



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:04 PM 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

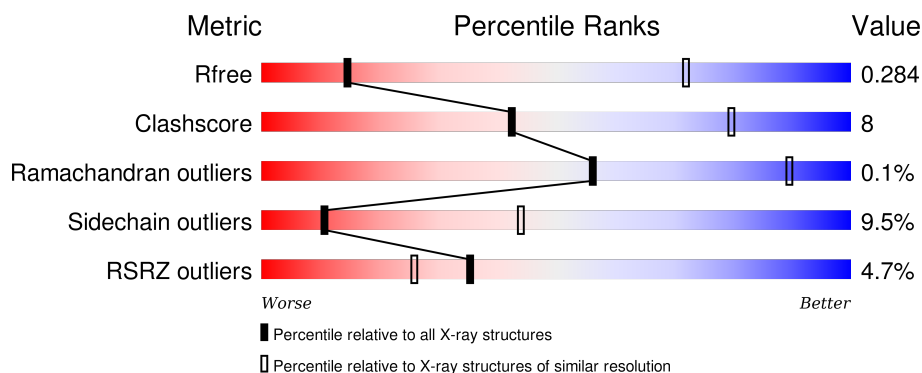
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.85 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}	91344	1000 (4.20-3.52)
Clashscore	102246	1090 (4.20-3.52)
Ramachandran outliers	100387	1046 (4.20-3.52)
Sidechain outliers	100360	1038 (4.20-3.52)
RSRZ outliers	91569	1004 (4.20-3.52)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	422	<div> <div>5%</div> <div>68%</div> <div>21%</div> <div>7%</div> </div>
1	B	422	<div> <div>2%</div> <div>69%</div> <div>20%</div> <div>8%</div> </div>
2	H	283	<div> <div>0%</div> <div>68%</div> <div>10%</div> <div>20%</div> </div>
2	I	283	<div> <div>7%</div> <div>60%</div> <div>11%</div> <div>28%</div> </div>
3	L	218	<div> <div>4%</div> <div>87%</div> <div>11%</div> <div>0%</div> </div>

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Mol	Chain	Length	Quality of chain
3	M	218	

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	A	506	-	-	-	X
4	MAN	B	506	-	-	-	X
5	NAG	A	567	-	-	-	X

2 Entry composition

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
1	A	391	Total	C	N	O	S	0	0	0
			3045	1919	524	578	24			
1	B	390	Total	C	N	O	S	0	0	0
			3036	1914	522	576	24			

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

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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
2	H	225	Total	C	N	O	S	0	0	0
			1718	1088	291	332	7			
2	I	203	Total	C	N	O	S	0	0	0
			1573	999	267	300	7			

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

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

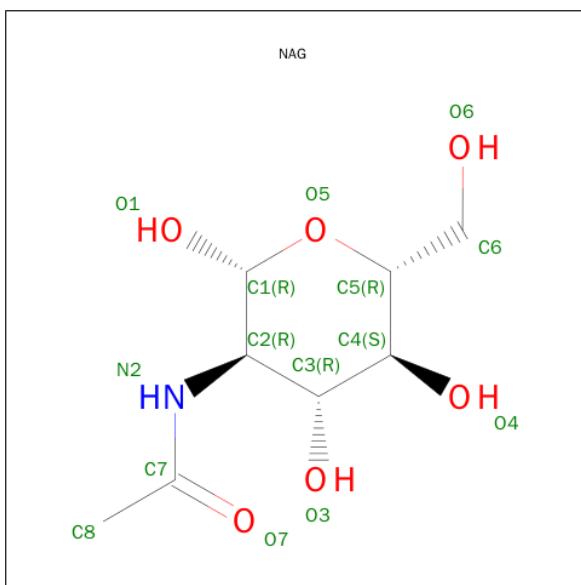
- Molecule 4 is a polymer of unknown type called SUGAR (6-MER).

