Full wwPDB X-ray Structure Validation Report (i)

Jan 2, 2024 – 09:57 am GMT

PDB ID	:	4UTB
Title	:	Crystal structure of dengue 2 virus envelope glycoprotein in complex with the Fab fragment of the broadly neutralizing human antibody EDE2 A11
Authors	:	Rouvinski, A.; Guardado-Calvo, P.; Barba-Spaeth, G.; Duquerroy, S.; Vaney, M.C.; Rey, F.A.
Deposited on	:	2014-07-18
Resolution	:	3.85 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 (i) 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 (i)) were used in the production of this report:

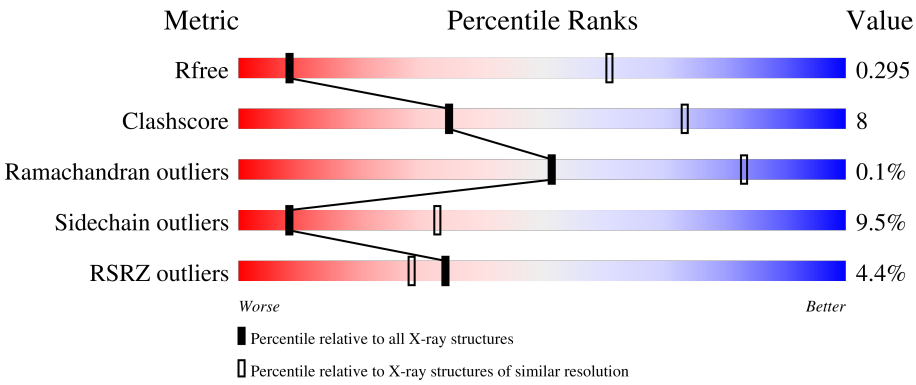
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.85 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1048 (4.10-3.62)
Clashscore	141614	1015 (4.08-3.64)
Ramachandran outliers	138981	1069 (4.10-3.62)
Sidechain outliers	138945	1062 (4.10-3.62)
RSRZ outliers	127900	1206 (4.12-3.60)

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 <=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	422	4% 68% 21% • 7%
1	B	422	2% 69% 20% • 8%
2	H	283	% 68% 10% • 20%
2	I	283	7% 60% 11% • 28%
3	L	218	3% 87% 11% •

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	M	218	7% 84% 11%
4	C	6	83% 17%
4	E	6	17% 50% 33%
5	D	2	50% 50%

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	MAN	C	4	-	-	-	X
4	FUC	C	6	-	-	-	X
5	NAG	D	1	-	-	-	X
5	NAG	D	2	-	-	-	X

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 12714 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called ENVELOPE GLYCOPROTEIN E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	391	3045	1919	524	578	24	0	0	0
1	B	390	3036	1914	522	576	24	0	0	0

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1392	LEU	-	expression tag	UNP Q68Y26
A	1393	ARG	-	expression tag	UNP Q68Y26
A	1394	PRO	-	expression tag	UNP Q68Y26
A	1395	LEU	-	expression tag	UNP Q68Y26
A	1396	GLU	-	expression tag	UNP Q68Y26
A	1397	SER	-	expression tag	UNP Q68Y26
A	1398	ARG	-	expression tag	UNP Q68Y26
A	1399	GLY	-	expression tag	UNP Q68Y26
A	1400	PRO	-	expression tag	UNP Q68Y26
A	1401	PHE	-	expression tag	UNP Q68Y26
A	1402	GLU	-	expression tag	UNP Q68Y26
A	1403	GLY	-	expression tag	UNP Q68Y26
A	1404	LYS	-	expression tag	UNP Q68Y26
A	1405	PRO	-	expression tag	UNP Q68Y26
A	1406	ILE	-	expression tag	UNP Q68Y26
A	1407	PRO	-	expression tag	UNP Q68Y26
A	1408	ASN	-	expression tag	UNP Q68Y26
A	1409	PRO	-	expression tag	UNP Q68Y26
A	1410	LEU	-	expression tag	UNP Q68Y26
A	1411	LEU	-	expression tag	UNP Q68Y26
A	1412	GLY	-	expression tag	UNP Q68Y26
A	1413	LEU	-	expression tag	UNP Q68Y26
A	1414	ASP	-	expression tag	UNP Q68Y26
A	1415	SER	-	expression tag	UNP Q68Y26
A	1416	THR	-	expression tag	UNP Q68Y26

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	1417	ARG	-	expression tag	UNP Q68Y26
A	1418	THR	-	expression tag	UNP Q68Y26
A	1419	GLY	-	expression tag	UNP Q68Y26
A	1420	HIS	-	expression tag	UNP Q68Y26
A	1421	HIS	-	expression tag	UNP Q68Y26
A	1422	HIS	-	expression tag	UNP Q68Y26
A	118	LYS	MET	conflict	UNP Q68Y26
B	1392	LEU	-	expression tag	UNP Q68Y26
B	1393	ARG	-	expression tag	UNP Q68Y26
B	1394	PRO	-	expression tag	UNP Q68Y26
B	1395	LEU	-	expression tag	UNP Q68Y26
B	1396	GLU	-	expression tag	UNP Q68Y26
B	1397	SER	-	expression tag	UNP Q68Y26
B	1398	ARG	-	expression tag	UNP Q68Y26
B	1399	GLY	-	expression tag	UNP Q68Y26
B	1400	PRO	-	expression tag	UNP Q68Y26
B	1401	PHE	-	expression tag	UNP Q68Y26
B	1402	GLU	-	expression tag	UNP Q68Y26
B	1403	GLY	-	expression tag	UNP Q68Y26
B	1404	LYS	-	expression tag	UNP Q68Y26
B	1405	PRO	-	expression tag	UNP Q68Y26
B	1406	ILE	-	expression tag	UNP Q68Y26
B	1407	PRO	-	expression tag	UNP Q68Y26
B	1408	ASN	-	expression tag	UNP Q68Y26
B	1409	PRO	-	expression tag	UNP Q68Y26
B	1410	LEU	-	expression tag	UNP Q68Y26
B	1411	LEU	-	expression tag	UNP Q68Y26
B	1412	GLY	-	expression tag	UNP Q68Y26
B	1413	LEU	-	expression tag	UNP Q68Y26
B	1414	ASP	-	expression tag	UNP Q68Y26
B	1415	SER	-	expression tag	UNP Q68Y26
B	1416	THR	-	expression tag	UNP Q68Y26
B	1417	ARG	-	expression tag	UNP Q68Y26
B	1418	THR	-	expression tag	UNP Q68Y26
B	1419	GLY	-	expression tag	UNP Q68Y26
B	1420	HIS	-	expression tag	UNP Q68Y26
B	1421	HIS	-	expression tag	UNP Q68Y26
B	1422	HIS	-	expression tag	UNP Q68Y26
B	118	LYS	MET	conflict	UNP Q68Y26

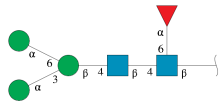
- Molecule 2 is a protein called BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	225	1718	1088	291	332	7	0	0	0
2	I	203	1573	999	267	300	7	0	0	0

- Molecule 3 is a protein called BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	L	213	1582	987	265	324	6	0	0	0
3	M	210	1563	975	262	320	6	0	0	0

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

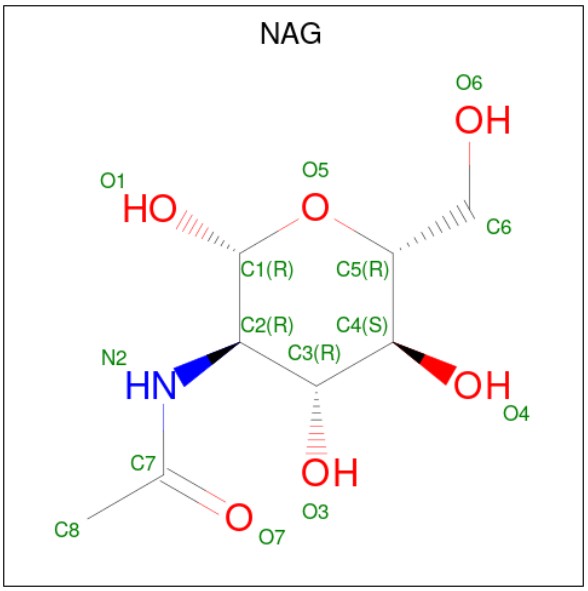


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	C	6	71	40	2	29	0	0	0
4	E	6	71	40	2	29	0	0	0

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

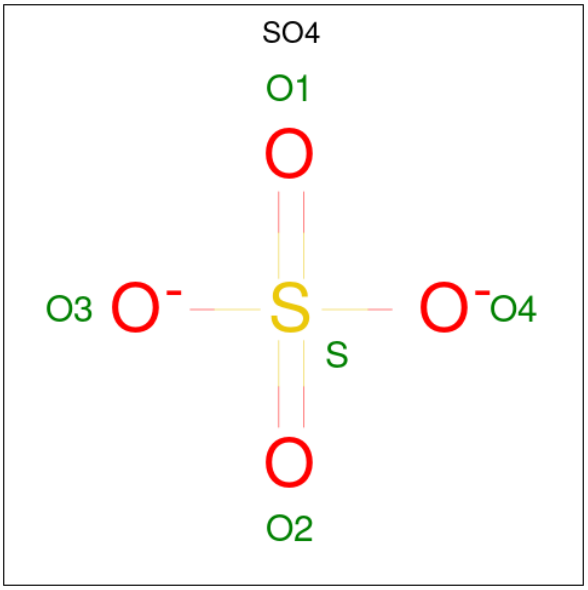
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	D	2	28	16	2	10	0	0	0

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
7	H	1	5	4	1	0	0
7	I	1	5	4	1	0	0

- Molecule 8 is water.

