



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2020 – 12:58 PM BST

PDB ID : 4LST
Title : Crystal structure of broadly and potently neutralizing antibody VRC01 in complex with HIV-1 clade C strain ZM176.66 gp120
Authors : Zhou, T.; Moquin, S.; Kwong, P.D.
Deposited on : 2013-07-23
Resolution : 2.55 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

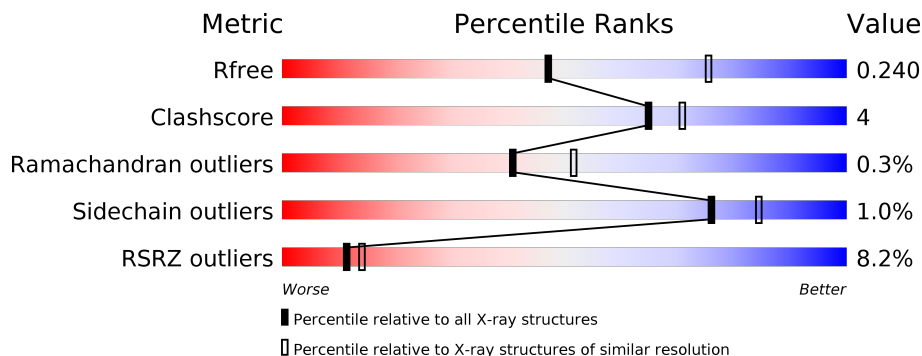
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	G	355	<div style="display: flex; align-items: center;"> <div style="width: 13%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 83%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div>
2	H	224	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 88%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div>
3	L	210	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 90%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow;"></div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6171 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 GP120 of HIV-1 clade C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	G	339	2657	1666	461	509	21	0	0	0

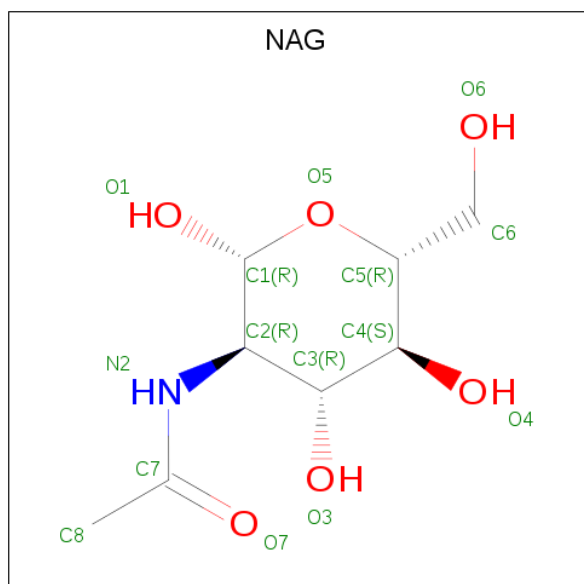
- Molecule 2 is a protein called HEAVY CHAIN OF ANTIBODY VRC01.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	222	1699	1072	295	321	11	0	0	0

- Molecule 3 is a protein called LIGHT CHAIN OF ANTIBODY VRC01.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	L	210	1632	1022	279	326	5	0	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

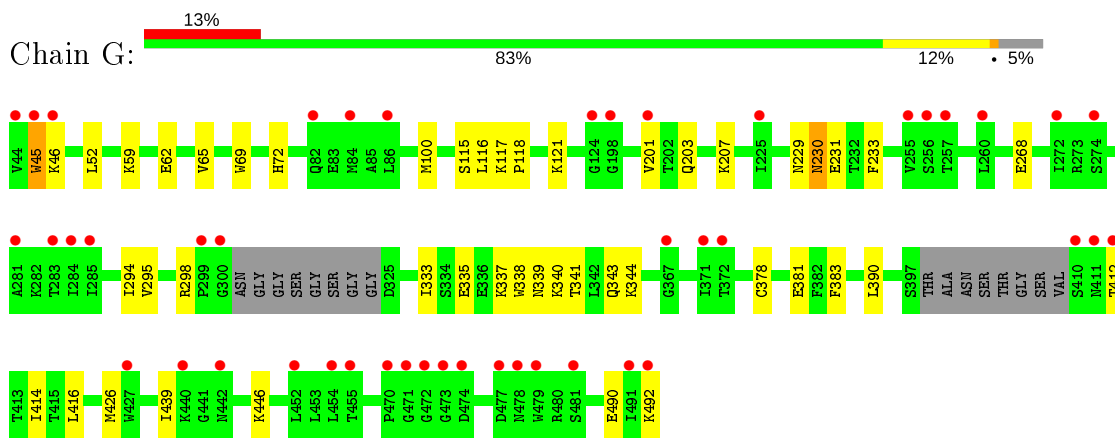
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	25	Total	O	0	0
			25	25		
5	H	27	Total	O	0	0
			27	27		
5	L	19	Total	O	0	0
			19	19		