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

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

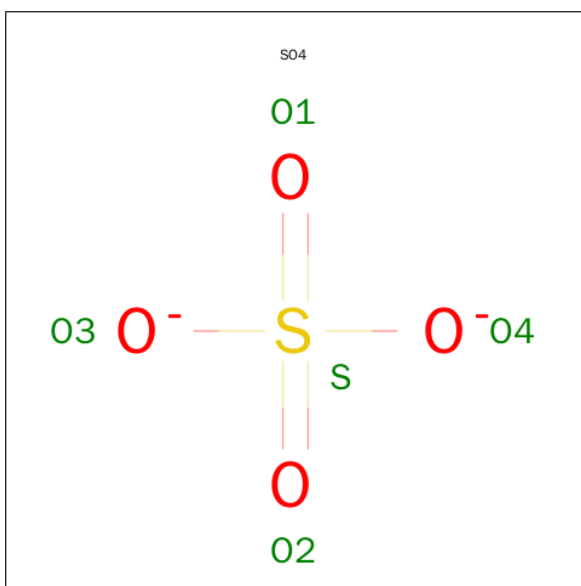
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 7 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

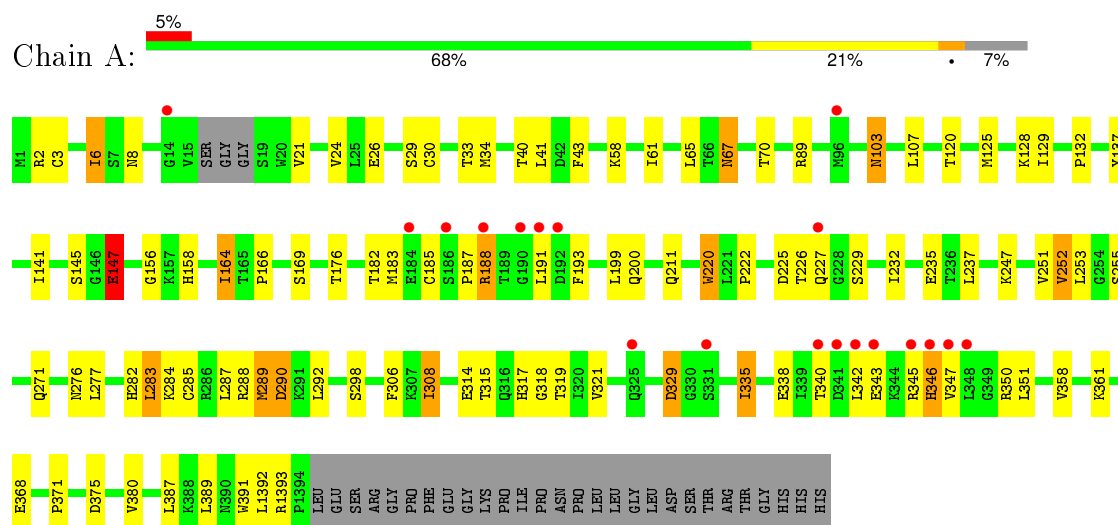
- Molecule 8 is water.

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

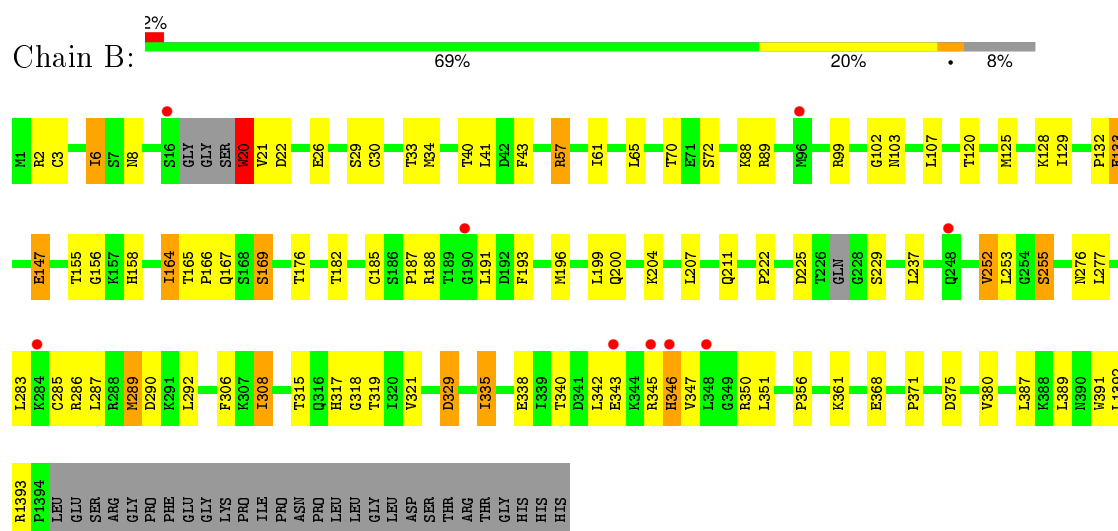
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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