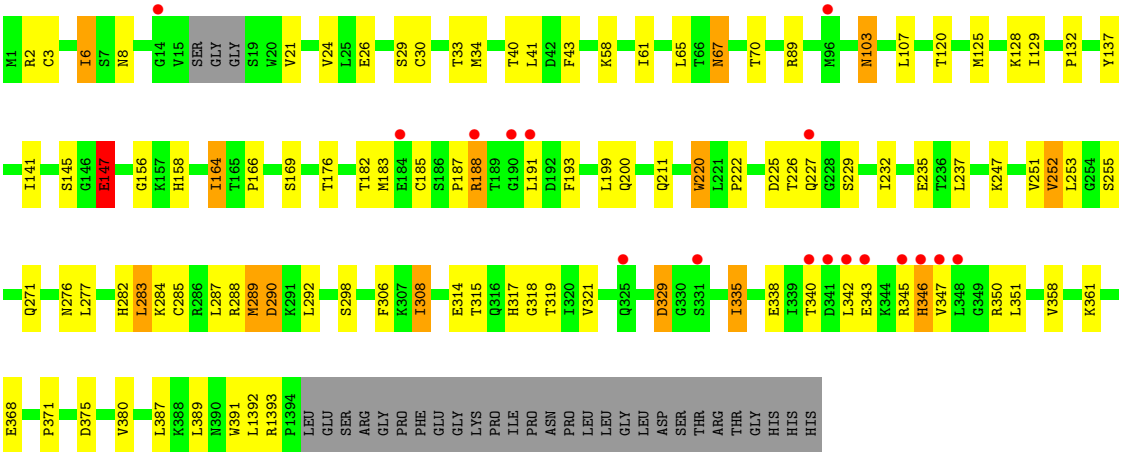
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	I	2	2	2	0	0
			Total	O		
8	M	1	1	1	0	0

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.

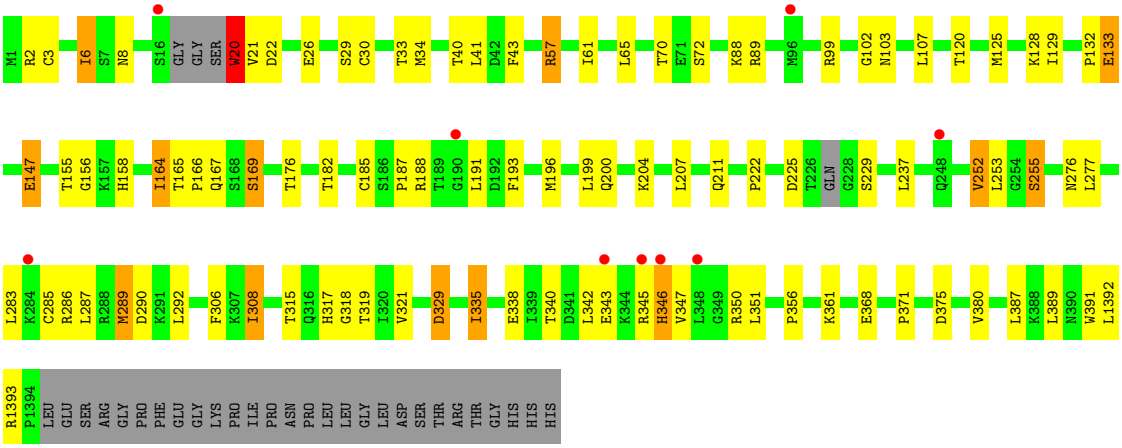
• Molecule 1: ENVELOPE GLYCOPROTEIN E

Chain A: 4% 68% 21% • 7%



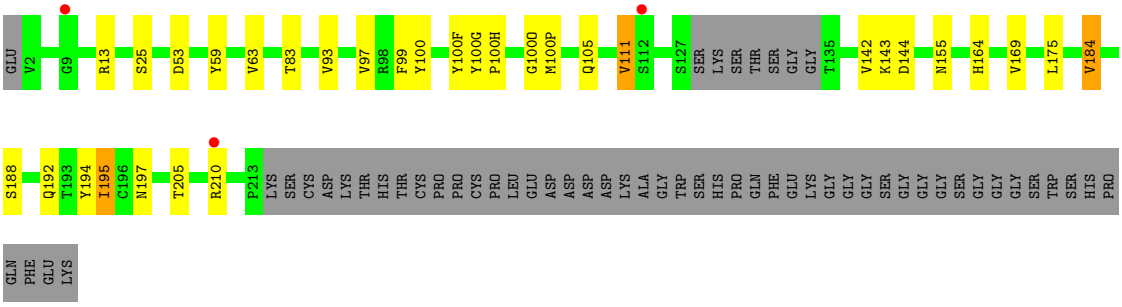
• Molecule 1: ENVELOPE GLYCOPROTEIN E

Chain B: 2% 69% 20% • 8%



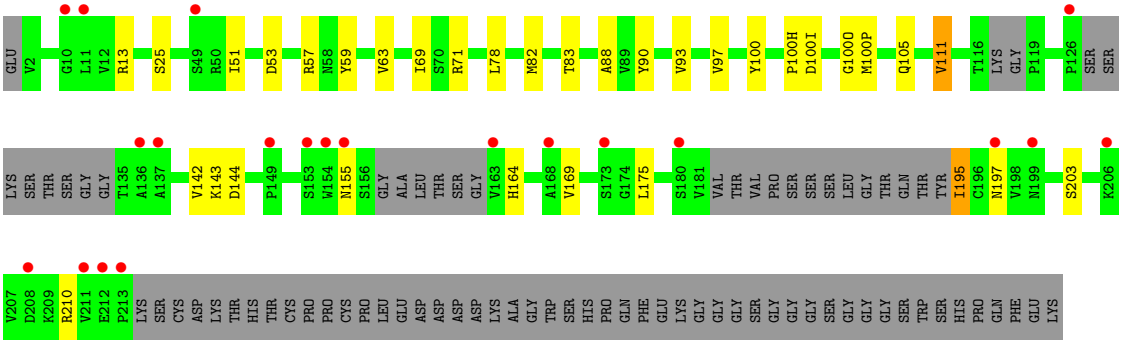
• Molecule 2: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11

Chain H: % 68% 10% • 20%



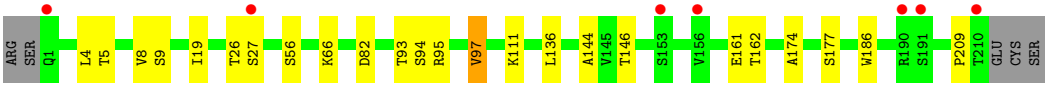
• Molecule 2: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11

Chain I:



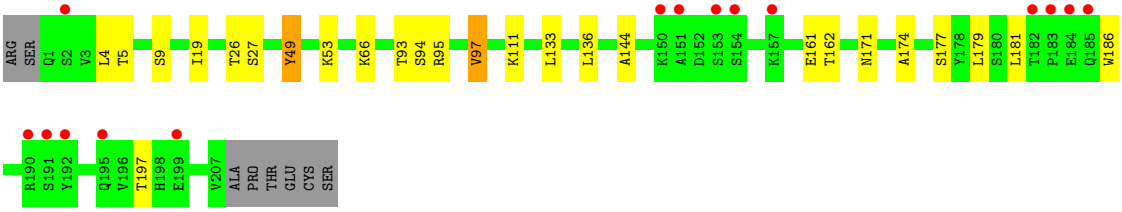
• Molecule 3: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11

Chain L:



• Molecule 3: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11

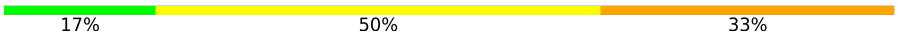
Chain M:



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

Chain C:

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

Chain E: 

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

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	58.78Å 181.85Å 204.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.85 29.74 – 3.85	Depositor EDS
% Data completeness (in resolution range)	97.9 (20.00-3.85) 97.9 (29.74-3.85)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.56 (at 3.86Å)	Xtriage
Refinement program	BUSTER 2.11.4	Depositor
R, R_{free}	0.230 , 0.257 0.278 , 0.295	Depositor DCC
R_{free} test set	1074 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	109.8	Xtriage
Anisotropy	0.530	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 124.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.39$, $< L^2 > = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	12714	wwPDB-VP
Average B, all atoms (Å^2)	129.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.23% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, SO4, NAG, FUC, BMA