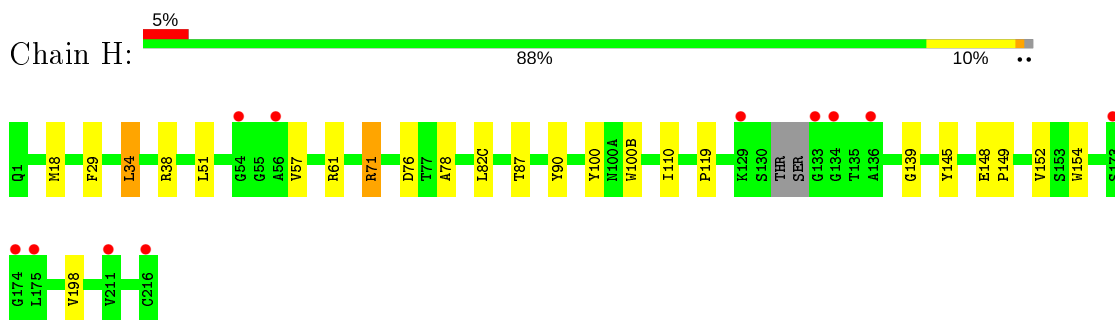
3 Residue-property plots [i](#)

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.

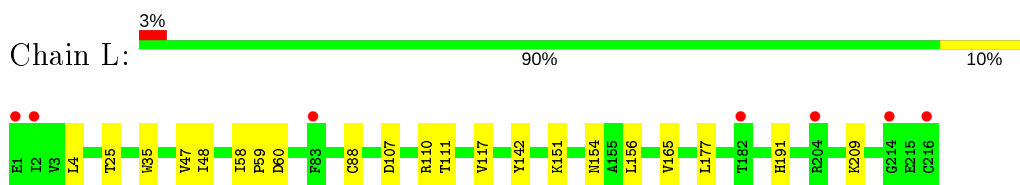
- Molecule 1: ENVELOPE GLYCOPROTEIN GP120 of HIV-1 clade C



- Molecule 2: HEAVY CHAIN OF ANTIBODY VRC01



- Molecule 3: LIGHT CHAIN OF ANTIBODY VRC01



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.11Å 77.98Å 200.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.68 – 2.55 40.68 – 2.55	Depositor EDS
% Data completeness (in resolution range)	91.9 (40.68-2.55) 91.9 (40.68-2.55)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.69 (at 2.54Å)	Xtrriage
Refinement program	PHENIX dev_998	Depositor
R, R_{free}	0.190 , 0.239 0.191 , 0.240	Depositor DCC
R_{free} test set	1653 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	74.0	Xtrriage
Anisotropy	0.527	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 79.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6171	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.95% 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 [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	G	0.27	0/2714	0.46	0/3682
2	H	0.27	0/1743	0.50	0/2369
3	L	0.28	0/1669	0.47	0/2265
All	All	0.27	0/6126	0.47	0/8316

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 [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	G	2657	0	2585	26	0
2	H	1699	0	1671	12	0
3	L	1632	0	1574	11	0
4	G	112	0	104	3	0
5	G	25	0	0	0	0
5	H	27	0	0	1	0
5	L	19	0	0	0	0
All	All	6171	0	5934	49	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 4.