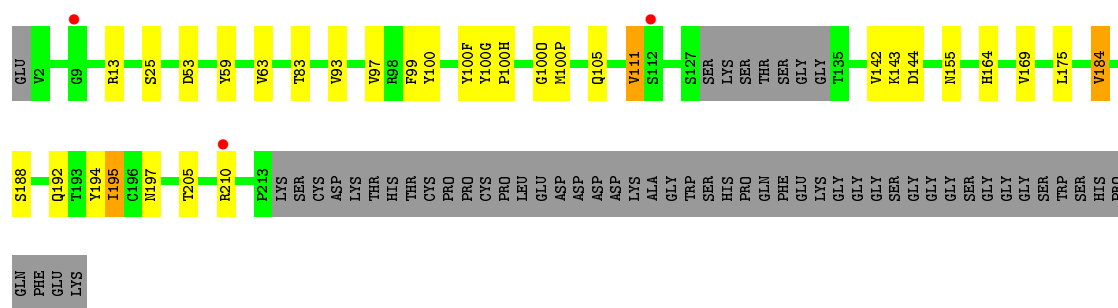


• Molecule 1: ENVELOPE GLYCOPROTEIN E

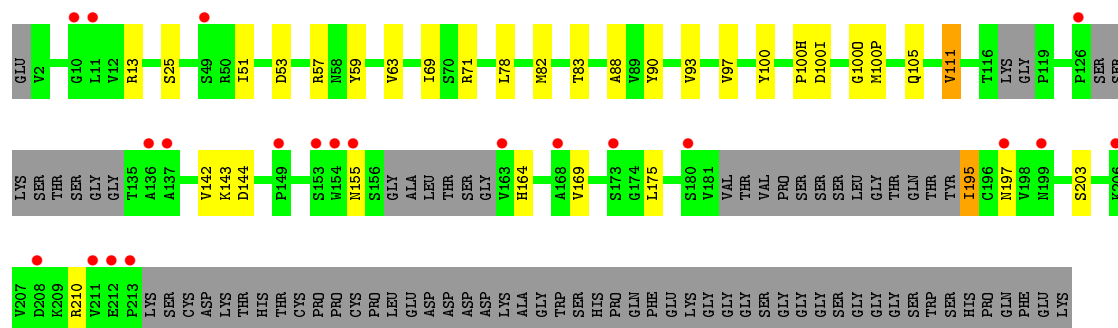


• Molecule 2: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11

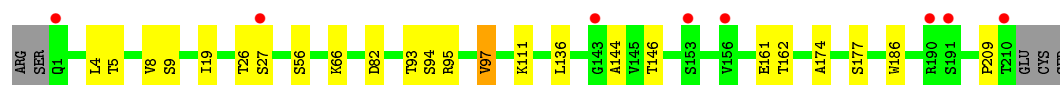
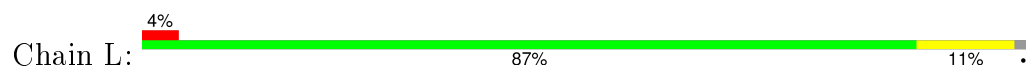