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| >5$	RMSZ	$\#\|Z\| >5$
1	A	0.46	0/3106	0.76	2/4199 (0.0%)
1	B	0.42	0/3096	0.72	1/4184 (0.0%)
2	H	0.36	0/1764	0.64	0/2406
2	I	0.38	0/1614	0.66	0/2195
3	L	0.38	0/1618	0.65	0/2209
3	M	0.40	0/1598	0.67	0/2180
All	All	0.41	0/12796	0.70	3/17373 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	67	ASN	CA-CB-CG	6.74	128.22	113.40
1	B	20	TRP	CA-CB-CG	5.28	123.74	113.70
1	A	103	ASN	N-CA-CB	-5.11	101.40	110.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3045	0	3044	68	0
1	B	3036	0	3035	65	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	1718	0	1644	25	0
2	I	1573	0	1498	36	0
3	L	1582	0	1553	12	0
3	M	1563	0	1534	23	0
4	C	71	0	61	2	0
4	E	71	0	61	2	0
5	D	28	0	25	0	0
6	B	14	0	13	0	0
7	H	5	0	0	0	0
7	I	5	0	0	0	0
8	I	2	0	0	0	0
8	M	1	0	0	0	0
All	All	12714	0	12468	209	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 8.

All (209) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ILE:CD1	1:A:358:VAL:HG23	1.64	1.25
1:A:335:ILE:HD11	1:A:358:VAL:CG2	1.70	1.19
1:B:57:ARG:HG3	1:B:57:ARG:HH11	1.17	1.02
1:B:34:MET:HG2	1:B:40:THR:HG22	1.49	0.94
1:A:371:PRO:HG2	1:A:391:TRP:CZ3	2.04	0.92
1:A:34:MET:HG2	1:A:40:THR:HG22	1.53	0.90
1:B:57:ARG:HH11	1:B:57:ARG:CG	1.85	0.88
1:A:335:ILE:HD11	1:A:358:VAL:HG23	0.88	0.87
1:B:335:ILE:HG12	1:B:356:PRO:HB2	1.53	0.87
1:B:335:ILE:CG1	1:B:356:PRO:HB2	2.05	0.87
2:H:169:VAL:HG21	3:L:161:GLU:HB3	1.55	0.86
2:I:51:ILE:HD11	2:I:78:LEU:CD1	2.06	0.86
3:M:49:TYR:CE1	3:M:53:LYS:HB2	2.12	0.84
1:A:34:MET:HG2	1:A:40:THR:CG2	2.08	0.83
1:A:141:ILE:HD11	1:A:183:MET:HE1	1.60	0.83
1:B:34:MET:HG2	1:B:40:THR:CG2	2.07	0.83
2:I:51:ILE:HG22	2:I:57:ARG:HG2	1.62	0.82
1:B:155:THR:HG21	2:H:99:PHE:HZ	1.46	0.80
2:I:169:VAL:HG21	3:M:161:GLU:HB3	1.64	0.79
1:A:220:TRP:CD1	1:A:220:TRP:C	2.57	0.77
1:B:155:THR:HG21	2:H:99:PHE:CZ	2.20	0.77

