I:11:LEU:HD22	2:I:149:PRO:HG3	1.98	0.45
2:I:51:ILE:HG23	2:I:51:ILE:O	2.15	0.45
1:C:947:MSE:O	1:C:950:PHE:HB3	2.15	0.45
1:C:1007:ILE:HD12	1:C:1041:SER:HB2	1.99	0.45
2:J:127:GLU:HG3	2:J:212:ILE:CG2	2.46	0.45
2:H:210:LYS:HB2	2:H:210:LYS:HZ3	1.80	0.45
2:J:86:HIS:C	2:J:88:ASP:N	2.65	0.45
1:A:1022:MSE:HE3	7:L:1103:HOH:O	2.17	0.45
2:I:169:PRO:HD2	3:M:162:SER:OG	2.16	0.45
2:J:87:THR:H	2:J:113:VAL:CG1	2.30	0.45
2:J:72:ASP:CG	2:J:75:ARG:HB2	2.42	0.45
3:L:151:ASP:OD2	3:L:189:HIS:HD2	2.00	0.45
1:B:1056:ARG:HH11	1:B:1056:ARG:HG2	1.82	0.45
1:C:1067:ARG:HG2	1:C:1067:ARG:HH11	1.82	0.45
2:H:24:VAL:HG11	2:H:27:PHE:CZ	2.52	0.45
2:I:172:LEU:HB2	2:I:177:TYR:CE1	2.52	0.45
1:C:944:PHE:O	1:C:948:LYS:HG3	2.17	0.45
2:J:22:CYS:HB2	2:J:36:TRP:CH2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:37:VAL:CG2	2:J:94:TYR:HB2	2.47	0.45
2:J:125:LYS:O	2:J:126:LEU:HD23	2.17	0.45
4:F:1:NAG:H61	4:F:2:FUC:H3	1.99	0.45
2:I:190:TRP:CG	2:I:191:PRO:HA	2.52	0.45
2:J:14:PRO:C	2:J:16:GLN:N	2.75	0.45
3:N:47:LEU:HD11	3:N:86:TYR:CE2	2.51	0.45
4:E:1:NAG:C6	4:E:2:FUC:H5	2.42	0.45
1:A:1086:VAL:HG21	1:A:1097:MSE:HE1	2.00	0.45
5:I:553:NAG:H62	7:I:629:HOH:O	2.16	0.45
2:J:179:LEU:HD23	2:J:179:LEU:C	2.41	0.45
2:H:154:LEU:HD23	2:H:154:LEU:C	2.42	0.44
1:C:971:GLN:HE22	1:C:1000:ARG:HA	1.82	0.44
3:N:45:ARG:HG3	3:N:45:ARG:NH1	2.32	0.44
1:A:960:ILE:HG22	1:A:961:GLY:N	2.31	0.44
1:A:1069:ASP:OD2	1:A:1069:ASP:C	2.59	0.44
1:B:960:ILE:HD11	1:B:967:VAL:HG23	1.99	0.44
1:B:1060:PHE:CE1	1:B:1108:LEU:HB3	2.52	0.44
3:N:45:ARG:HG3	3:N:45:ARG:HH11	1.82	0.44
1:A:922:ASP:HA	1:A:1110:SER:H	1.82	0.44
1:A:950:PHE:HB2	1:A:1094:LEU:HG	2.00	0.44
1:A:1068:TYR:OH	1:A:1087:LYS:HD3	2.17	0.44
2:H:190:TRP:CG	2:H:191:PRO:HA	2.52	0.44
1:B:972:TYR:HE2	1:B:1035:ILE:HD11	1.81	0.44
3:M:113:PRO:HB3	3:M:139:PHE:HB3	2.00	0.44
3:M:116:SER:O	3:M:134:CYS:HA	2.17	0.44
3:N:33:LEU:HG	3:N:71:TYR:CG	2.53	0.44
3:L:108:ARG:HG2	3:L:109:ALA:H	1.79	0.44
2:I:9:PRO:HG2	2:I:10:GLY:H	1.82	0.44
3:M:61:ARG:NH1	3:M:82:ASP:OD2	2.50	0.44
1:C:963:ARG:CZ	3:N:30:GLY:HA2	2.47	0.44
3:N:89:GLN:NE2	3:N:90:ASN:O	2.50	0.44
1:A:983:ASN:ND2	1:A:983:ASN:N	2.65	0.44
2:I:24:VAL:HG12	2:I:25:SER:N	2.33	0.44
2:J:13:ALA:O	2:J:16:GLN:HB2	2.18	0.44
2:J:24:VAL:CG1	2:J:25:SER:N	2.80	0.44
3:N:5:THR:HG23	7:N:235:HOH:O	2.18	0.44
1:B:950:PHE:HB2	1:B:1094:LEU:HG	1.99	0.43
3:M:14:SER:CB	3:M:17:GLN:NE2	2.81	0.43
2:J:66:ARG:HH21	2:J:82:LEU:HG	1.83	0.43
3:N:38:GLN:O	3:N:84:ALA:HB1	2.18	0.43
1:A:1100:LEU:HG	1:A:1101:GLY:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:24:VAL:CG1	2:H:25:SER:N	2.81	0.43
2:H:51:ILE:HG23	2:H:51:ILE:O	2.18	0.43
3:M:34:ASN:OD1	3:M:49:TYR:HA	2.18	0.43
1:C:1094:LEU:HB3	1:C:1095:PRO:CD	2.47	0.43
1:B:984:VAL:O	1:B:985:VAL:C	2.60	0.43
3:L:10:SER:HA	3:L:103:LYS:O	2.18	0.43
2:J:59:TYR:HB2	2:J:64:ILE:HG12	2.00	0.43
2:J:127:GLU:HG3	2:J:212:ILE:HG23	2.00	0.43
1:A:1046:ASP:OD1	1:A:1075:ILE:HD13	2.18	0.43
2:H:157:ASN:HA	7:H:628:HOH:O	2.18	0.43
2:I:24:VAL:CG1	2:I:25:SER:N	2.81	0.43
2:J:35:SER:O	2:J:95:CYS:HA	2.19	0.43
2:H:9:PRO:HG2	2:H:10:GLY:H	1.82	0.43
1:A:1023:HIS:HE1	3:N:81:GLU:OE2	2.02	0.43
1:A:1090:ARG:HG3	1:A:1092:GLU:HB2	1.99	0.43
1:B:948:LYS:O	1:B:952:LYS:HG3	2.19	0.43
2:J:68:SER:OG	2:J:81:LYS:HB2	2.17	0.43
3:N:130:ALA:HB3	3:N:181:LEU:O	2.19	0.43
1:A:945:ASP:HB3	7:A:509:HOH:O	2.19	0.43
2:H:6:GLN:HE21	2:H:6:GLN:N	2.12	0.43
2:H:121:PRO:CB	2:H:147:TYR:HB3	2.34	0.43
2:H:75:ARG:O	2:H:77:GLN:HG3	2.19	0.43
3:L:14:SER:H	3:L:17:GLN:NE2	2.17	0.43
3:M:44:VAL:HG22	7:M:312:HOH:O	2.18	0.43
1:C:936:SER:C	1:C:938:SER:H	2.26	0.43
1:A:1080:ALA:O	1:A:1083:SER:HB2	2.19	0.42
2:H:157:ASN:CB	2:H:160:SER:HB3	2.49	0.42
2:H:208:VAL:CG1	2:H:209:ASP:N	2.81	0.42
3:N:181:LEU:HD23	3:N:181:LEU:H	1.83	0.42
3:M:115:THR:HA	3:M:135:PHE:O	2.20	0.42
1:C:983:ASN:N	1:C:983:ASN:ND2	2.65	0.42
1:A:985:VAL:HG11	3:L:96:TRP:HZ2	1.83	0.42
2:H:38:ARG:HG3	2:H:46:GLU:HB2	2.01	0.42
2:H:126:LEU:HB2	2:H:141:GLY:C	2.44	0.42
2:H:202:PRO:C	2:H:204:SER:N	2.76	0.42
2:I:189:THR:O	2:I:193:GLN:HB2	2.19	0.42
1:C:983:ASN:N	1:C:983:ASN:HD22	2.16	0.42