All (49) 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)
2:H:61:ARG:NH1	5:H:306:HOH:O	2.22	0.73
3:L:47:VAL:HG12	3:L:48:ILE:HG12	1.75	0.68
1:G:335:GLU:HG2	1:G:414:ILE:HG13	1.77	0.67
3:L:110:ARG:NH1	3:L:111:THR:O	2.27	0.67
1:G:340:LYS:HG2	1:G:344:LYS:HD2	1.79	0.64
2:H:18:MET:HG2	2:H:82(C):LEU:HD11	1.80	0.64
1:G:46:LYS:HD2	1:G:490:GLU:O	1.97	0.64
2:H:119:PRO:HB3	2:H:145:TYR:HB3	1.80	0.62
2:H:51:LEU:HD21	2:H:71:ARG:HB3	1.82	0.61
1:G:231:GLU:HB2	1:G:268:GLU:HG3	1.83	0.60
1:G:295:VAL:HG22	1:G:446:LYS:HG2	1.88	0.55
1:G:294:ILE:HD12	1:G:333:ILE:HD11	1.90	0.53
2:H:148:GLU:HB2	2:H:149:PRO:HA	1.90	0.53
1:G:230:ASN:HB3	1:G:233:PHE:HB2	1.91	0.51
2:H:152:VAL:HG22	2:H:198:VAL:HG22	1.93	0.50
1:G:69:TRP:HA	1:G:72:HIS:CE1	2.47	0.49
1:G:117:LYS:HB3	1:G:203:GLN:HE21	1.78	0.48
3:L:60:ASP:N	3:L:60:ASP:OD1	2.47	0.48
1:G:339:ASN:O	1:G:343:GLN:HG3	2.14	0.47
1:G:233:PHE:HD1	4:G:503:NAG:H82	1.78	0.47
1:G:390:LEU:HD21	1:G:416:LEU:HD11	1.96	0.47
2:H:100:TYR:CE2	2:H:100(B):TRP:HB2	2.50	0.47
1:G:298:ARG:NH2	1:G:439:ILE:O	2.43	0.46
3:L:151:LYS:NZ	3:L:154:ASN:HA	2.30	0.46
1:G:59:LYS:HB2	1:G:62:GLU:HB2	1.98	0.46
2:H:34:LEU:HD23	2:H:78:ALA:HB2	1.99	0.45
1:G:378:CYS:HB3	1:G:383:PHE:CE1	2.52	0.45
3:L:107:ASP:OD2	3:L:142:TYR:OH	2.29	0.44
1:G:233:PHE:CD1	4:G:503:NAG:H82	2.52	0.44
2:H:38:ARG:HD3	2:H:90:TYR:CE2	2.53	0.44
3:L:35:TRP:CZ3	3:L:88:CYS:HB3	2.52	0.43
1:G:207:LYS:NZ	1:G:381:GLU:OE2	2.42	0.43
2:H:139:GLY:HA2	2:H:154:TRP:CH2	2.53	0.43
1:G:65:VAL:HB	1:G:115:SER:HB2	2.00	0.43
3:L:58:ILE:HA	3:L:59:PRO:HD3	1.82	0.43
1:G:45:TRP:O	1:G:46:LYS:HG3	2.18	0.42
3:L:4:LEU:HD12	3:L:25:THR:HB	2.01	0.42
2:H:87:THR:HG23	2:H:110:ILE:HA	2.01	0.42
1:G:52:LEU:HD11	1:G:100:MET:HG2	2.00	0.42
3:L:151:LYS:HG2	3:L:156:LEU:HD23	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:337:LYS:O	1:G:341:THR:HG23	2.20	0.42
1:G:121:LYS:HE3	1:G:426:MET:HE1	2.01	0.41
1:G:118:PRO:HD2	1:G:203:GLN:HE22	1.86	0.41
1:G:121:LYS:HB3	1:G:201:VAL:HG23	2.03	0.41
3:L:117:VAL:O	3:L:209:LYS:HE2	2.20	0.41
1:G:338:TRP:CZ2	1:G:390:LEU:HB3	2.56	0.41
2:H:29:PHE:CD2	2:H:76:ASP:HA	2.56	0.40
3:L:165:VAL:HG22	3:L:177:LEU:HD12	2.04	0.40
1:G:229:ASN:HB3	4:G:504:NAG:HN2	1.87	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	G	333/355 (94%)	319 (96%)	13 (4%)	1 (0%)	41	51
2	H	218/224 (97%)	208 (95%)	10 (5%)	0	100	100
3	L	208/210 (99%)	203 (98%)	4 (2%)	1 (0%)	29	40
All	All	759/789 (96%)	730 (96%)	27 (4%)	2 (0%)	41	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	412	THR
3	L	191	HIS