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:LEU:HD12	1:A:285:CYS:SG	2.24	0.77
1:A:220:TRP:C	1:A:220:TRP:HD1	1.89	0.76
1:B:57:ARG:HG3	1:B:57:ARG:NH1	1.95	0.75
3:M:179:LEU:HD21	3:M:181:LEU:HD21	1.67	0.75
1:A:188:ARG:HG3	1:A:284:LYS:HE3	1.69	0.75
1:A:371:PRO:CG	1:A:391:TRP:CZ3	2.71	0.74
1:B:283:LEU:HD22	1:B:285:CYS:SG	2.27	0.74
3:L:186:TRP:CH2	3:L:209:PRO:HA	2.23	0.73
1:B:133:GLU:HG2	1:B:167:GLN:HB2	1.71	0.73
1:A:391:TRP:CZ3	1:A:1393:ARG:HG3	2.24	0.72
1:A:220:TRP:HD1	1:A:220:TRP:O	1.76	0.69
1:A:166:PRO:HB3	1:A:187:PRO:HG3	1.75	0.68
3:M:179:LEU:CD2	3:M:181:LEU:HG	2.24	0.67
3:M:133:LEU:HD12	3:M:179:LEU:HD22	1.77	0.67
1:A:26:GLU:HG2	1:A:29:SER:HB2	1.77	0.66
3:M:49:TYR:CE1	3:M:53:LYS:CB	2.78	0.66
2:I:51:ILE:HG23	2:I:69:ILE:CG2	2.25	0.66
1:A:306:PHE:O	1:A:387:LEU:HD11	1.96	0.65
1:B:166:PRO:HB3	1:B:187:PRO:HG3	1.78	0.65
3:L:26:THR:HG23	3:L:27:SER:H	1.62	0.65
1:B:164:ILE:HD11	1:B:185:CYS:HB2	1.77	0.65
3:M:49:TYR:CD1	3:M:53:LYS:HB2	2.32	0.65
2:H:83:THR:O	2:H:111:VAL:HG11	1.97	0.64
3:M:26:THR:HG23	3:M:27:SER:H	1.61	0.64
1:A:185:CYS:HB3	1:A:283:LEU:HD11	1.79	0.64
1:A:220:TRP:CD1	1:A:220:TRP:O	2.51	0.64
3:L:4:LEU:HD13	3:L:97:VAL:HG12	1.79	0.64
3:M:179:LEU:HD21	3:M:181:LEU:CD2	2.28	0.64
3:M:4:LEU:HD13	3:M:97:VAL:HG12	1.81	0.63
1:A:391:TRP:CH2	1:A:1393:ARG:HG3	2.34	0.63
2:I:51:ILE:HD11	2:I:78:LEU:HD12	1.80	0.63
1:B:306:PHE:O	1:B:387:LEU:HD11	1.98	0.62
2:I:83:THR:O	2:I:111:VAL:HG11	1.99	0.62
1:A:314:GLU:OE2	1:A:391:TRP:HH2	1.83	0.62
1:B:41:LEU:HD11	1:B:292:LEU:HD11	1.80	0.62
2:I:51:ILE:CG2	2:I:69:ILE:HG23	2.30	0.61
1:A:314:GLU:OE2	1:A:391:TRP:CH2	2.53	0.61
1:A:3:CYS:O	1:A:6:ILE:HG23	2.01	0.61
1:B:26:GLU:HG2	1:B:29:SER:HB2	1.82	0.61
1:A:188:ARG:HG3	1:A:284:LYS:CE	2.31	0.61
1:B:185:CYS:HB3	1:B:283:LEU:HD21	1.82	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:SER:HB2	1:B:255:SER:HB2	1.82	0.60
1:A:220:TRP:HE1	1:A:232:ILE:HD13	1.67	0.60
1:A:41:LEU:HD11	1:A:292:LEU:HD11	1.83	0.59
1:B:315:THR:HG21	1:B:319:THR:OG1	2.03	0.59
1:A:342:LEU:HD11	1:A:375:ASP:HB2	1.84	0.59
1:B:342:LEU:HD11	1:B:375:ASP:HB2	1.84	0.59
2:I:100:TYR:CE1	4:C:2:NAG:H4	2.38	0.58
1:A:371:PRO:HG2	1:A:391:TRP:HZ3	1.60	0.58
1:B:3:CYS:O	1:B:6:ILE:HG23	2.02	0.58
1:A:315:THR:HG21	1:A:319:THR:OG1	2.02	0.58
1:B:20:TRP:CE2	1:B:286:ARG:HD3	2.38	0.57
1:B:315:THR:CG2	1:B:319:THR:OG1	2.52	0.57
1:B:57:ARG:CG	1:B:57:ARG:NH1	2.55	0.57
1:B:72:SER:O	2:I:100(I):ASP:HA	2.05	0.57
2:I:51:ILE:HG23	2:I:69:ILE:HG23	1.87	0.57
1:A:315:THR:CG2	1:A:319:THR:OG1	2.52	0.57
1:A:335:ILE:CD1	1:A:358:VAL:CG2	2.53	0.57
2:H:93:VAL:HG11	2:H:100(P):MET:HB3	1.86	0.57
2:I:93:VAL:HG11	2:I:100(P):MET:HB3	1.86	0.56
1:A:335:ILE:N	1:A:335:ILE:HD13	2.21	0.56
1:B:289:MET:HG3	1:B:292:LEU:HD12	1.87	0.56
3:M:179:LEU:HD23	3:M:181:LEU:HG	1.86	0.56
1:A:391:TRP:CZ3	1:A:1393:ARG:HB2	2.40	0.56
1:A:185:CYS:HB3	1:A:283:LEU:CD1	2.35	0.56
1:A:251:VAL:HG21	1:B:204:LYS:HE2	1.88	0.56
1:B:185:CYS:HB3	1:B:283:LEU:CD2	2.35	0.56
1:A:89:ARG:HG2	1:A:229:SER:HB3	1.89	0.55
2:I:195:ILE:HD12	2:I:197:ASN:HD21	1.71	0.55
1:B:34:MET:CG	1:B:40:THR:HG22	2.32	0.55
2:I:51:ILE:HG23	2:I:69:ILE:HG21	1.88	0.55
1:A:289:MET:HG3	1:A:292:LEU:HD12	1.87	0.55
2:H:195:ILE:HD12	2:H:197:ASN:HD21	1.72	0.54
3:M:49:TYR:HE1	3:M:53:LYS:CB	2.20	0.54
1:B:315:THR:HG23	1:B:317:HIS:H	1.72	0.54
1:A:34:MET:CG	1:A:40:THR:HG22	2.34	0.54
1:A:315:THR:HG23	1:A:317:HIS:H	1.73	0.54
1:B:133:GLU:HG2	1:B:167:GLN:CB	2.38	0.54
2:I:164:HIS:HE1	3:M:174:ALA:HB3	1.72	0.53
1:A:220:TRP:CD1	1:A:232:ILE:HB	2.44	0.53
2:I:51:ILE:HG12	2:I:69:ILE:CD1	2.38	0.53
3:M:179:LEU:HD23	3:M:179:LEU:C	2.28	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:9:SER:HB3	3:M:144:ALA:CB	2.39	0.53
2:I:59:TYR:H	3:M:95:ARG:HH21	1.56	0.53
3:L:9:SER:HB3	3:L:144:ALA:CB	2.38	0.53
3:M:179:LEU:HD21	3:M:181:LEU:CG	2.39	0.52
3:M:179:LEU:CD2	3:M:181:LEU:CG	2.88	0.52
1:B:20:TRP:CZ2	1:B:286:ARG:HD3	2.45	0.52
2:H:59:TYR:H	3:L:95:ARG:HH21	1.57	0.52
1:B:72:SER:OG	2:I:100(H):PRO:HD2	2.10	0.52
1:B:371:PRO:HG2	1:B:391:TRP:HE1	1.75	0.51
3:M:49:TYR:C	3:M:49:TYR:HD1	2.14	0.51
1:A:137:TYR:HB2	1:A:164:ILE:HG13	1.93	0.51
1:A:220:TRP:NE1	1:A:232:ILE:HB	2.26	0.51
1:B:89:ARG:HG2	1:B:229:SER:HB3	1.93	0.51
2:H:164:HIS:CE1	3:L:174:ALA:HB3	2.46	0.51
2:H:155:ASN:OD1	2:H:195:ILE:HG13	2.10	0.51
1:B:133:GLU:HG2	1:B:167:GLN:CG	2.42	0.50
1:A:288:ARG:HB3	1:A:290:ASP:OD1	2.12	0.49
1:B:125:MET:HG2	1:B:199:LEU:HD11	1.94	0.49
3:M:49:TYR:CD1	3:M:49:TYR:C	2.85	0.49
1:B:371:PRO:CG	1:B:391:TRP:HE1	2.25	0.49
2:H:97:VAL:CG1	2:H:100(O):GLY:HA3	2.43	0.49
1:A:125:MET:HG2	1:A:199:LEU:HD11	1.93	0.49
2:H:192:GLN:HG2	2:H:194:TYR:CZ	2.48	0.48
1:B:380:VAL:HG13	1:B:387:LEU:HB2	1.96	0.48
1:A:222:PRO:HD2	1:A:225:ASP:HB3	1.96	0.48
2:H:205:THR:HG21	2:I:203:SER:HB2	1.96	0.48
2:I:88:ALA:HB3	2:I:90:TYR:CE1	2.48	0.48
2:I:155:ASN:OD1	2:I:195:ILE:HG13	2.13	0.48
1:A:318:GLY:HA3	1:A:1393:ARG:HH21	1.79	0.47
1:B:222:PRO:HD2	1:B:225:ASP:HB3	1.96	0.47
2:I:97:VAL:CG1	2:I:100(O):GLY:HA3	2.44	0.47
1:A:34:MET:HG2	1:A:40:THR:HG21	1.92	0.47
2:I:51:ILE:HD12	2:I:71:ARG:HD2	1.94	0.47
2:I:88:ALA:HB3	2:I:90:TYR:HE1	1.79	0.47
1:A:380:VAL:HG13	1:A:387:LEU:HB2	1.95	0.47
2:I:144:ASP:HB3	2:I:175:LEU:HD23	1.97	0.47
2:H:100:TYR:CE1	4:E:2:NAG:H4	2.50	0.47
1:A:24:VAL:HG22	1:A:284:LYS:HG2	1.97	0.46
2:H:97:VAL:HG13	2:H:100(O):GLY:HA3	1.97	0.46
1:B:308:ILE:H	1:B:308:ILE:HG13	1.52	0.46
1:B:318:GLY:HA3	1:B:1393:ARG:HH21	1.80	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:TRP:CZ3	1:A:1393:ARG:CG	2.98	0.46
1:B:99:ARG:NH2	2:I:100(I):ASP:OD2	2.49	0.46
1:B:102:GLY:HA2	4:C:1:NAG:H81	1.98	0.46
1:B:164:ILE:HD11	1:B:169:SER:HA	1.98	0.46
2:I:164:HIS:CE1	3:M:174:ALA:HB3	2.49	0.46
1:B:34:MET:HG2	1:B:40:THR:HG21	1.95	0.46
3:M:162:THR:HG22	3:M:177:SER:OG	2.16	0.46
2:H:164:HIS:HE1	3:L:174:ALA:HB3	1.81	0.45
3:L:8:VAL:HG21	3:L:146:THR:OG1	2.15	0.45
1:B:164:ILE:CD1	1:B:169:SER:HA	2.47	0.45
1:A:141:ILE:HD11	1:A:183:MET:CE	2.39	0.45
1:B:164:ILE:CD1	1:B:185:CYS:HB2	2.46	0.45
1:B:340:THR:HG22	1:B:347:VAL:HA	1.98	0.45
2:I:82:MET:HE3	2:I:90:TYR:CE2	2.52	0.45
2:I:97:VAL:HG13	2:I:100(O):GLY:HA3	1.99	0.45
1:B:342:LEU:HD11	1:B:375:ASP:CB	2.46	0.45
1:B:196:MET:HB3	1:B:207:LEU:HG	1.99	0.45
1:A:65:LEU:HD12	1:A:252:VAL:HG22	1.99	0.44
2:H:205:THR:CG2	2:I:203:SER:HB2	2.47	0.44
2:H:97:VAL:HG13	2:H:100(O):GLY:CA	2.48	0.44
1:B:65:LEU:HD12	1:B:252:VAL:HG22	2.00	0.44
2:I:82:MET:CE	2:I:90:TYR:CE2	3.00	0.44
1:A:247:LYS:HB2	2:H:100(F):TYR:CZ	2.52	0.44
2:I:51:ILE:CG2	2:I:57:ARG:HG2	2.42	0.44
1:A:340:THR:HG22	1:A:347:VAL:HA	2.00	0.44
3:L:162:THR:HG22	3:L:177:SER:OG	2.18	0.43
1:A:24:VAL:HG13	1:A:282:HIS:HB3	2.00	0.43
1:A:342:LEU:HD11	1:A:375:ASP:CB	2.46	0.43
2:I:97:VAL:HG13	2:I:100(O):GLY:CA	2.48	0.43
1:A:345:ARG:HD2	1:A:346:HIS:NE2	2.34	0.43
1:B:133:GLU:OE1	1:B:167:GLN:HB2	2.18	0.43
1:B:345:ARG:HD2	1:B:346:HIS:NE2	2.34	0.43
1:A:308:ILE:H	1:A:308:ILE:HG13	1.48	0.43
1:B:164:ILE:HG13	1:B:165:THR:N	2.33	0.43
2:I:51:ILE:HD11	2:I:78:LEU:HD11	1.95	0.43
3:M:133:LEU:HD21	3:M:186:TRP:CZ3	2.53	0.43
1:A:33:THR:HG21	1:A:43:PHE:HE1	1.84	0.43
1:B:33:THR:HG22	1:B:41:LEU:HB2	2.01	0.43
2:H:184:VAL:HG21	2:H:194:TYR:CZ	2.54	0.43
2:I:93:VAL:CG1	2:I:100(P):MET:HB3	2.48	0.42
1:B:329:ASP:HA	1:B:361:LYS:HE2	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:VAL:HG22	1:A:368:GLU:HG3	2.01	0.42
1:B:33:THR:HG21	1:B:43:PHE:HE1	1.85	0.42
2:H:93:VAL:CG1	2:H:100(P):MET:HB3	2.48	0.42
1:B:371:PRO:CG	1:B:391:TRP:NE1	2.83	0.42
3:L:56:SER:OG	4:E:3:BMA:H3	2.20	0.42
1:A:329:ASP:HA	1:A:361:LYS:HE2	2.02	0.42
1:B:133:GLU:HG2	1:B:167:GLN:HG2	2.01	0.42
1:A:33:THR:HG22	1:A:41:LEU:HB2	2.03	0.41
1:B:321:VAL:HG22	1:B:368:GLU:HG3	2.01	0.41
1:B:133:GLU:CG	1:B:167:GLN:HB2	2.46	0.41
1:A:58:LYS:NZ	1:A:226:THR:HG23	2.36	0.41
1:B:156:GLY:C	1:B:158:HIS:H	2.24	0.41
2:H:169:VAL:CG2	3:L:161:GLU:HB3	2.39	0.41
2:I:51:ILE:HG12	2:I:69:ILE:HD13	2.03	0.41
2:H:100(G):TYR:HA	2:H:100(H):PRO:HD3	1.99	0.41
1:A:156:GLY:C	1:A:158:HIS:H	2.23	0.40
2:I:82:MET:HE3	2:I:90:TYR:HE2	1.84	0.40
1:A:141:ILE:CD1	1:A:183:MET:HE1	2.42	0.40
1:B:155:THR:CG2	2:H:99:PHE:CZ	3.00	0.40
2:H:184:VAL:HG21	2:H:194:TYR:CE1	2.56	0.40
1:B:132:PRO:HG3	1:B:193:PHE:HB2	2.04	0.40
1:A:132:PRO:HG3	1:A:193:PHE:HB2	2.03	0.40
1:A:145:SER:C	1:A:147:GLU:H	2.25	0.40
2:H:144:ASP:HB3	2:H:175:LEU:HD13	2.03	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	387/422 (92%)	370 (96%)	16 (4%)	1 (0%)	41 74