1:C:1067:ARG:O	1:C:1067:ARG:HG3	2.19	0.42
1:A:962:PRO:HB3	3:L:29:ILE:O	2.19	0.42
2:H:14:PRO:O	2:H:85:LEU:HB2	2.19	0.42
2:H:129:VAL:HG11	7:L:1160:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1026:ARG:HA	1:C:1027:PRO:HD3	1.86	0.42
2:J:121:PRO:HB3	2:J:147:TYR:CB	2.42	0.42
3:N:119:PRO:HG3	3:N:209:PHE:CE2	2.54	0.42
2:H:4:LEU:N	2:H:4:LEU:HD12	2.34	0.42
2:H:149:PRO:HD2	2:H:203:ALA:HB1	2.00	0.42
2:H:202:PRO:O	2:H:204:SER:N	2.53	0.42
1:C:1105:LEU:HD22	1:C:1105:LEU:H	1.84	0.42
2:H:121:PRO:HB3	2:H:147:TYR:CB	2.34	0.42
1:B:968:SER:HB2	1:B:982:TRP:CZ3	2.54	0.42
1:B:1018:LEU:HD11	1:B:1033:VAL:CG2	2.50	0.42
1:B:1089:GLN:HE21	1:B:1089:GLN:HB3	1.56	0.42
3:M:61:ARG:CG	3:M:76:SER:O	2.68	0.42
3:M:207:LYS:HD3	3:M:207:LYS:HA	1.89	0.42
2:J:60:HIS:O	2:J:61:SER:C	2.63	0.42
2:H:127:GLU:HG3	2:H:212:ILE:HG23	2.02	0.42
2:I:92:THR:HA	2:I:110:SER:HA	2.01	0.42
3:M:118:PHE:HA	3:M:119:PRO:HD3	1.93	0.42
1:C:1089:GLN:HG2	7:C:616:HOH:O	2.18	0.42
1:A:1096:THR:HA	1:A:1099:THR:OG1	2.19	0.42
3:L:14:SER:H	3:L:17:GLN:CD	2.27	0.42
2:J:14:PRO:HD3	2:J:114:SER:C	2.43	0.42
2:H:67:LEU:HD11	2:H:80:LEU:CD2	2.49	0.42
3:N:4:MET:HE3	3:N:23:CYS:SG	2.60	0.42
3:L:14:SER:CB	3:L:17:GLN:NE2	2.83	0.41
1:B:940:PRO:O	1:B:943:TYR:HB2	2.20	0.41
3:M:44:VAL:HG12	3:M:45:ARG:N	2.35	0.41
2:J:80:LEU:HD22	2:J:81:LYS:H	1.84	0.41
3:N:41:ASP:OD1	3:N:43:THR:HG23	2.20	0.41
3:L:175:MET:O	3:L:175:MET:HG3	2.20	0.41
3:M:14:SER:HB2	3:M:17:GLN:HE21	1.85	0.41
3:M:37:GLN:HB2	3:M:47:LEU:HD11	2.01	0.41
1:A:1069:ASP:OD2	1:A:1071:ALA:HB3	2.20	0.41
3:L:136:LEU:HD12	3:L:136:LEU:N	2.36	0.41
1:B:963:ARG:CZ	3:M:30:GLY:HA2	2.51	0.41
2:J:60:HIS:NE2	3:N:1:ASP:OD2	2.47	0.41
3:L:144:ILE:HG22	3:L:163:TRP:CZ3	2.56	0.41
2:I:34:VAL:HG21	2:I:78:VAL:HG21	2.02	0.41
3:M:36:TYR:HB2	3:M:87:PHE:CE1	2.55	0.41
3:M:170:ASP:O	3:M:171:SER:HB2	2.20	0.41
2:H:3:LYS:C	2:H:4:LEU:HD12	2.46	0.41
2:H:38:ARG:CD	2:H:46:GLU:OE1	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:145:LYS:HE3	2:H:145:LYS:HB2	1.92	0.41
1:C:1062:ILE:HA	1:C:1086:VAL:O	2.21	0.41
2:J:34:VAL:HG11	2:J:78:VAL:HG21	2.02	0.41
2:H:201:HIS:HB3	2:H:206:THR:HB	2.03	0.41
2:I:14:PRO:O	2:I:15:SER:HB2	2.21	0.41
1:A:1074:ARG:HH11	1:A:1074:ARG:HG2	1.86	0.41
1:B:1102:ASN:O	1:B:1103:SER:C	2.64	0.41
3:N:115:THR:HA	3:N:135:PHE:O	2.20	0.41
2:H:140:LEU:HG	2:H:212:ILE:HG21	2.03	0.41
3:L:184:ASP:HB3	3:L:188:ARG:NH1	2.34	0.41
3:M:3:ALA:HB2	7:M:348:HOH:O	2.21	0.41
3:M:54:LEU:CD1	3:M:60:SER:HA	2.48	0.41
2:J:80:LEU:HD13	2:J:81:LYS:N	2.36	0.41
2:J:121:PRO:CB	2:J:147:TYR:HB3	2.41	0.41
2:J:125:LYS:HE2	3:N:123:GLU:OE2	2.21	0.41
2:J:127:GLU:HB2	2:J:215:ARG:HG3	2.03	0.41
2:I:179:LEU:C	2:I:179:LEU:HD23	2.46	0.41
3:M:160:LEU:HD12	3:M:160:LEU:HA	1.94	0.41
4:D:1:NAG:H62	4:D:2:FUC:C5	2.49	0.41
2:I:12:VAL:HG11	2:I:85:LEU:HD12	2.04	0.40
1:C:1033:VAL:CG1	1:C:1059:VAL:HG22	2.52	0.40
2:J:56:ASN:HA	4:F:1:NAG:C7	2.51	0.40
2:J:62:ALA:O	2:J:63:LEU:HD23	2.20	0.40
1:A:983:ASN:N	1:A:983:ASN:HD22	2.18	0.40
1:A:1069:ASP:CG	1:A:1072:GLN:HG3	2.47	0.40
1:C:944:PHE:CD1	1:C:947:MSE:HE3	2.56	0.40
3:N:33:LEU:C	3:N:33:LEU:HD13	2.46	0.40
1:A:1069:ASP:OD2	1:A:1071:ALA:N	2.50	0.40
3:L:29:ILE:HG12	7:L:1009:HOH:O	2.21	0.40
2:I:24:VAL:HG11	2:I:27:PHE:CE1	2.56	0.40
1:C:932:LEU:HD21	1:C:1010:ALA:HB1	2.04	0.40
1:C:1022:MSE:HE2	7:C:162:HOH:O	2.20	0.40
1:C:1096:THR:HA	1:C:1099:THR:OG1	2.21	0.40
3:N:167:ASP:OD2	3:N:170:ASP:HB2	2.20	0.40
1:C:925:GLN:OE1	1:C:925:GLN:HA	2.21	0.40
3:N:136:LEU:N	3:N:136:LEU:HD12	2.35	0.40
3:N:193:THR:OG1	3:N:208:SER:HB3	2.21	0.40
1:B:1056:ARG:HG2	1:B:1056:ARG:NH1	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	188/196 (96%)	182 (97%)	5 (3%)	1 (0%)	24	17
1	B	188/196 (96%)	180 (96%)	7 (4%)	1 (0%)	24	17
1	C	189/196 (96%)	176 (93%)	10 (5%)	3 (2%)	7	2
2	H	206/210 (98%)	189 (92%)	15 (7%)	2 (1%)	12	5
2	I	206/210 (98%)	194 (94%)	11 (5%)	1 (0%)	24	17
2	J	206/210 (98%)	184 (89%)	21 (10%)	1 (0%)	24	17
3	L	209/211 (99%)	204 (98%)	5 (2%)	0	100	100
3	M	209/211 (99%)	206 (99%)	3 (1%)	0	100	100
3	N	209/211 (99%)	202 (97%)	7 (3%)	0	100	100
All	All	1810/1851 (98%)	1717 (95%)	84 (5%)	9 (0%)	24	17