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	G	300/309 (97%)	296 (99%)	4 (1%)	69	80
2	H	191/193 (99%)	188 (98%)	3 (2%)	62	77
3	L	182/182 (100%)	182 (100%)	0	100	100
All	All	673/684 (98%)	666 (99%)	7 (1%)	76	84

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	45	TRP
1	G	116	LEU
1	G	230	ASN
1	G	492	LYS
2	H	34	LEU
2	H	57	VAL
2	H	71	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	72	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](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry

8 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	502	1	14,14,15	0.58	0	17,19,21	0.88	1 (5%)
4	NAG	G	506	1	14,14,15	0.52	0	17,19,21	0.76	1 (5%)
4	NAG	G	507	1	14,14,15	0.51	0	17,19,21	0.86	1 (5%)
4	NAG	G	504	1	14,14,15	0.61	0	17,19,21	1.29	2 (11%)
4	NAG	G	503	1	14,14,15	1.09	1 (7%)	17,19,21	2.68	3 (17%)
4	NAG	G	508	1	14,14,15	0.58	0	17,19,21	0.70	0
4	NAG	G	501	1	14,14,15	0.47	0	17,19,21	1.25	2 (11%)
4	NAG	G	505	1	14,14,15	0.46	0	17,19,21	0.92	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	502	1	-	2/6/23/26	0/1/1/1
4	NAG	G	506	1	-	4/6/23/26	0/1/1/1
4	NAG	G	507	1	-	2/6/23/26	0/1/1/1
4	NAG	G	504	1	-	4/6/23/26	0/1/1/1
4	NAG	G	503	1	-	4/6/23/26	0/1/1/1
4	NAG	G	508	1	-	4/6/23/26	0/1/1/1
4	NAG	G	501	1	-	0/6/23/26	0/1/1/1
4	NAG	G	505	1	-	3/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	503	NAG	O7-C7	-3.59	1.15	1.23

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	503	NAG	C1-C2-N2	-9.38	94.47	110.49
4	G	501	NAG	C2-N2-C7	-3.43	118.02	122.90
4	G	504	NAG	C1-O5-C5	3.22	116.56	112.19
4	G	503	NAG	C4-C3-C2	-3.22	106.30	111.02
4	G	504	NAG	O5-C1-C2	2.81	115.72	111.29
4	G	503	NAG	O5-C5-C6	2.49	111.11	107.20
4	G	501	NAG	C1-O5-C5	2.48	115.55	112.19
4	G	506	NAG	C1-O5-C5	2.37	115.40	112.19
4	G	507	NAG	C1-O5-C5	2.22	115.20	112.19
4	G	502	NAG	O5-C5-C6	2.13	110.54	107.20

