



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 8, 2016 – 11:25 AM EST

PDB ID : 5T33
Title : Crystal structure of strain-specific glycan-dependent CD4 binding site-directed neutralizing antibody CAP257-RH1, in complex with HIV-1 strain RHPA gp120 core with an oligomannose N276 glycan.
Authors : Wibmer, C.K.; Gorman, J.; Kwong, P.D.
Deposited on : 2016-08-24
Resolution : 3.21 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028320
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) : rb-20028320

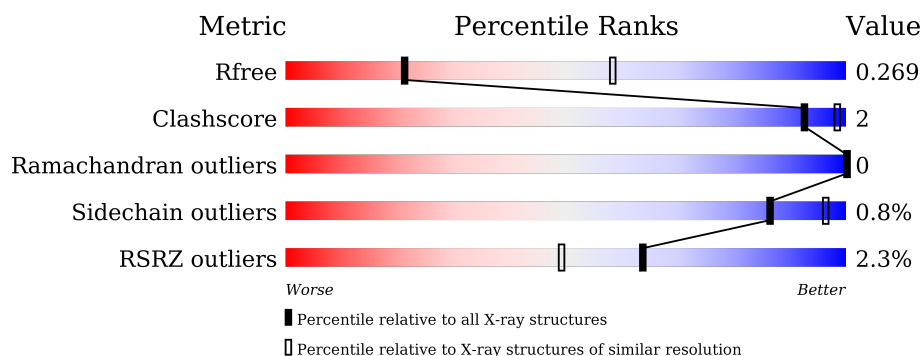
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.21 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	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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	H	223	<div> <div style="width: 94%;"></div> <div style="width: 6%;"></div> </div> <div>94% 6%</div>
2	L	214	<div> <div style="width: 95%;"></div> <div style="width: 5%;"></div> </div> <div>95% . .</div>
3	G	357	<div> <div style="width: 4%;"></div> <div style="width: 83%;"></div> <div style="width: 6%;"></div> <div style="width: 11%;"></div> </div> <div>4% 83% 6% 11%</div>

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	G	1088	-	-	-	X
6	MAN	G	1281	-	-	-	X
6	MAN	G	1282	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 11978 atoms, of which 5904 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 CAP257-RH1 heavy chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	H	223	Total	C	H	N	O	S	0	0	0
			3330	1060	1650	289	324	7			

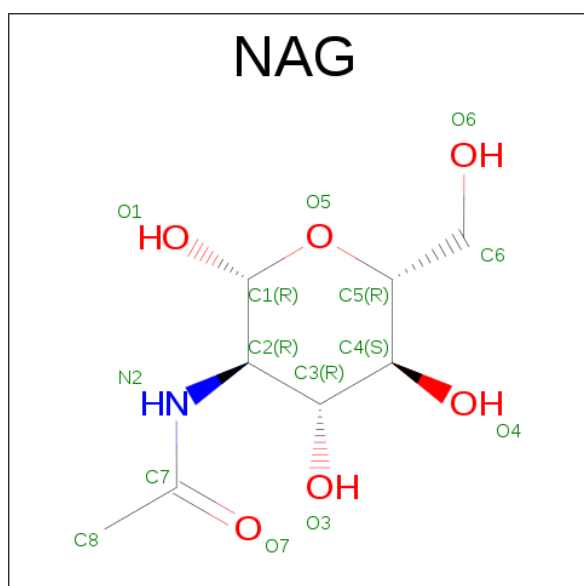
- Molecule 2 is a protein called CAP257-RH1 light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	L	212	Total	C	H	N	O	S	0	0	0
			3139	1006	1538	264	324	7			

- Molecule 3 is a protein called RHPA gp120 core.

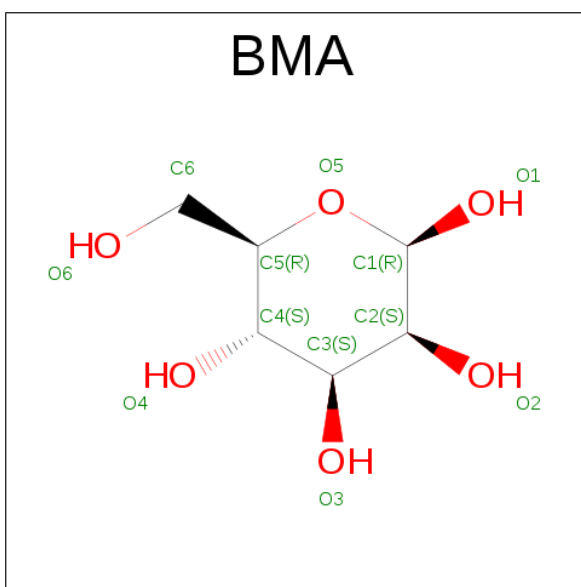
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	G	319	Total	C	H	N	O	S	0	1	0
			4976	1591	2456	436	472	21			

- Molecule 4 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C₈H₁₅NO₆).