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1100	LEU
1	C	1102	ASN
2	J	15	SER
1	B	1100	LEU
2	I	41	PRO
1	C	1100	LEU
2	H	14	PRO
2	H	203	ALA
1	C	1079	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	157/155 (101%)	153 (98%)	4 (2%)	42	40
1	B	157/155 (101%)	154 (98%)	3 (2%)	50	50
1	C	157/155 (101%)	154 (98%)	3 (2%)	50	50
2	H	183/183 (100%)	175 (96%)	8 (4%)	25	18
2	I	183/183 (100%)	173 (94%)	10 (6%)	19	12
2	J	183/183 (100%)	179 (98%)	4 (2%)	45	45
3	L	186/186 (100%)	182 (98%)	4 (2%)	45	45
3	M	186/186 (100%)	182 (98%)	4 (2%)	45	45
3	N	186/186 (100%)	182 (98%)	4 (2%)	45	45
All	All	1578/1572 (100%)	1534 (97%)	44 (3%)	38	35

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	945	ASP
1	A	983	ASN
1	A	1036	LEU
1	A	1067	ARG
2	H	6	GLN
2	H	14	PRO
2	H	80	LEU
2	H	102	MET
2	H	140	LEU
2	H	149	PRO
2	H	198	ASN
2	H	210	LYS
3	L	79	GLU
3	L	90	ASN
3	L	107	LYS
3	L	175	MET
1	B	1089	GLN
1	B	1108	LEU
1	B	1109	CYS
2	I	6	GLN
2	I	41	PRO
2	I	80	LEU
2	I	107	GLN