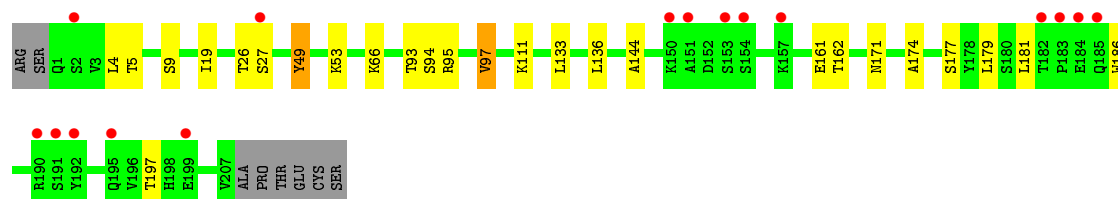
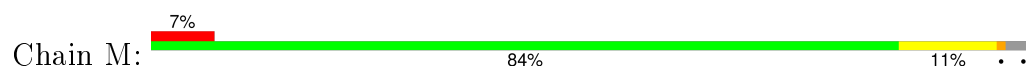
• Molecule 2: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11



• Molecule 3: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11



• Molecule 3: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11



4 Data and refinement statistics

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
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.86Å)	Xtriage
Refinement program	BUSTER 2.11.4	Depositor
R, R_{free}	0.230 , 0.257 0.268 , 0.284	Depositor DCC
R_{free} test set	1072 reflections (5.08%)	DCC
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 , 124.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Outliers	1 of 21161 reflections (0.005%)	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.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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, BMA, NAG, FUC

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(°)	Ideal(°)
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

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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	A	71	0	61	2	0
4	B	71	0	61	2	0
5	A	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:HH11	1:B:57:ARG:HG3	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:B:34:MET:HG2	1:B:40:THR:CG2	2.07	0.83
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
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

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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:O	1:A:220:TRP:HD1	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
3:L:26:THR:HG23	3:L:27:SER:H	1.62	0.65
1:B:166:PRO:HB3	1:B:187:PRO:HG3	1.78	0.65
1:A:306:PHE:O	1:A:387:LEU:HD11	1.96	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

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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
4:A:503:NAG:H4	2:I:100:TYR:CE1	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:315:THR:CG2	1:B:319:THR:OG1	2.52	0.57
1:B:20:TRP:CE2	1:B:286:ARG:HD3	2.38	0.57
1:B:57:ARG:NH1	1:B:57:ARG:CG	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:335:ILE:CD1	1:A:358:VAL:CG2	2.53	0.57
1:A:315:THR:CG2	1:A:319:THR:OG1	2.52	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:HD13	1:A:335:ILE:N	2.21	0.56
3:M:179:LEU:HD23	3:M:181:LEU:HG	1.86	0.56
1:B:289:MET:HG3	1:B:292:LEU:HD12	1.87	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:B:133:GLU:HG2	1:B:167:GLN:CB	2.38	0.54
1:A:315:THR:HG23	1:A:317:HIS:H	1.73	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

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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:220:TRP:NE1	1:A:232:ILE:HB	2.26	0.51
1:A:137:TYR:HB2	1:A:164:ILE:HG13	1.93	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:B:125:MET:HG2	1:B:199:LEU:HD11	1.94	0.49
1:A:288:ARG:HB3	1:A:290:ASP:OD1	2.12	0.49
3:M:49:TYR:C	3:M:49:TYR:CD1	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:I:155:ASN:OD1	2:I:195:ILE:HG13	2.13	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
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
4:B:503:NAG:H4	2:H:100:TYR:CE1	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:318:GLY:HA3	1:B:1393:ARG:HH21	1.80	0.46
1:B:308:ILE:HG13	1:B:308:ILE:H	1.52	0.46