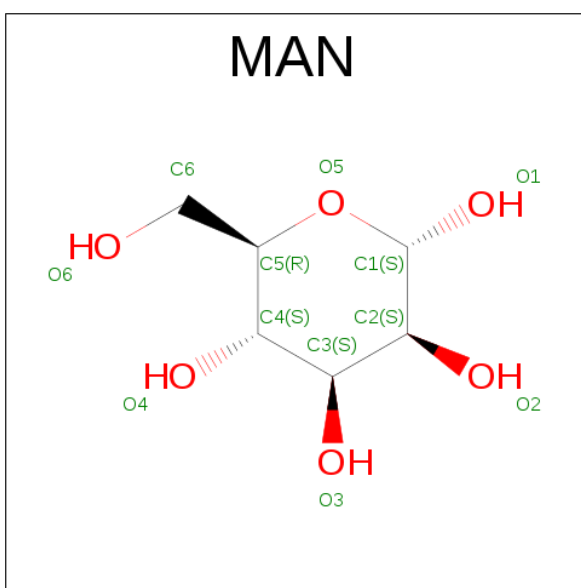
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	G	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
4	G	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
4	G	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
4	G	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
4	G	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
4	G	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
4	G	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
4	G	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
4	G	1	Total	C	H	N	O	0	0
			27	8	13	1	5		
4	G	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
4	G	1	Total	C	H	N	O	0	0
			27	8	13	1	5		

- Molecule 5 is BETA-D-MANNOSE (three-letter code: BMA) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	G	1	Total	C	H	O	0	0
			21	6	10	5		
5	G	1	Total	C	H	O	0	0
			20	6	9	5		

- Molecule 6 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	G	1	Total	C	H	O	0	0
			22	6	11	5		
6	G	1	Total	C	H	O	0	0
			22	6	11	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	G	1	Total	C	H	O	0	0
			20	6	9	5		
6	G	1	Total	C	H	O	0	0
			22	6	11	5		
6	G	1	Total	C	H	O	0	0
			22	6	11	5		

3 Residue-property plots [i](#)

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: CAP257-RH1 heavy chain

Chain H:  94% 6%




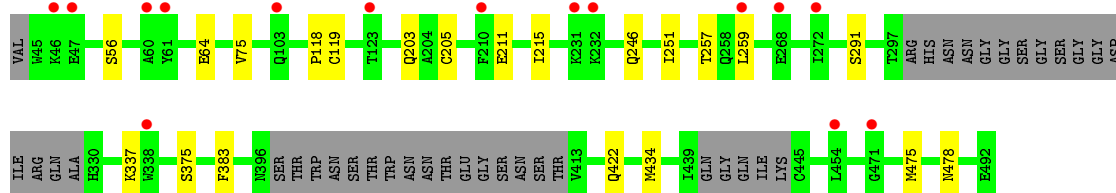
- Molecule 2: CAP257-RH1 light chain

Chain L:  95% ..



- Molecule 3: RHPA gp120 core

Chain G:  4% 83% 6% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.18Å 71.14Å 190.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.66 – 3.21 47.66 – 3.21	Depositor EDS
% Data completeness (in resolution range)	82.0 (47.66-3.21) 79.5 (47.66-3.21)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 3.19Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.229 , 0.269 0.224 , 0.269	Depositor DCC
R_{free} test set	630 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	86.0	Xtriage
Anisotropy	0.385	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.054 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11978	wwPDB-VP
Average B, all atoms (Å ²)	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	H	0.26	0/1721	0.46	0/2341
2	L	0.26	0/1643	0.43	0/2243
3	G	0.24	0/2575	0.42	0/3495
All	All	0.25	0/5939	0.44	0/8079

There are no bond length outliers.

There are no bond angle outliers.

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	H	1680	1650	1650	6	0
2	L	1601	1538	1538	6	0
3	G	2520	2456	2459	12	0
4	G	196	188	176	2	0
5	G	22	19	17	0	0
6	G	55	53	48	0	0
All	All	6074	5904	5888	24	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 2.