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Mol	Chain	Res	Type
2	I	125	LYS
2	I	127	GLU
2	I	140	LEU
2	I	143	LEU
2	I	150	GLU
2	I	151	PRO
3	M	8	THR
3	M	90	ASN
3	M	147	LYS
3	M	163	TRP
1	C	923	CYS
1	C	1079	PRO
1	C	1109	CYS
2	J	6	GLN
2	J	43	LYS
2	J	143	LEU
2	J	151	PRO
3	N	79	GLU
3	N	90	ASN
3	N	160	LEU
3	N	175	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	959	ASN
2	H	6	GLN
2	H	39	GLN
2	H	77	GLN
2	H	98	ASN
2	H	173	GLN
3	L	17	GLN
3	L	38	GLN
3	L	94	ASN
3	L	137	ASN
3	L	189	HIS
3	L	210	ASN
1	B	1006	GLN
1	B	1055	ASN
1	B	1089	GLN
2	I	6	GLN
2	I	107	GLN

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Mol	Chain	Res	Type
3	M	17	GLN
3	M	37	GLN
3	M	77	ASN
3	M	156	GLN
3	M	189	HIS
1	C	959	ASN
1	C	971	GLN
1	C	990	HIS
1	C	1023	HIS
1	C	1084	ASN
2	J	6	GLN
2	J	77	GLN
2	J	173	GLN
3	N	137	ASN
3	N	156	GLN

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 ⓘ

6 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$
4	NAG	D	1	2,4	14,14,15	0.58	0	17,19,21	0.82	1 (5%)
4	FUC	D	2	4	10,10,11	0.57	0	14,14,16	0.74	0
4	NAG	E	1	2,4	14,14,15	0.62	0	17,19,21	0.90	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FUC	E	2	4	10,10,11	0.64	0	14,14,16	0.64	0
4	NAG	F	1	2,4	14,14,15	0.57	0	17,19,21	0.73	0
4	FUC	F	2	4	10,10,11	0.71	0	14,14,16	0.72	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
4	NAG	D	1	2,4	-	1/6/23/26	0/1/1/1
4	FUC	D	2	4	-	-	0/1/1/1
4	NAG	E	1	2,4	-	0/6/23/26	0/1/1/1
4	FUC	E	2	4	-	-	0/1/1/1
4	NAG	F	1	2,4	-	0/6/23/26	0/1/1/1
4	FUC	F	2	4	-	-	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(°)
4	E	1	NAG	C2-N2-C7	-2.41	119.67	122.90
4	D	1	NAG	C2-N2-C7	-2.11	120.07	122.90

There are no chirality outliers.

All (1) torsion outliers are listed below:

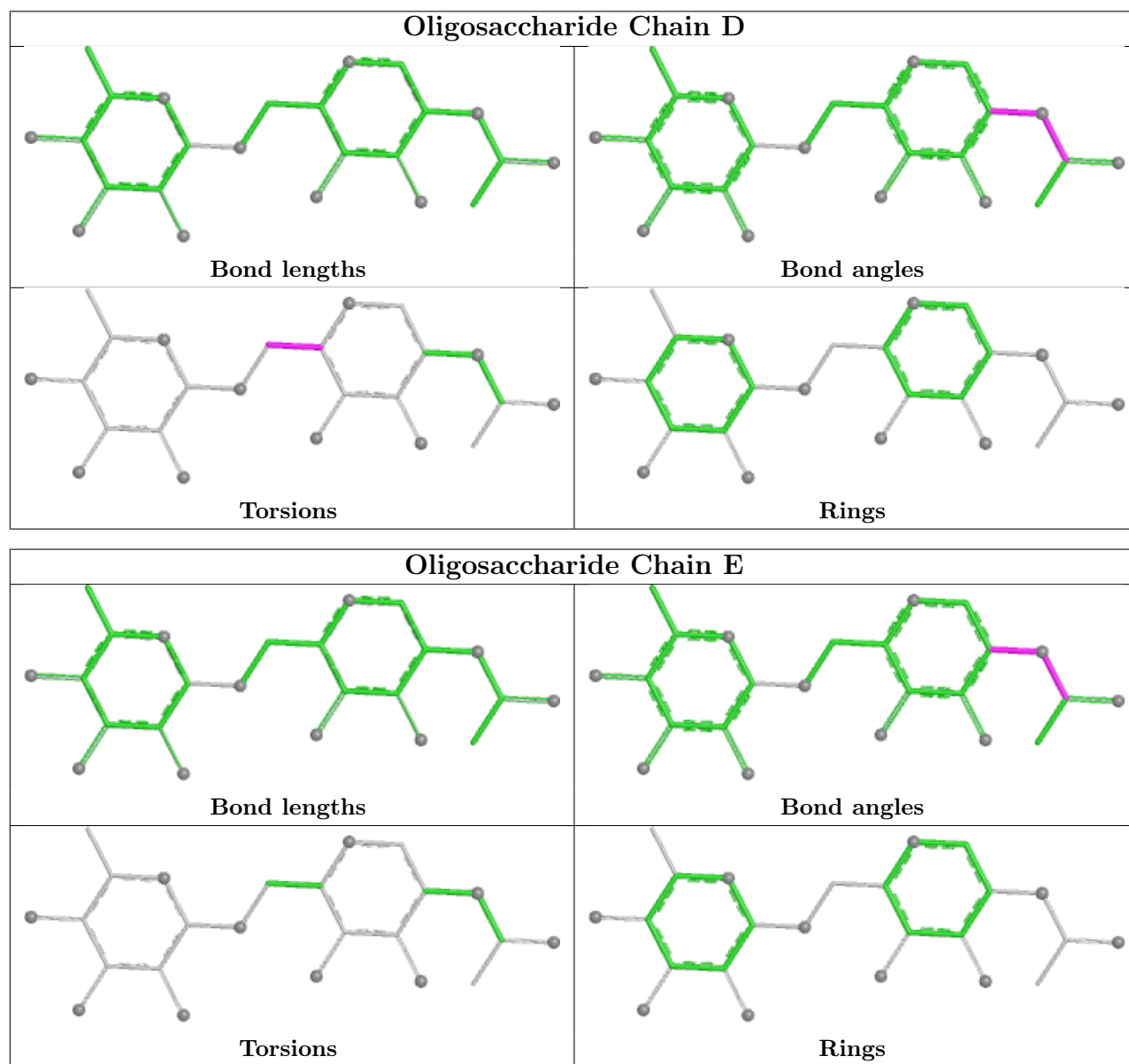
Mol	Chain	Res	Type	Atoms
4	D	1	NAG	C4-C5-C6-O6

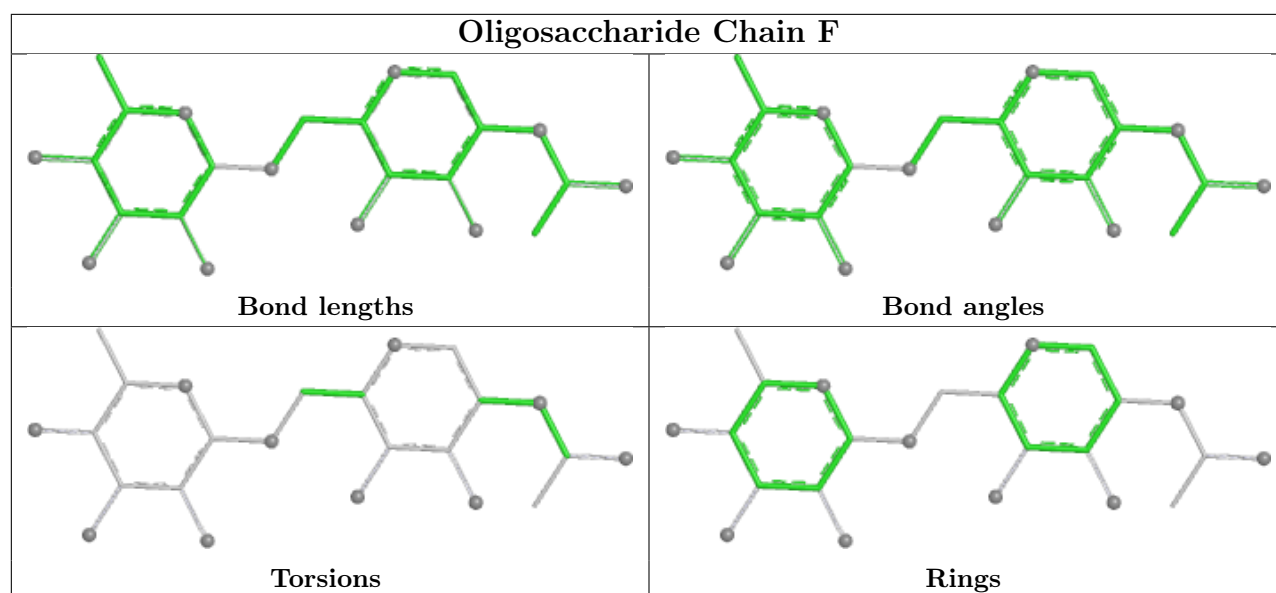
There are no ring outliers.