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	384/422 (91%)	367 (96%)	16 (4%)	1 (0%)	41	74
2	H	221/283 (78%)	211 (96%)	10 (4%)	0	100	100
2	I	193/283 (68%)	185 (96%)	8 (4%)	0	100	100
3	L	211/218 (97%)	205 (97%)	6 (3%)	0	100	100
3	M	208/218 (95%)	203 (98%)	5 (2%)	0	100	100
All	All	1604/1846 (87%)	1541 (96%)	61 (4%)	2 (0%)	51	83

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	147	GLU
1	B	147	GLU

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	341/366 (93%)	295 (86%)	46 (14%)	4	22
1	B	340/366 (93%)	295 (87%)	45 (13%)	4	22
2	H	190/236 (80%)	178 (94%)	12 (6%)	18	47
2	I	174/236 (74%)	164 (94%)	10 (6%)	20	50
3	L	180/185 (97%)	171 (95%)	9 (5%)	24	53
3	M	178/185 (96%)	167 (94%)	11 (6%)	18	48
All	All	1403/1574 (89%)	1270 (90%)	133 (10%)	8	33

All (133) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	6	ILE
1	A	8	ASN
1	A	21	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	30	CYS
1	A	61	ILE
1	A	67	ASN
1	A	70	THR
1	A	103	ASN
1	A	107	LEU
1	A	120	THR
1	A	128	LYS
1	A	129	ILE
1	A	147	GLU
1	A	164	ILE
1	A	169	SER
1	A	176	THR
1	A	182	THR
1	A	188	ARG
1	A	191	LEU
1	A	200	GLN
1	A	211	GLN
1	A	220	TRP
1	A	227	GLN
1	A	235	GLU
1	A	237	LEU
1	A	252	VAL
1	A	253	LEU
1	A	271	GLN
1	A	276	ASN
1	A	277	LEU
1	A	283	LEU
1	A	287	LEU
1	A	289	MET
1	A	290	ASP
1	A	298	SER
1	A	308	ILE
1	A	329	ASP
1	A	335	ILE
1	A	338	GLU
1	A	343	GLU
1	A	346	HIS
1	A	350	ARG
1	A	351	LEU
1	A	389	LEU
1	A	1392	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	2	ARG
1	B	6	ILE
1	B	8	ASN
1	B	20	TRP
1	B	21	VAL
1	B	22	ASP
1	B	30	CYS
1	B	57	ARG
1	B	61	ILE
1	B	70	THR
1	B	88	LYS
1	B	103	ASN
1	B	107	LEU
1	B	120	THR
1	B	128	LYS
1	B	129	ILE
1	B	133	GLU
1	B	147	GLU
1	B	164	ILE
1	B	169	SER
1	B	176	THR
1	B	182	THR
1	B	188	ARG
1	B	191	LEU
1	B	200	GLN
1	B	211	GLN
1	B	237	LEU
1	B	252	VAL
1	B	253	LEU
1	B	255	SER
1	B	276	ASN
1	B	277	LEU
1	B	287	LEU
1	B	289	MET
1	B	290	ASP
1	B	308	ILE
1	B	329	ASP
1	B	335	ILE
1	B	338	GLU
1	B	343	GLU
1	B	346	HIS
1	B	350	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	351	LEU
1	B	389	LEU
1	B	1392	LEU
2	H	13	ARG
2	H	25	SER
2	H	53	ASP
2	H	63	VAL
2	H	105	GLN
2	H	111	VAL
2	H	142	VAL
2	H	143	LYS
2	H	184	VAL
2	H	188	SER
2	H	195	ILE
2	H	210	ARG
2	I	13	ARG
2	I	25	SER
2	I	53	ASP
2	I	63	VAL
2	I	105	GLN
2	I	111	VAL
2	I	142	VAL
2	I	143	LYS
2	I	195	ILE
2	I	210	ARG
3	L	5	THR
3	L	19	ILE
3	L	66	LYS
3	L	82	ASP
3	L	93	THR
3	L	94	SER
3	L	97	VAL
3	L	111	LYS
3	L	136	LEU
3	M	5	THR
3	M	19	ILE
3	M	49	TYR
3	M	66	LYS
3	M	93	THR
3	M	94	SER
3	M	97	VAL
3	M	111	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	M	136	LEU
3	M	171	ASN
3	M	197	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	233	GLN
1	A	276	ASN
1	B	8	ASN
1	B	276	ASN
2	H	39	GLN
2	H	105	GLN
2	H	164	HIS
2	I	39	GLN
2	I	105	GLN
2	I	164	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

14 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	C	1	4,1	14,14,15	0.30	0	17,19,21	0.52	0
4	NAG	C	2	4	14,14,15	0.30	0	17,19,21	1.67	3 (17%)
4	BMA	C	3	4	11,11,12	0.38	0	15,15,17	1.15	1 (6%)
4	MAN	C	4	4	11,11,12	0.41	0	15,15,17	1.12	1 (6%)
4	MAN	C	5	4	11,11,12	0.30	0	15,15,17	0.75	1 (6%)
4	FUC	C	6	4	10,10,11	0.48	0	14,14,16	1.13	1 (7%)
5	NAG	D	1	5,1	14,14,15	0.26	0	17,19,21	1.06	2 (11%)
5	NAG	D	2	5	14,14,15	0.35	0	17,19,21	0.70	0
4	NAG	E	1	4,1	14,14,15	0.30	0	17,19,21	0.58	0
4	NAG	E	2	4	14,14,15	0.28	0	17,19,21	1.70	3 (17%)
4	BMA	E	3	4	11,11,12	0.42	0	15,15,17	1.21	1 (6%)
4	MAN	E	4	4	11,11,12	0.40	0	15,15,17	1.06	1 (6%)
4	MAN	E	5	4	11,11,12	0.28	0	15,15,17	0.84	1 (6%)
4	FUC	E	6	4	10,10,11	0.47	0	14,14,16	1.13	1 (7%)

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	C	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	C	2	4	-	0/6/23/26	0/1/1/1
4	BMA	C	3	4	-	2/2/19/22	1/1/1/1
4	MAN	C	4	4	-	1/2/19/22	1/1/1/1
4	MAN	C	5	4	-	0/2/19/22	0/1/1/1
4	FUC	C	6	4	-	-	0/1/1/1
5	NAG	D	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	D	2	5	-	0/6/23/26	0/1/1/1
4	NAG	E	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	2/2/19/22	1/1/1/1
4	MAN	E	4	4	-	1/2/19/22	1/1/1/1
4	MAN	E	5	4	-	0/2/19/22	0/1/1/1
4	FUC	E	6	4	-	-	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	C	2	NAG	C1-O5-C5	4.63	118.46	112.19
4	E	2	NAG	C1-O5-C5	4.31	118.03	112.19
4	E	2	NAG	O5-C1-C2	4.12	117.80	111.29
4	C	4	MAN	C1-O5-C5	3.95	117.54	112.19
4	E	4	MAN	C1-O5-C5	3.83	117.38	112.19
4	C	2	NAG	O5-C1-C2	3.69	117.12	111.29
4	C	3	BMA	C1-O5-C5	3.52	116.97	112.19
4	E	3	BMA	C1-O5-C5	3.44	116.85	112.19
4	C	6	FUC	C1-O5-C5	2.99	119.56	112.78
4	E	6	FUC	C1-O5-C5	2.87	119.29	112.78
4	E	2	NAG	O4-C4-C5	2.65	115.88	109.30
4	E	5	MAN	C1-O5-C5	2.65	115.78	112.19
5	D	1	NAG	O5-C1-C2	2.64	115.46	111.29
4	C	2	NAG	O4-C4-C5	2.54	115.61	109.30
5	D	1	NAG	C1-O5-C5	2.43	115.48	112.19
4	C	5	MAN	C1-O5-C5	2.25	115.25	112.19