All (24) 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)
3:G:118:PRO:O	3:G:203:GLN:NE2	2.36	0.58
1:H:6:GLU:N	1:H:6:GLU:OE1	2.35	0.58
3:G:246:GLN:OE1	3:G:246:GLN:N	2.42	0.52
3:G:211:GLU:OE2	4:G:1262:NAG:O6	2.29	0.50
4:G:1088:NAG:H83	4:G:1088:NAG:H3	1.94	0.50
1:H:66:ARG:NH1	1:H:82(B):SER:O	2.48	0.47
1:H:201:LYS:N	1:H:202:PRO:CD	2.79	0.46
3:G:475:MET:O	3:G:478:ASN:N	2.49	0.46
2:L:48:ILE:CG2	2:L:51:GLY:O	2.64	0.46
2:L:52:ASN:OD1	2:L:52:ASN:N	2.45	0.46
1:H:143:LYS:HZ1	2:L:131:THR:HG21	1.82	0.45
3:G:119:CYS:HB2	3:G:434:MET:O	2.18	0.44
1:H:156:SER:N	1:H:197:ASN:OD1	2.40	0.44
2:L:131:THR:HG22	2:L:179:SER:HA	1.99	0.44
3:G:422:GLN:OE1	3:G:422:GLN:N	2.50	0.43
3:G:215:ILE:N	3:G:251:ILE:O	2.42	0.43
3:G:64:GLU:HG3	3:G:211:GLU:HB3	2.01	0.43
3:G:257:THR:O	3:G:259:LEU:N	2.48	0.42
1:H:5:VAL:O	1:H:22:CYS:HA	2.19	0.41
3:G:375:SER:HA	3:G:383:PHE:O	2.19	0.41
2:L:132:LEU:HD12	2:L:132:LEU:N	2.35	0.41
3:G:291:SER:O	3:G:337:LYS:NZ	2.40	0.41
3:G:56:SER:N	3:G:75:VAL:O	2.53	0.40
2:L:184:GLN:O	2:L:191:TYR:OH	2.39	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	221/223 (99%)	206 (93%)	15 (7%)	0	100	100
2	L	210/214 (98%)	193 (92%)	17 (8%)	0	100	100
3	G	311/357 (87%)	282 (91%)	29 (9%)	0	100	100
All	All	742/794 (94%)	681 (92%)	61 (8%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	187/187 (100%)	184 (98%)	3 (2%)	70	91
2	L	180/182 (99%)	179 (99%)	1 (1%)	90	97
3	G	286/316 (90%)	285 (100%)	1 (0%)	94	98
All	All	653/685 (95%)	648 (99%)	5 (1%)	86	96

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	92	CYS
1	H	150	VAL
1	H	196	CYS
2	L	55	PHE
3	G	205	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