6 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	2	FUC	6	0
4	E	2	FUC	5	0
4	F	1	NAG	6	0
4	D	1	NAG	7	0
4	E	1	NAG	6	0
4	F	2	FUC	6	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	I	553	-	14,14,15	0.58	0	17,19,21	0.69	0
6	CAC	L	1001	-	2,4,4	1.61	1 (50%)	4,6,6	1.30	1 (25%)
5	NAG	H	553	-	14,14,15	0.64	0	17,19,21	0.64	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	I	553	-	-	3/6/23/26	0/1/1/1
5	NAG	H	553	-	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L	1001	CAC	AS-C1	2.03	1.95	1.90

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
6	L	1001	CAC	O1-AS-C1	2.11	114.13	111.50

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	I	553	NAG	C1-C2-N2-C7
5	H	553	NAG	C8-C7-N2-C2
5	H	553	NAG	O7-C7-N2-C2
5	I	553	NAG	C3-C2-N2-C7
5	I	553	NAG	C8-C7-N2-C2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	I	553	NAG	1	0
5	H	553	NAG	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	J	1
2	H	1
2	I	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	J	129:VAL	C	136:SER	N	14.08
1	H	129:VAL	C	136:SER	N	14.07
1	I	129:VAL	C	136:SER	N	13.60

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	186/196 (94%)	0.43	14 (7%)	20 19	21, 33, 56, 61	0
1	B	186/196 (94%)	0.58	10 (5%)	31 30	24, 35, 56, 61	0
1	C	187/196 (95%)	0.98	26 (13%)	6 5	26, 43, 56, 61	0
2	H	210/210 (100%)	0.80	19 (9%)	15 14	22, 40, 54, 57	0
2	I	210/210 (100%)	0.49	10 (4%)	35 35	24, 36, 48, 59	0
2	J	210/210 (100%)	1.45	44 (20%)	2 2	30, 50, 57, 60	0
3	L	211/211 (100%)	0.01	3 (1%)	73 73	18, 28, 46, 53	0
3	M	211/211 (100%)	0.29	9 (4%)	40 39	23, 33, 45, 55	0
3	N	211/211 (100%)	0.55	14 (6%)	24 23	25, 37, 52, 56	0
All	All	1822/1851 (98%)	0.62	149 (8%)	17 17	18, 37, 54, 61	0

All (149) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	127	GLU	5.0
1	B	923	CYS	4.8
1	C	1080	ALA	4.6
1	B	1108	LEU	4.3
1	C	1081	GLY	4.2
1	C	939	PHE	3.9
2	H	129	VAL	3.8
2	I	127	GLU	3.8
1	A	1100	LEU	3.7
2	J	7	SER	3.6
1	C	923	CYS	3.6
2	J	16	GLN	3.5
2	J	127	GLU	3.5
2	H	125	LYS	3.5
1	C	1109	CYS	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	922	ASP	3.4
1	B	1109	CYS	3.4
2	J	114	SER	3.4
1	B	1106	HIS	3.3
2	J	24	VAL	3.3
2	J	216	GLY	3.3
3	M	170	ASP	3.3
1	A	921	PRO	3.3
1	A	1105	LEU	3.3
3	M	163	TRP	3.2
1	C	937	SER	3.2
2	J	123	VAL	3.2
2	J	22	CYS	3.2
3	N	132	VAL	3.2
2	J	42	GLY	3.2
2	H	64	ILE	3.1
1	C	938	SER	3.1
2	H	136	SER	3.1
2	J	15	SER	3.1
2	H	203	ALA	3.1
2	H	81	LYS	3.0
1	B	1100	LEU	3.0
2	J	129	VAL	3.0
2	J	87	THR	3.0
2	J	64	ILE	3.0
1	A	1106	HIS	3.0
3	N	163	TRP	3.0
1	C	1110	SER	2.9
3	N	170	ASP	2.9
3	L	128	GLY	2.9
2	J	5	VAL	2.9
1	C	920	ALA	2.9
2	I	129	VAL	2.9
2	H	42	GLY	2.9
1	A	923	CYS	2.8
2	J	67	LEU	2.8
2	H	197	CYS	2.8
2	H	14	PRO	2.8
1	A	1042	VAL	2.7
2	J	1	ASP	2.7
1	C	1100	LEU	2.7
2	J	189	THR	2.7

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Mol	Chain	Res	Type	RSRZ
3	N	182	THR	2.7
1	A	1109	CYS	2.7
1	C	1105	LEU	2.7
2	J	136	SER	2.7
2	J	54	ASP	2.7
1	A	1101	GLY	2.7
1	B	1105	LEU	2.6
1	C	997	VAL	2.6
1	C	1089	GLN	2.6
2	J	86	HIS	2.6
1	A	1099	THR	2.6
1	C	1099	THR	2.6
3	M	45	ARG	2.6
1	C	960	ILE	2.6
2	I	216	GLY	2.6
3	N	168	SER	2.6
2	J	79	PHE	2.5
2	J	88	ASP	2.5
3	N	184	ASP	2.5
2	J	78	VAL	2.5
3	N	128	GLY	2.5
2	I	205	SER	2.5
3	N	125	LEU	2.5
2	J	125	LYS	2.5
2	J	138	VAL	2.5
2	H	118	THR	2.5
1	B	921	PRO	2.5
1	C	1079	PRO	2.5
2	J	203	ALA	2.5
2	I	44	GLY	2.5
2	J	83	ASN	2.5
1	C	922	ASP	2.4
2	H	173	GLN	2.4
1	C	1067	ARG	2.4
1	A	1108	LEU	2.4
3	N	108	ARG	2.4
2	H	160	SER	2.4
2	I	136	SER	2.4
2	J	156	TRP	2.4
2	J	85	LEU	2.3
1	A	1081	GLY	2.3
2	J	70	SER	2.3