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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:164:ILE:HD11	1:B:169:SER:HA	1.98	0.46
1:B:99:ARG:NH2	2:I:100(I):ASP:OD2	2.49	0.46
2:I:164:HIS:CE1	3:M:174:ALA:HB3	2.49	0.46
4:A:501:NAG:H81	1:B:102:GLY:HA2	1.98	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
2:I:82:MET:HE3	2:I:90:TYR:CE2	2.52	0.45
1:B:340:THR:HG22	1:B:347:VAL:HA	1.98	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
2:H:205:THR:CG2	2:I:203:SER:HB2	2.47	0.44
1:A:65:LEU:HD12	1:A:252:VAL:HG22	1.99	0.44
2:H:97:VAL:HG13	2:H:100(O):GLY:CA	2.48	0.44
2:I:82:MET:CE	2:I:90:TYR:CE2	3.00	0.44
1:B:65:LEU:HD12	1:B:252:VAL:HG22	2.00	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:B:133:GLU:OE1	1:B:167:GLN:HB2	2.18	0.43
1:A:345:ARG:HD2	1:A:346:HIS:NE2	2.34	0.43
1:B:345:ARG:HD2	1:B:346:HIS:NE2	2.34	0.43
2:I:51:ILE:HD11	2:I:78:LEU:HD11	1.95	0.43
1:B:164:ILE:HG13	1:B:165:THR:N	2.33	0.43
1:A:308:ILE:H	1:A:308:ILE:HG13	1.48	0.43
3:M:133:LEU:HD21	3:M:186:TRP:CZ3	2.53	0.43
1:B:33:THR:HG22	1:B:41:LEU:HB2	2.01	0.43
1:A:33:THR:HG21	1:A:43:PHE:HE1	1.84	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:THR:HG21	1:B:43:PHE:HE1	1.85	0.42
1:A:321:VAL:HG22	1:A:368:GLU:HG3	2.01	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
4:B:504:BMA:H3	3:L:56:SER:OG	2.20	0.42
1:B:133:GLU:HG2	1:B:167:GLN:HG2	2.01	0.42
1:A:329:ASP:HA	1:A:361:LYS:HE2	2.02	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
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
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:100(G):TYR:HA	2:H:100(H):PRO:HD3	1.99	0.41
2:I:82:MET:HE3	2:I:90:TYR:HE2	1.84	0.40
1:A:156:GLY:C	1:A:158:HIS:H	2.23	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: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
1:A:132:PRO:HG3	1:A:193:PHE:HB2	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%)	46 82

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	384/422 (91%)	367 (96%)	16 (4%)	1 (0%)	46	82
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%)	56	89

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%)	5	29
1	B	340/366 (93%)	295 (87%)	45 (13%)	5	30
2	H	190/236 (80%)	178 (94%)	12 (6%)	22	61
2	I	174/236 (74%)	164 (94%)	10 (6%)	25	65
3	L	180/185 (97%)	171 (95%)	9 (5%)	30	68
3	M	178/185 (96%)	167 (94%)	11 (6%)	23	62
All	All	1403/1574 (89%)	1270 (90%)	133 (10%)	11	44

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

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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

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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

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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

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

Some 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 carbohydrates 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	A	501	1,4	14,14,15	0.29	0	15,19,21	0.48	0
4	FUC	A	502	4	10,10,11	0.46	0	14,14,16	1.50	1 (7%)
4	NAG	A	503	4	14,14,15	0.29	0	15,19,21	1.50	2 (13%)
4	BMA	A	504	4	11,11,12	0.37	0	14,15,17	1.26	2 (14%)
4	MAN	A	505	4	11,11,12	0.30	0	14,15,17	0.81	1 (7%)
4	MAN	A	506	4	11,11,12	0.39	0	14,15,17	1.22	1 (7%)
5	NAG	A	567	1,5	14,14,15	0.25	0	15,19,21	0.81	1 (6%)
5	NAG	A	568	5	14,14,15	0.34	0	15,19,21	0.63	0
4	NAG	B	501	1,4	14,14,15	0.30	0	15,19,21	0.49	0
4	FUC	B	502	4	10,10,11	0.46	0	14,14,16	1.48	1 (7%)
4	NAG	B	503	4	14,14,15	0.27	0	15,19,21	1.43	2 (13%)
4	BMA	B	504	4	11,11,12	0.42	0	14,15,17	1.32	2 (14%)
4	MAN	B	505	4	11,11,12	0.27	0	14,15,17	0.90	1 (7%)
4	MAN	B	506	4	11,11,12	0.39	0	14,15,17	1.16	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	A	501	1,4	-	0/6/23/26	0/1/1/1
4	FUC	A	502	4	-	0/0/17/20	0/1/1/1
4	NAG	A	503	4	-	0/6/23/26	0/1/1/1
4	BMA	A	504	4	-	0/2/19/22	1/1/1/1
4	MAN	A	505	4	-	0/2/19/22	0/1/1/1
4	MAN	A	506	4	-	0/2/19/22	1/1/1/1
5	NAG	A	567	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	568	5	-	0/6/23/26	0/1/1/1
4	NAG	B	501	1,4	-	0/6/23/26	0/1/1/1
4	FUC	B	502	4	-	0/0/17/20	0/1/1/1
4	NAG	B	503	4	-	0/6/23/26	0/1/1/1
4	BMA	B	504	4	-	0/2/19/22	1/1/1/1
4	MAN	B	505	4	-	0/2/19/22	0/1/1/1
4	MAN	B	506	4	-	0/2/19/22	1/1/1/1