21 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$
4	NAG	G	1049	3	14,14,15	1.40	1 (7%)	15,19,21	1.05	1 (6%)
4	NAG	G	1088	3	14,14,15	1.38	2 (14%)	15,19,21	1.88	4 (26%)
4	NAG	G	1234	3,4	14,14,15	1.45	1 (7%)	15,19,21	0.98	1 (6%)
4	NAG	G	1235	4	14,14,15	1.40	1 (7%)	15,19,21	1.01	0
4	NAG	G	1241	3	14,14,15	1.42	1 (7%)	15,19,21	1.21	1 (6%)
4	NAG	G	1262	3,4	14,14,15	1.36	1 (7%)	15,19,21	1.11	1 (6%)
4	NAG	G	1263	5,4	14,14,15	1.43	1 (7%)	15,19,21	1.13	1 (6%)
5	BMA	G	1264	4,6	11,11,12	1.63	2 (18%)	15,15,17	0.85	0
6	MAN	G	1265	5	11,11,12	1.74	2 (18%)	15,15,17	1.32	1 (6%)
4	NAG	G	1276	3,4	14,14,15	1.36	1 (7%)	15,19,21	1.19	1 (6%)
4	NAG	G	1277	5,4	14,14,15	1.33	1 (7%)	15,19,21	1.23	1 (6%)
5	BMA	G	1278	4,6	11,11,12	1.66	2 (18%)	15,15,17	0.77	0
6	MAN	G	1279	5	11,11,12	1.71	2 (18%)	15,15,17	0.82	0
6	MAN	G	1280	5,6	11,11,12	1.70	2 (18%)	15,15,17	0.76	0
6	MAN	G	1281	6	11,11,12	1.64	2 (18%)	15,15,17	0.90	0
6	MAN	G	1282	6	11,11,12	1.70	2 (18%)	15,15,17	0.90	0
4	NAG	G	1289	3	14,14,15	1.39	1 (7%)	15,19,21	1.23	1 (6%)
4	NAG	G	1295	3,4	14,14,15	1.37	1 (7%)	15,19,21	1.12	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	G	1296	4	14,14,15	1.38	1 (7%)	15,19,21	1.08	2 (13%)
4	NAG	G	1386	3	14,14,15	1.41	1 (7%)	15,19,21	1.18	1 (6%)
4	NAG	G	1448	3	14,14,15	1.38	1 (7%)	15,19,21	1.01	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	G	1049	3	-	0/6/23/26	0/1/1/1
4	NAG	G	1088	3	-	0/6/23/26	0/1/1/1
4	NAG	G	1234	3,4	-	0/6/23/26	0/1/1/1
4	NAG	G	1235	4	-	0/6/23/26	0/1/1/1
4	NAG	G	1241	3	-	0/6/23/26	0/1/1/1
4	NAG	G	1262	3,4	-	0/6/23/26	0/1/1/1
4	NAG	G	1263	5,4	-	0/6/23/26	0/1/1/1
5	BMA	G	1264	4,6	-	0/2/19/22	0/1/1/1
6	MAN	G	1265	5	-	0/2/19/22	0/1/1/1
4	NAG	G	1276	3,4	-	0/6/23/26	0/1/1/1
4	NAG	G	1277	5,4	-	0/6/23/26	0/1/1/1
5	BMA	G	1278	4,6	-	0/2/19/22	0/1/1/1
6	MAN	G	1279	5	-	0/2/19/22	0/1/1/1
6	MAN	G	1280	5,6	-	0/2/19/22	0/1/1/1
6	MAN	G	1281	6	-	0/2/19/22	0/1/1/1
6	MAN	G	1282	6	-	0/2/19/22	0/1/1/1
4	NAG	G	1289	3	-	0/6/23/26	0/1/1/1
4	NAG	G	1295	3,4	-	0/6/23/26	0/1/1/1
4	NAG	G	1296	4	-	0/6/23/26	0/1/1/1
4	NAG	G	1386	3	-	0/6/23/26	0/1/1/1
4	NAG	G	1448	3	-	0/6/23/26	0/1/1/1

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	G	1282	MAN	C2-C3	-3.34	1.48	1.52
6	G	1279	MAN	C2-C3	-3.34	1.48	1.52
6	G	1280	MAN	C2-C3	-3.28	1.48	1.52
6	G	1281	MAN	C2-C3	-3.18	1.48	1.52
6	G	1265	MAN	C2-C3	-3.08	1.48	1.52
5	G	1278	BMA	C2-C3	-3.03	1.48	1.52
5	G	1264	BMA	C2-C3	-3.02	1.48	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	1088	NAG	C7-N2	2.07	1.42	1.34
4	G	1277	NAG	O5-C1	2.57	1.47	1.43
4	G	1276	NAG	O5-C1	2.75	1.48	1.43
4	G	1262	NAG	O5-C1	2.81	1.48	1.43
4	G	1241	NAG	O5-C1	2.85	1.48	1.43
4	G	1295	NAG	O5-C1	2.86	1.48	1.43
4	G	1448	NAG	O5-C1	2.88	1.48	1.43
4	G	1296	NAG	O5-C1	2.90	1.48	1.43
4	G	1386	NAG	O5-C1	2.93	1.48	1.43
4	G	1289	NAG	O5-C1	2.94	1.48	1.43
4	G	1263	NAG	O5-C1	2.99	1.48	1.43
4	G	1049	NAG	O5-C1	3.01	1.48	1.43
4	G	1088	NAG	O5-C1	3.01	1.48	1.43
4	G	1235	NAG	O5-C1	3.03	1.48	1.43
4	G	1234	NAG	O5-C1	3.10	1.48	1.43
6	G	1281	MAN	O5-C1	3.54	1.49	1.43
6	G	1280	MAN	O5-C1	3.63	1.49	1.43
6	G	1279	MAN	O5-C1	3.67	1.49	1.43
6	G	1282	MAN	O5-C1	3.67	1.49	1.43
5	G	1278	BMA	O5-C1	3.71	1.49	1.43
5	G	1264	BMA	O5-C1	3.72	1.49	1.43
6	G	1265	MAN	O5-C1	4.02	1.50	1.43