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Mol	Chain	Res	Type	RSRZ
3	N	208	SER	2.3
2	H	87	THR	2.3
3	M	81	GLU	2.3
3	M	77	ASN	2.3
1	A	1074	ARG	2.3
2	H	15	SER	2.3
2	J	205	SER	2.3
2	H	16	GLN	2.3
2	I	1	ASP	2.3
2	J	117	GLU	2.3
2	J	69	ILE	2.3
2	J	81	LYS	2.3
2	J	206	THR	2.2
1	B	922	ASP	2.2
2	H	116	ALA	2.2
2	J	82	LEU	2.2
1	C	943	TYR	2.2
3	M	79	GLU	2.2
2	H	119	THR	2.2
2	J	23	THR	2.2
1	C	996	ASP	2.2
2	J	44	GLY	2.2
1	C	1070	ALA	2.2
2	H	120	ALA	2.2
1	C	1088	LEU	2.2
2	J	163	SER	2.2
3	M	87	PHE	2.2
2	J	215	ARG	2.2
3	M	103	LYS	2.2
3	L	130	ALA	2.1
1	A	1067	ARG	2.1
2	J	126	LEU	2.1
2	I	64	ILE	2.1
1	C	1056	ARG	2.1
2	J	80	LEU	2.1
2	I	88	ASP	2.1
2	J	196	THR	2.1
1	C	944	PHE	2.1
1	B	1110	SER	2.1
3	L	127	SER	2.1
3	N	77	ASN	2.1
1	B	1067	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
2	J	109	THR	2.1
3	N	116	SER	2.1
3	M	11	LEU	2.0
1	C	1098	VAL	2.0
2	I	208	VAL	2.0
3	N	148	TRP	2.0
3	N	190	ASN	2.0
1	C	1094	LEU	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](#)

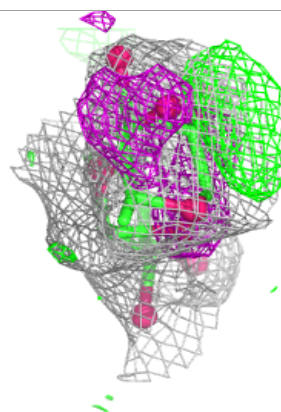
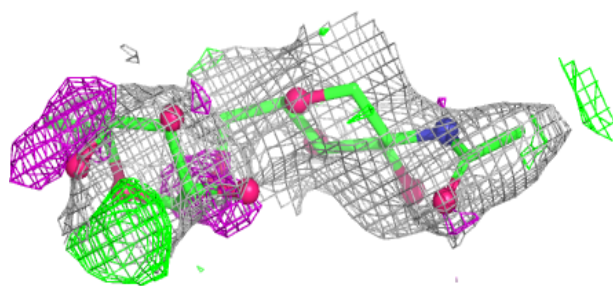
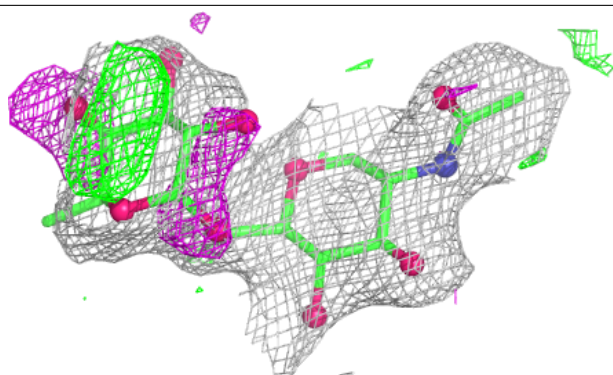
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	D	1	14/15	-	-	46,49,53,56	0
4	FUC	D	2	10/11	-	-	55,57,57,57	0
4	NAG	E	1	14/15	-	-	47,50,54,56	0
4	FUC	E	2	10/11	-	-	58,58,59,59	0
4	NAG	F	1	14/15	-	-	56,57,58,59	0
4	FUC	F	2	10/11	-	-	60,60,61,61	0