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	1	NAG	O5-C5-C6-O6
4	C	3	BMA	O5-C5-C6-O6
4	C	3	BMA	C4-C5-C6-O6
4	E	3	BMA	C4-C5-C6-O6
4	E	3	BMA	O5-C5-C6-O6
4	E	1	NAG	O5-C5-C6-O6
4	E	1	NAG	C4-C5-C6-O6
4	C	1	NAG	C4-C5-C6-O6
4	C	4	MAN	O5-C5-C6-O6
4	E	4	MAN	O5-C5-C6-O6
5	D	1	NAG	O5-C5-C6-O6

All (4) ring outliers are listed below:

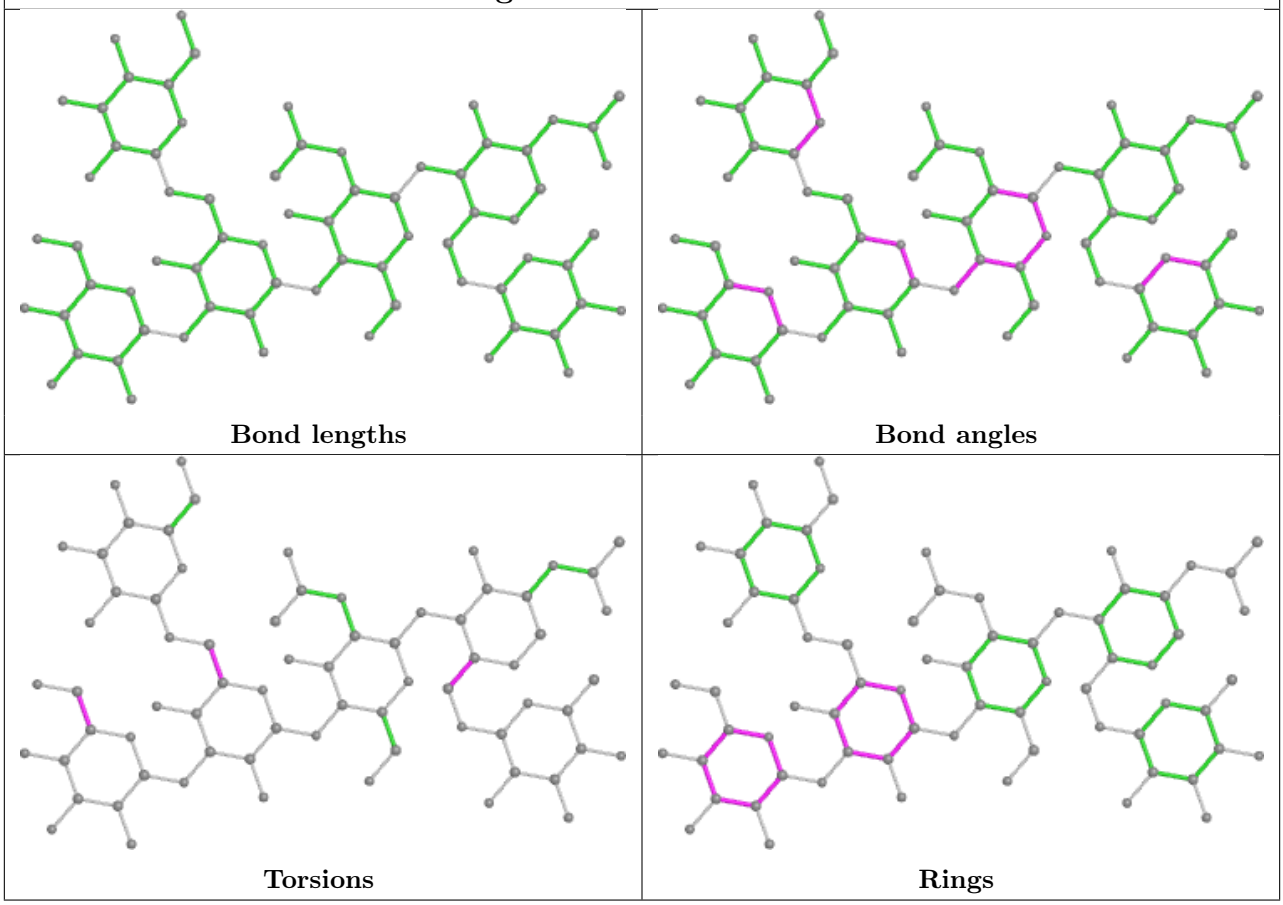
Mol	Chain	Res	Type	Atoms
4	C	3	BMA	C1-C2-C3-C4-C5-O5
4	E	3	BMA	C1-C2-C3-C4-C5-O5
4	E	4	MAN	C1-C2-C3-C4-C5-O5
4	C	4	MAN	C1-C2-C3-C4-C5-O5

4 monomers are involved in 4 short contacts:

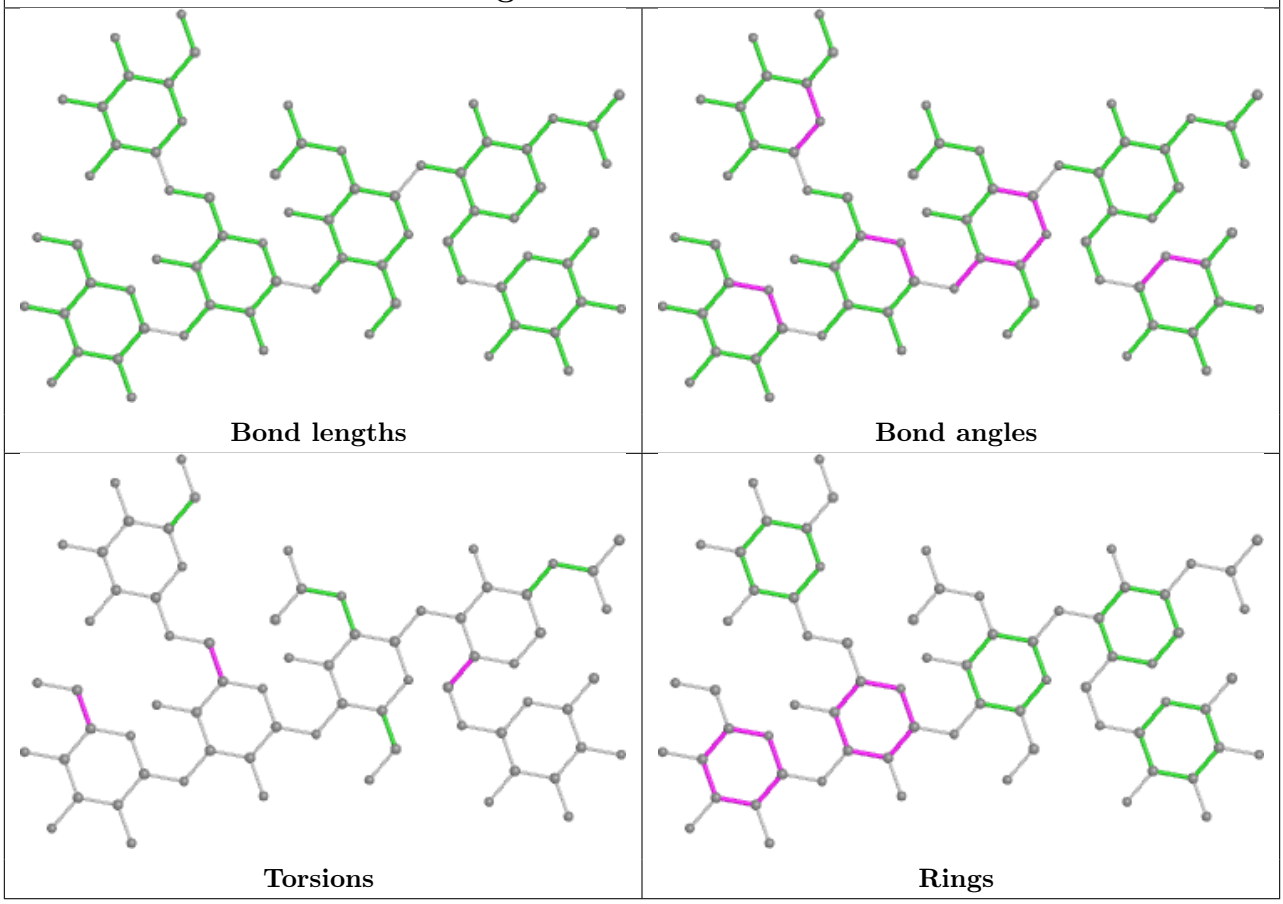
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	2	NAG	1	0
4	C	2	NAG	1	0
4	E	3	BMA	1	0
4	C	1	NAG	1	0