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	1088	NAG	O7-C7-C8	-2.77	116.97	122.07
4	G	1088	NAG	O7-C7-N2	-2.42	116.91	121.84
4	G	1295	NAG	C2-N2-C7	-2.28	120.14	123.11
4	G	1296	NAG	C2-N2-C7	-2.27	120.16	123.11
4	G	1234	NAG	C2-N2-C7	-2.14	120.32	123.11
4	G	1049	NAG	C2-N2-C7	-2.09	120.39	123.11
4	G	1296	NAG	C8-C7-N2	2.09	120.11	116.10
4	G	1295	NAG	C8-C7-N2	2.13	120.17	116.10
4	G	1088	NAG	C2-N2-C7	2.37	126.19	123.11
4	G	1276	NAG	C8-C7-N2	2.94	121.73	116.10
4	G	1277	NAG	C8-C7-N2	2.99	121.83	116.10
4	G	1262	NAG	C8-C7-N2	3.02	121.88	116.10
4	G	1241	NAG	C8-C7-N2	3.02	121.89	116.10
4	G	1263	NAG	C8-C7-N2	3.06	121.96	116.10
4	G	1289	NAG	C8-C7-N2	3.09	122.02	116.10
4	G	1386	NAG	C8-C7-N2	3.14	122.12	116.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	G	1265	MAN	C1-C2-C3	3.76	114.11	109.55
4	G	1088	NAG	C8-C7-N2	5.23	126.12	116.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	1088	NAG	1	0
4	G	1262	NAG	1	0

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	H	223/223 (100%)	-0.04	1 (0%) 93 90	68, 85, 110, 154	0
2	L	212/214 (99%)	-0.10	1 (0%) 91 87	65, 83, 105, 127	0
3	G	319/357 (89%)	0.26	15 (4%) 35 22	79, 129, 164, 191	0
All	All	754/794 (94%)	0.07	17 (2%) 64 49	65, 98, 153, 191	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	61	TYR	6.1
3	G	232	LYS	4.7
3	G	47	GLU	4.3
1	H	189	LEU	3.6
3	G	268	GLU	3.6
3	G	123	THR	2.9
3	G	60	ALA	2.9
3	G	103	GLN	2.7
3	G	46	LYS	2.6
3	G	210	PHE	2.4
2	L	2	TYR	2.3
3	G	471	GLY	2.3
3	G	259	LEU	2.1
3	G	454	LEU	2.1
3	G	272	ILE	2.1
3	G	338	TRP	2.0
3	G	231	LYS	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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
6	MAN	G	1282	11/12	0.78	0.51	10.66	124,128,153,153	0
4	NAG	G	1088	14/15	0.65	0.77	6.61	173,182,218,218	0
6	MAN	G	1281	11/12	0.78	0.32	2.97	99,105,126,127	0
4	NAG	G	1386	14/15	0.79	0.30	1.06	135,144,170,172	0
4	NAG	G	1448	14/15	0.83	0.22	0.61	148,153,183,183	0
4	NAG	G	1295	14/15	0.88	0.30	0.48	156,164,194,196	0
4	NAG	G	1049	14/15	0.81	0.36	0.32	130,148,176,176	0
4	NAG	G	1241	14/15	0.81	0.28	-0.06	151,162,194,194	0
6	MAN	G	1279	11/12	0.90	0.21	-0.10	117,120,143,145	0
4	NAG	G	1289	14/15	0.84	0.29	-0.16	145,153,182,183	0
4	NAG	G	1263	14/15	0.85	0.17	-0.51	143,153,180,184	0
4	NAG	G	1234	14/15	0.82	0.18	-0.58	130,135,161,162	0
4	NAG	G	1262	14/15	0.94	0.14	-0.90	133,147,172,176	0
6	MAN	G	1265	11/12	0.58	0.37	-	135,153,180,183	0
4	NAG	G	1235	14/15	0.82	0.29	-	145,158,190,190	0
5	BMA	G	1264	11/12	0.66	0.20	-	141,146,175,176	0
4	NAG	G	1276	14/15	0.91	0.17	-	109,119,143,145	0
4	NAG	G	1277	14/15	0.93	0.17	-	111,115,138,138	0
6	MAN	G	1280	11/12	0.88	0.14	-	108,113,136,136	0
5	BMA	G	1278	11/12	0.91	0.12	-	108,114,137,137	0
4	NAG	G	1296	14/15	0.84	0.44	-	167,172,207,207	0

6.5 Other polymers ⓘ

There are no such residues in this entry.