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	504	BMA	C1-C2-C3	2.10	112.03	109.54
4	B	504	BMA	C1-C2-C3	2.11	112.03	109.54
4	A	505	MAN	C1-O5-C5	2.36	115.25	112.25
4	A	503	NAG	O4-C4-C5	2.40	115.61	109.24
4	B	503	NAG	O4-C4-C5	2.50	115.88	109.24
5	A	567	NAG	C1-O5-C5	2.55	115.48	112.25
4	B	505	MAN	C1-O5-C5	2.78	115.78	112.25
4	B	504	BMA	C1-O5-C5	3.63	116.85	112.25
4	A	504	BMA	C1-O5-C5	3.72	116.97	112.25
4	B	506	MAN	C1-O5-C5	4.04	117.38	112.25
4	A	506	MAN	C1-O5-C5	4.17	117.54	112.25
4	B	502	FUC	C1-O5-C5	4.48	119.29	112.38
4	B	503	NAG	C1-O5-C5	4.55	118.03	112.25
4	A	502	FUC	C1-O5-C5	4.65	119.56	112.38
4	A	503	NAG	C1-O5-C5	4.89	118.46	112.25

There are no chirality outliers.

There are no torsion outliers.

All (4) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	504	BMA	C1-C2-C3-C4-C5-O5
4	B	504	BMA	C1-C2-C3-C4-C5-O5
4	B	506	MAN	C1-C2-C3-C4-C5-O5
4	A	506	MAN	C1-C2-C3-C4-C5-O5

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	501	NAG	1	0
4	A	503	NAG	1	0
4	B	503	NAG	1	0
4	B	504	BMA	1	0