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

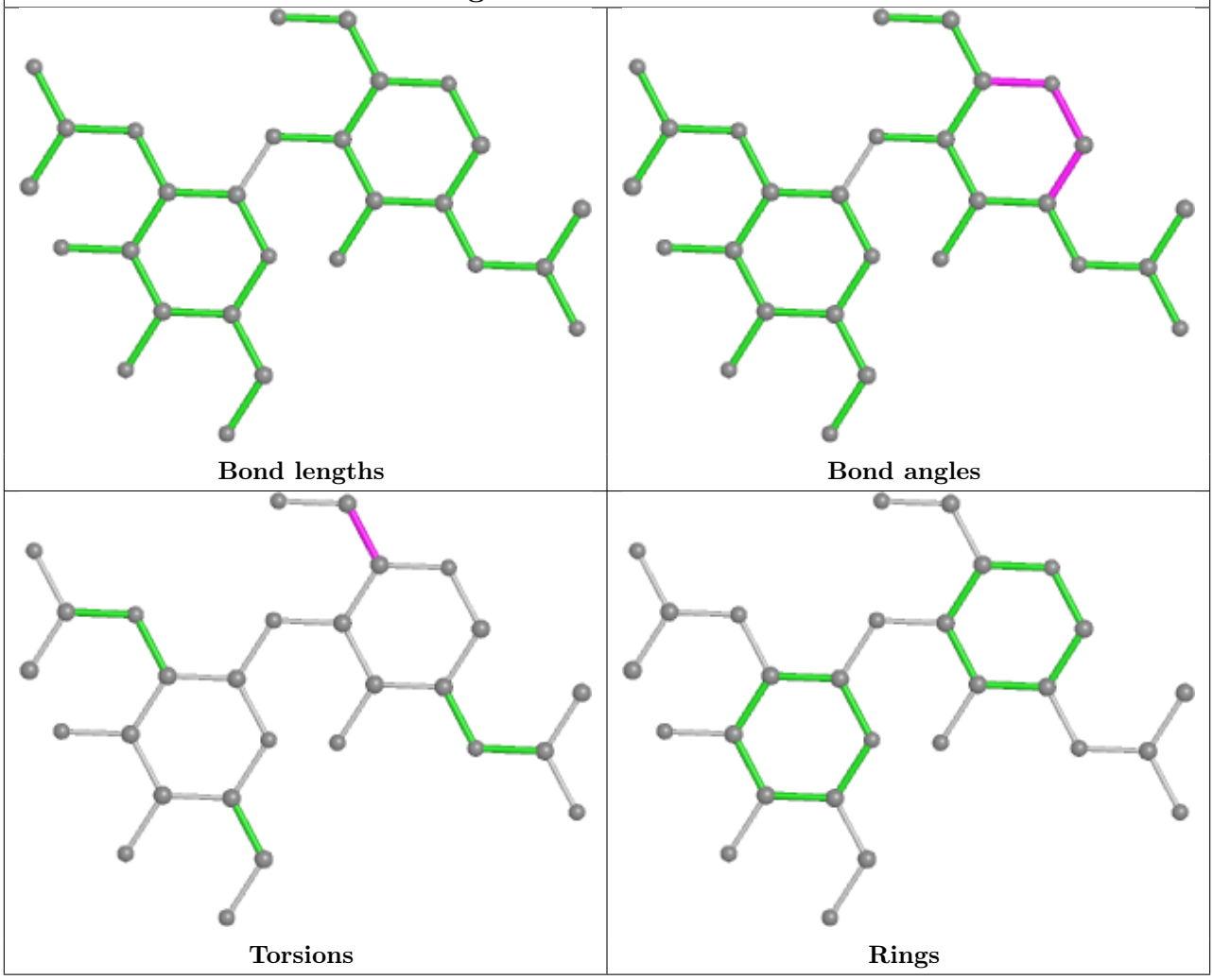
Oligosaccharide Chain C



Oligosaccharide Chain E



Oligosaccharide Chain D



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	$\#\|Z\| > 2$
7	SO4	H	581	-	4,4,4	0.22	0	6,6,6	0.16	0
7	SO4	I	581	-	4,4,4	0.23	0	6,6,6	0.14	0
6	NAG	B	567	1	14,14,15	0.38	0	17,19,21	0.66	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
6	NAG	B	567	1	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	567	NAG	C1-C2-N2-C7

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	391/422 (92%)	0.28	17 (4%)	35	29	100, 125, 158, 180	0
1	B	390/422 (92%)	0.19	9 (2%)	60	51	94, 125, 158, 177	0
2	H	225/283 (79%)	0.16	3 (1%)	77	69	94, 126, 146, 162	0
2	I	203/283 (71%)	0.48	21 (10%)	6	6	91, 122, 187, 206	0
3	L	213/218 (97%)	0.25	7 (3%)	46	37	90, 126, 160, 184	0
3	M	210/218 (96%)	0.38	15 (7%)	16	12	92, 128, 178, 193	0
All	All	1632/1846 (88%)	0.28	72 (4%)	34	28	90, 125, 172, 206	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	213	PRO	6.7
1	A	191	LEU	4.9
2	I	212	GLU	4.4
1	A	346	HIS	4.4
1	A	347	VAL	4.3
3	L	210	THR	3.9
1	A	348	LEU	3.9
2	H	112	SER	3.8
3	M	153	SER	3.6
2	I	206	LYS	3.5
2	I	208	ASP	3.5
3	M	184	GLU	3.5
1	B	96	MET	3.4
1	A	190	GLY	3.4
3	L	153	SER	3.2
3	M	151	ALA	3.2
1	B	16	SER	3.2
1	A	342	LEU	3.1
2	I	163	VAL	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
3	M	2	SER	3.0
1	A	343	GLU	2.9
3	M	157	LYS	2.9
1	A	14	GLY	2.8
2	I	211	VAL	2.8
3	M	182	THR	2.8
1	B	190	GLY	2.8
1	A	341	ASP	2.8
2	I	197	ASN	2.7
3	M	190	ARG	2.7
3	M	183	PRO	2.6
1	A	227	GLN	2.6
2	I	168	ALA	2.6
3	L	1	GLN	2.6
2	I	136	ALA	2.5
1	B	346	HIS	2.5
2	I	154	TRP	2.5
1	A	184	GLU	2.4
2	I	149	PRO	2.4
1	B	345	ARG	2.4
2	I	137	ALA	2.4
2	I	173	SER	2.3
2	I	10	GLY	2.3
1	A	331	SER	2.3
3	L	191	SER	2.3
3	L	156	VAL	2.3
3	M	191	SER	2.3
3	M	195	GLN	2.2
1	B	284	LYS	2.2
1	A	345	ARG	2.2
1	B	248	GLN	2.2
2	I	180	SER	2.2
2	I	153	SER	2.2
2	I	199	ASN	2.2
2	H	210	ARG	2.2
1	A	96	MET	2.1
3	L	190	ARG	2.1
1	A	188	ARG	2.1
3	M	150	LYS	2.1
1	A	325	GLN	2.1
3	M	185	GLN	2.1
1	B	343	GLU	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	I	49	SER	2.1
3	M	192	TYR	2.1
2	H	9	GLY	2.1
1	B	348	LEU	2.1
2	I	126	PRO	2.1
3	M	154	SER	2.1
2	I	155	ASN	2.1
2	I	11	LEU	2.0
3	L	27	SER	2.0
3	M	199	GLU	2.0
1	A	340	THR	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)

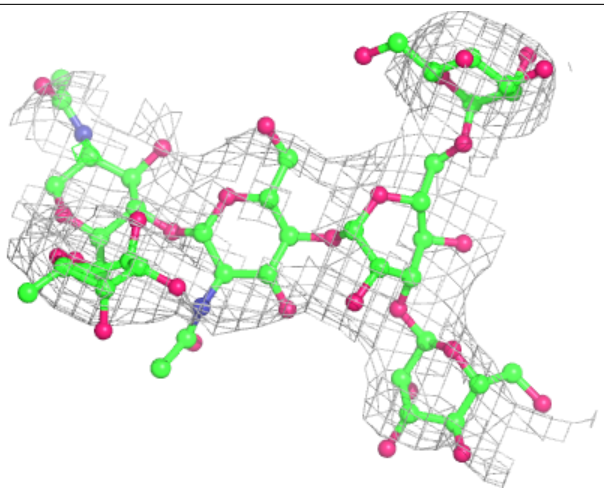
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