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	506	NAG	C8-C7-N2-C2
4	G	507	NAG	C8-C7-N2-C2
4	G	507	NAG	O7-C7-N2-C2
4	G	504	NAG	C8-C7-N2-C2
4	G	504	NAG	O7-C7-N2-C2
4	G	503	NAG	C1-C2-N2-C7
4	G	508	NAG	C8-C7-N2-C2
4	G	508	NAG	O7-C7-N2-C2
4	G	505	NAG	O7-C7-N2-C2
4	G	506	NAG	O7-C7-N2-C2
4	G	505	NAG	C8-C7-N2-C2
4	G	508	NAG	O5-C5-C6-O6
4	G	503	NAG	C8-C7-N2-C2
4	G	503	NAG	O7-C7-N2-C2
4	G	506	NAG	O5-C5-C6-O6
4	G	508	NAG	C4-C5-C6-O6
4	G	502	NAG	C8-C7-N2-C2
4	G	504	NAG	C1-C2-N2-C7
4	G	505	NAG	O5-C5-C6-O6
4	G	502	NAG	O7-C7-N2-C2
4	G	504	NAG	C3-C2-N2-C7
4	G	503	NAG	C3-C2-N2-C7
4	G	506	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	504	NAG	1	0
4	G	503	NAG	2	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

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	G	339/355 (95%)	0.84	45 (13%) 3 4	70, 99, 153, 185	0
2	H	222/224 (99%)	0.20	11 (4%) 28 34	67, 98, 148, 208	0
3	L	210/210 (100%)	0.09	7 (3%) 46 53	73, 102, 154, 199	0
All	All	771/789 (97%)	0.45	63 (8%) 11 14	67, 99, 154, 208	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	472	GLY	6.9
1	G	45	TRP	6.7
1	G	124	GLY	6.0
2	H	216	CYS	5.9
2	H	175	LEU	5.8
2	H	211	VAL	5.5
1	G	410	SER	5.4
3	L	216	CYS	5.1
1	G	492	LYS	5.0
1	G	473	GLY	4.7
3	L	1	GLU	4.5
1	G	471	GLY	4.2
1	G	440	LYS	4.1
3	L	2	ILE	3.5
1	G	491	ILE	3.5
1	G	283	THR	3.3
1	G	300	GLY	3.3
1	G	372	THR	3.1
1	G	284	ILE	3.1
1	G	481	SER	3.1
2	H	136	ALA	3.1
1	G	82	GLN	3.0
1	G	257	THR	3.0

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Mol	Chain	Res	Type	RSRZ
3	L	83	PHE	2.9
3	L	204	ARG	2.9
1	G	44	VAL	2.9
1	G	46	LYS	2.9
1	G	299	PRO	2.9
1	G	412	THR	2.8
2	H	54	GLY	2.8
1	G	411	ASN	2.8
1	G	455	THR	2.7
2	H	56	ALA	2.6
1	G	371	ILE	2.6
1	G	474	ASP	2.6
1	G	84	MET	2.6
1	G	285	ILE	2.5
2	H	174	GLY	2.5
1	G	256	SER	2.5
2	H	134	GLY	2.5
1	G	454	LEU	2.5
1	G	274	SER	2.4
1	G	198	GLY	2.4
1	G	272	ILE	2.4
1	G	470	PRO	2.4
2	H	133	GLY	2.3
1	G	86	LEU	2.3
3	L	182	THR	2.3
1	G	479	TRP	2.3
2	H	129	LYS	2.2
1	G	477	ASP	2.2
1	G	442	ASN	2.2
1	G	427	TRP	2.2
1	G	281	ALA	2.2
1	G	478	ASN	2.2
1	G	367	GLY	2.1
1	G	260	LEU	2.1
3	L	214	GLY	2.1
1	G	255	VAL	2.1
1	G	225	ILE	2.1
1	G	452	LEU	2.1
2	H	173	SER	2.1
1	G	201	VAL	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](#)

There are no monosaccharides in this entry.

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
4	NAG	G	503	14/15	0.79	0.17	106,128,132,136	0
4	NAG	G	506	14/15	0.82	0.42	154,173,184,191	0
4	NAG	G	504	14/15	0.84	0.31	130,148,161,161	0
4	NAG	G	507	14/15	0.92	0.23	139,160,163,164	0
4	NAG	G	508	14/15	0.92	0.30	144,160,167,168	0
4	NAG	G	505	14/15	0.93	0.19	92,115,138,140	0
4	NAG	G	502	14/15	0.94	0.18	107,139,155,170	0
4	NAG	G	501	14/15	0.95	0.13	66,94,107,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.