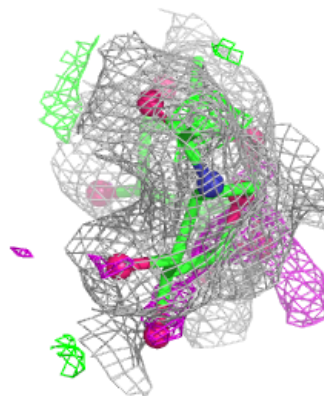
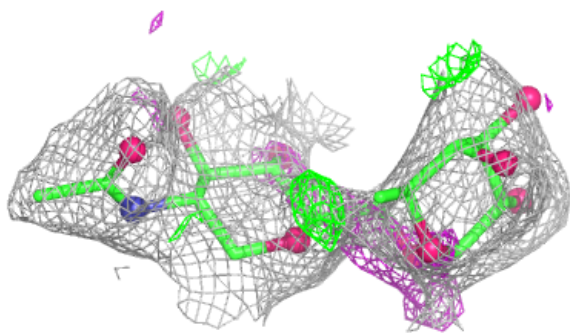
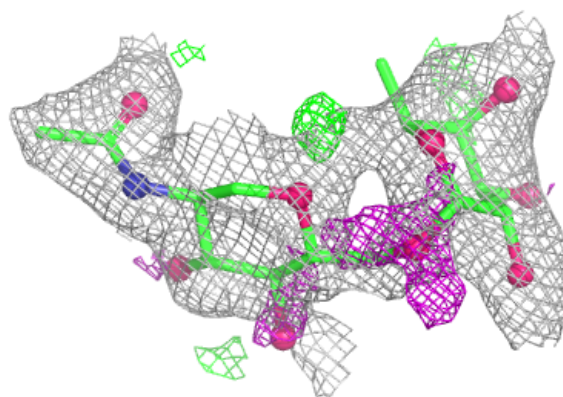
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

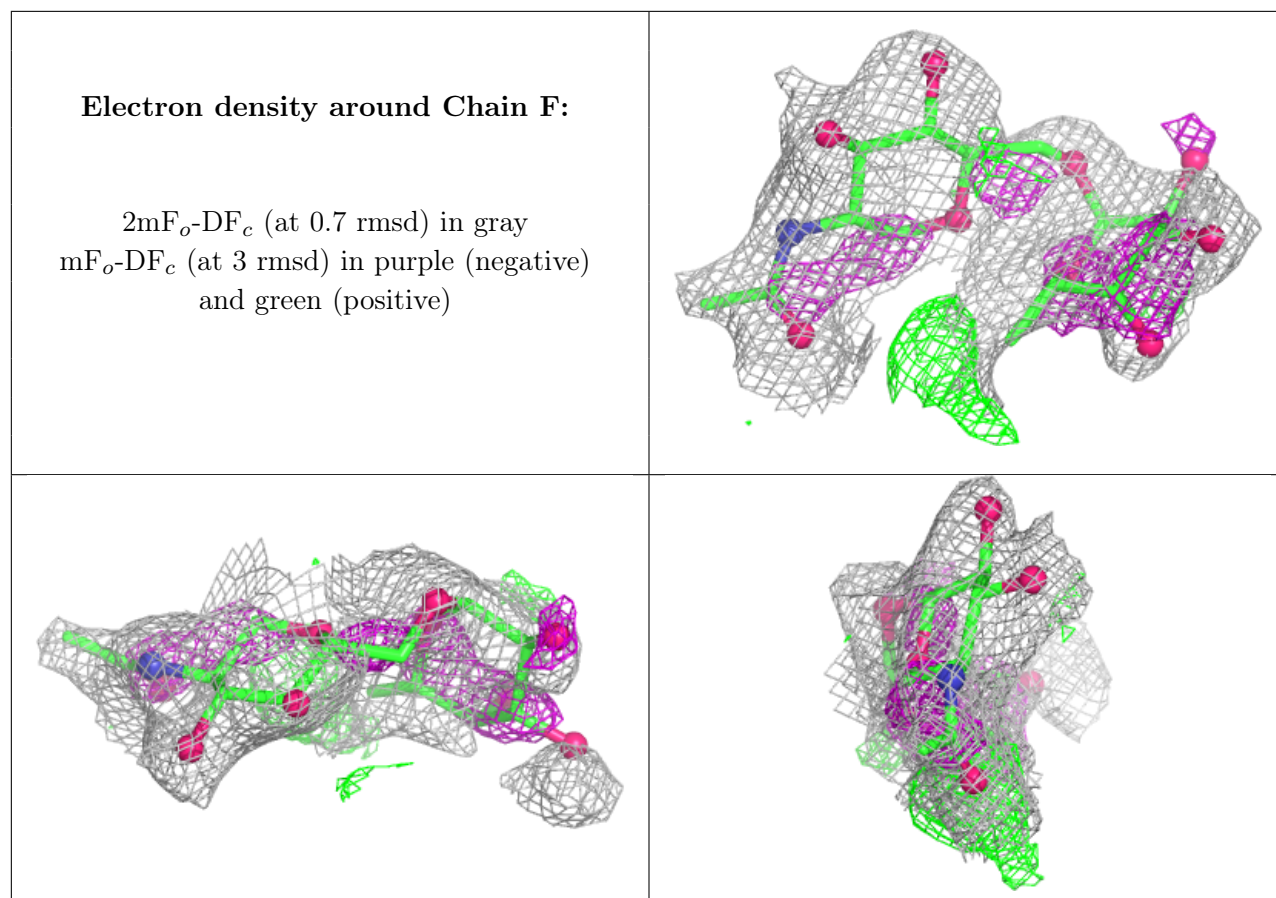
Electron density around Chain D:

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

**Electron density around Chain E:**

$2mF_o-DF_c$ (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	I	553	14/15	0.51	0.19	59,60,62,62	0
5	NAG	H	553	14/15	0.56	0.16	57,58,59,59	0
6	CAC	L	1001	5/5	0.96	0.19	43,44,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.