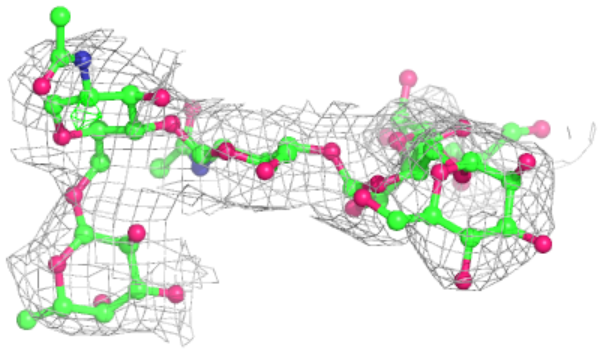
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	NAG	D	2	14/15	0.20	0.76	169,174,177,177	0
5	NAG	D	1	14/15	0.47	0.62	169,172,173,174	0
4	MAN	C	4	11/12	0.76	0.50	151,155,160,161	0
4	FUC	C	6	10/11	0.77	0.43	130,133,135,137	0
4	MAN	E	4	11/12	0.83	0.43	157,161,163,163	0
4	NAG	C	1	14/15	0.85	0.28	110,112,124,128	0
4	FUC	E	6	10/11	0.88	0.32	129,133,135,137	0
4	BMA	E	3	11/12	0.89	0.18	132,133,143,151	0
4	NAG	C	2	14/15	0.89	0.30	117,121,130,131	0
4	MAN	E	5	11/12	0.89	0.27	133,136,137,138	0
4	MAN	C	5	11/12	0.90	0.32	125,127,130,131	0
4	NAG	E	2	14/15	0.90	0.24	112,115,118,125	0
4	NAG	E	1	14/15	0.92	0.23	102,108,118,124	0
4	BMA	C	3	11/12	0.93	0.15	125,127,135,144	0

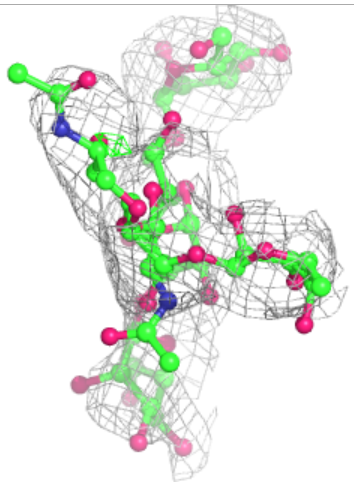
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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

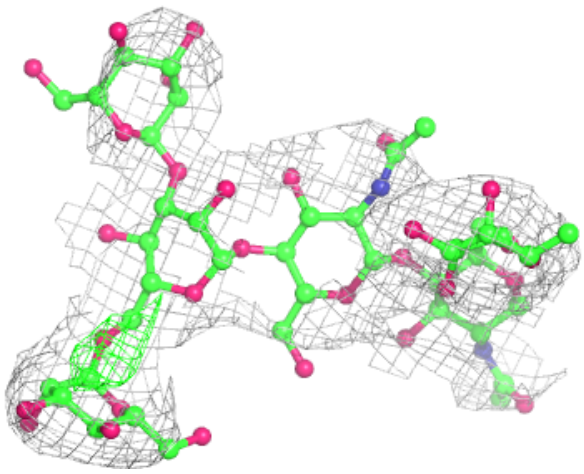


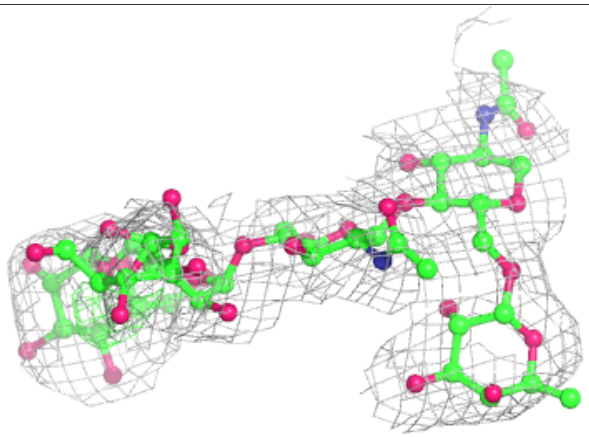


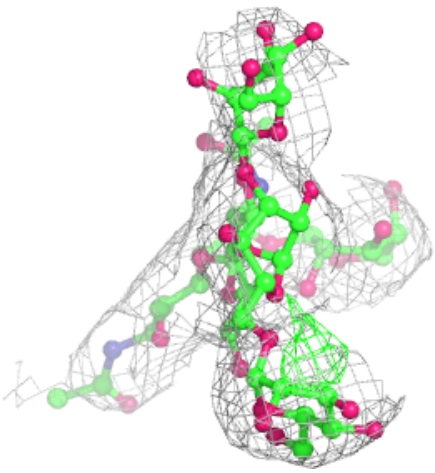


Electron density around Chain E:

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

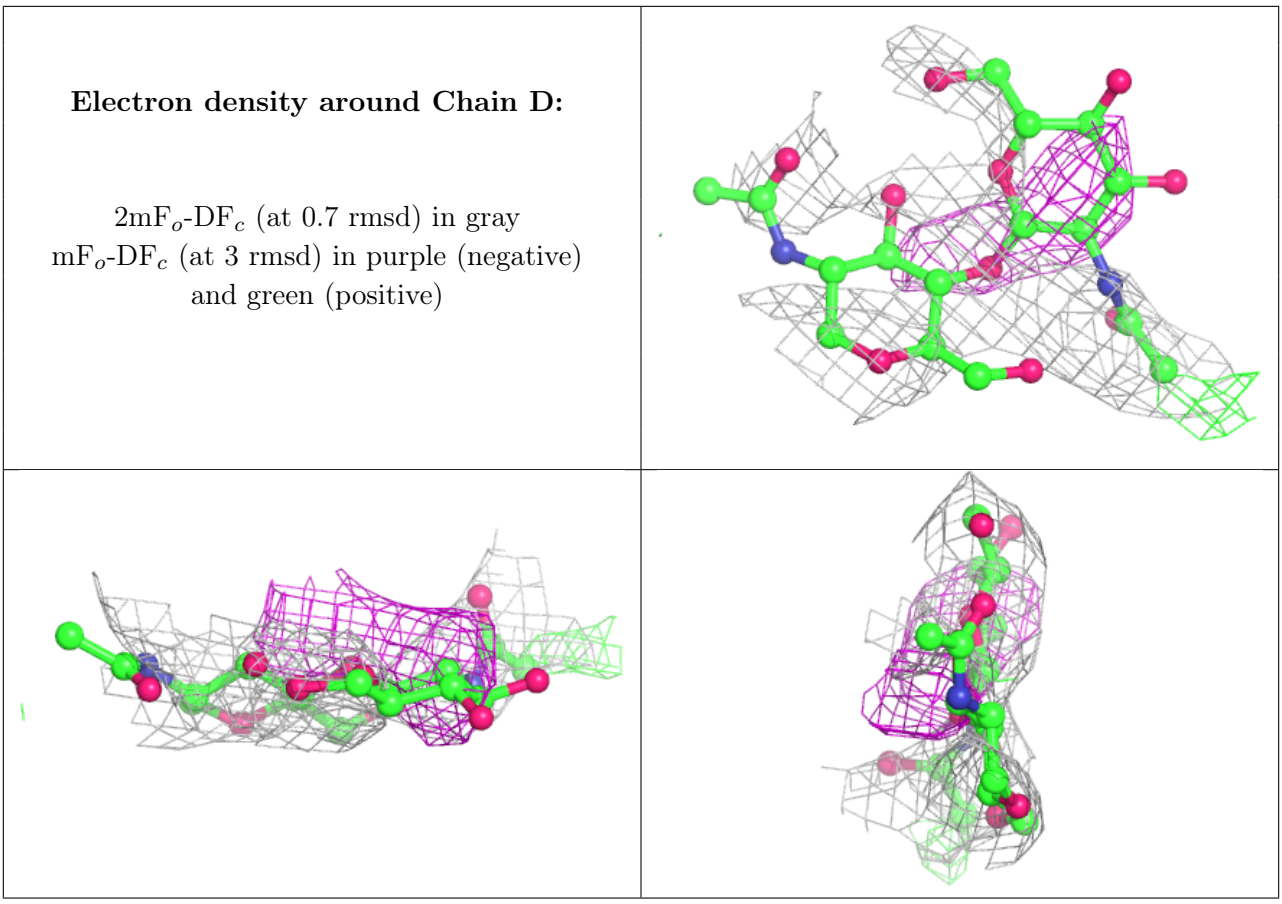






Electron density around Chain D:

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



6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
7	SO4	I	581	5/5	0.59	0.26	169,169,170,170	0
7	SO4	H	581	5/5	0.72	0.24	144,144,145,146	0
6	NAG	B	567	14/15	0.74	0.26	154,156,158,158	0

6.5 Other polymers (i)

There are no such residues in this entry.