5.6 Ligand geometry

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	$\# Z > 2$
6	NAG	B	567	1	14,14,15	0.36	0	15,19,21	0.59	0
7	SO4	H	581	-	4,4,4	0.20	0	6,6,6	0.21	0
7	SO4	I	581	-	4,4,4	0.19	0	6,6,6	0.15	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	-	0/6/23/26	0/1/1/1
7	SO4	H	581	-	-	0/0/0/0	0/0/0/0
7	SO4	I	581	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	391/422 (92%)	0.30	19 (4%) 33 24	100, 125, 158, 180	0
1	B	390/422 (92%)	0.22	9 (2%) 64 52	94, 125, 158, 177	0
2	H	225/283 (79%)	0.19	3 (1%) 79 70	94, 126, 146, 162	0
2	I	203/283 (71%)	0.51	21 (10%) 9 6	91, 122, 187, 206	0
3	L	213/218 (97%)	0.28	8 (3%) 44 33	90, 126, 160, 184	0
3	M	210/218 (96%)	0.41	16 (7%) 17 11	92, 128, 178, 193	0
All	All	1632/1846 (88%)	0.30	76 (4%) 35 26	90, 125, 172, 206	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	213	PRO	6.7
1	A	191	LEU	4.8
2	I	212	GLU	4.5
1	A	346	HIS	4.2
1	A	347	VAL	4.2
2	H	112	SER	3.9
3	L	210	THR	3.9
1	A	348	LEU	3.8
3	M	153	SER	3.7
3	M	184	GLU	3.6
2	I	208	ASP	3.5
2	I	206	LYS	3.4
1	B	96	MET	3.4
1	A	190	GLY	3.3
3	L	153	SER	3.3
1	B	16	SER	3.3
3	M	151	ALA	3.2
3	M	2	SER	3.0
1	A	342	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
2	I	163	VAL	3.0
1	A	343	GLU	3.0
3	M	157	LYS	2.9
1	A	341	ASP	2.8
1	B	190	GLY	2.8
3	M	182	THR	2.8
2	I	197	ASN	2.8
1	A	14	GLY	2.8
2	I	211	VAL	2.8
2	I	168	ALA	2.7
1	A	227	GLN	2.6
2	I	136	ALA	2.6
3	M	183	PRO	2.6
3	L	1	GLN	2.6
3	M	190	ARG	2.6
2	I	154	TRP	2.5
2	I	149	PRO	2.4
1	B	346	HIS	2.4
2	I	173	SER	2.4
1	A	184	GLU	2.4
2	I	137	ALA	2.4
3	L	191	SER	2.4
1	B	345	ARG	2.4
3	M	191	SER	2.3
2	I	10	GLY	2.3
1	A	331	SER	2.3
3	M	195	GLN	2.3
3	L	156	VAL	2.3
1	B	248	GLN	2.2
2	I	153	SER	2.2
2	I	199	ASN	2.2
2	I	180	SER	2.2
1	B	284	LYS	2.2
2	I	49	SER	2.2
1	A	345	ARG	2.2
1	B	343	GLU	2.2
3	M	192	TYR	2.1
3	M	185	GLN	2.1
2	I	155	ASN	2.1
1	A	96	MET	2.1
3	L	27	SER	2.1
1	A	325	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
3	L	190	ARG	2.1
3	M	154	SER	2.1
2	H	210	ARG	2.1
3	M	150	LYS	2.1
2	I	126	PRO	2.1
1	A	186	SER	2.1
3	M	199	GLU	2.1
1	B	348	LEU	2.1
2	H	9	GLY	2.1
1	A	188	ARG	2.1
2	I	11	LEU	2.0
1	A	340	THR	2.0
3	L	143	GLY	2.0
3	M	27	SER	2.0
1	A	192	ASP	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	NAG	A	567	14/15	0.49	0.63	3.00	169,172,173,174	0
4	MAN	A	506	11/12	0.76	0.50	1.83	151,155,160,161	0
4	MAN	B	506	11/12	0.83	0.43	1.30	157,161,163,163	0
4	NAG	A	501	14/15	0.85	0.28	0.03	110,112,124,128	0
4	NAG	B	501	14/15	0.92	0.23	-0.46	102,108,118,124	0
4	FUC	B	502	10/11	0.88	0.32	-	129,133,135,137	0
4	NAG	A	503	14/15	0.89	0.30	-	117,121,130,131	0
5	NAG	A	568	14/15	0.21	0.76	-	169,174,177,177	0
4	NAG	B	503	14/15	0.90	0.24	-	112,115,118,125	0
4	BMA	B	504	11/12	0.89	0.19	-	132,133,143,151	0
4	MAN	A	505	11/12	0.90	0.32	-	125,127,130,131	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	FUC	A	502	10/11	0.77	0.43	-	130,133,135,137	0
4	MAN	B	505	11/12	0.88	0.27	-	133,136,137,138	0
4	BMA	A	504	11/12	0.93	0.15	-	125,127,135,144	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
7	SO4	I	581	5/5	0.59	0.26	-0.39	169,169,170,170	0
7	SO4	H	581	5/5	0.71	0.25	-0.88	144,144,145,146	0
6	NAG	B	567	14/15	0.75	0.26	-	154,156,158,158	0

6.5 Other polymers [i](#)

There are no such residues in this entry.