



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:30 PM GMT

PDB ID : 4LSV
Title : Crystal structure of broadly and potently neutralizing antibody 3BNC117 in complex with HIV-1 clade C C1086 gp120
Authors : Zhou, T.; Moquin, S.; Kwong, P.D.
Deposited on : 2013-07-23
Resolution : 3.00 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

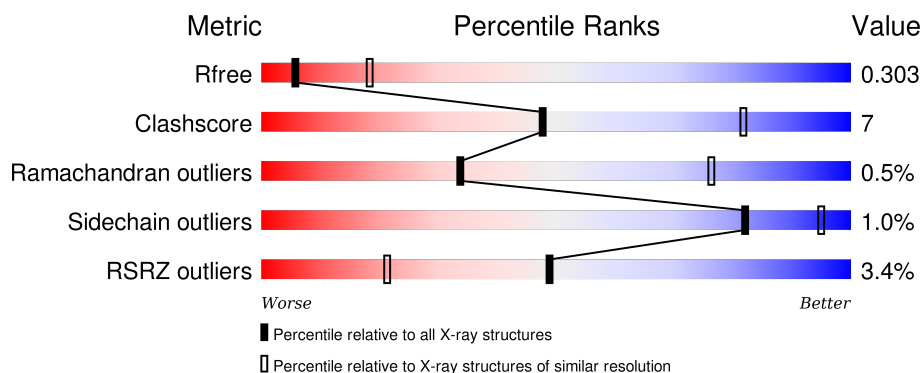
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	G	358	<div> <div>6%</div> <div> <div></div> <div>77%</div> <div>18%</div> <div>• •</div> </div> </div>
2	H	226	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div>• •</div> </div> </div>
3	L	206	<div> <div></div> <div> <div></div> <div>81%</div> <div>17%</div> <div>•</div> </div> </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	505	-	-	X	-
4	NAG	G	506	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6232 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	343	Total	C	N	O	S	0	0	0
			2687	1681	468	518	20			

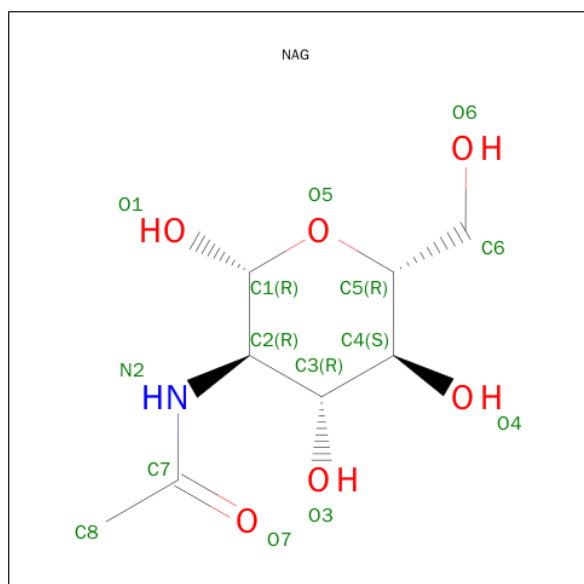
- Molecule 2 is a protein called HEAVY CHAIN OF ANTIBODY 3BNC117.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	221	Total	C	N	O	S	0	0	0
			1710	1084	295	326	5			

- Molecule 3 is a protein called LIGHT CHAIN OF ANTIBODY 3BNC117.

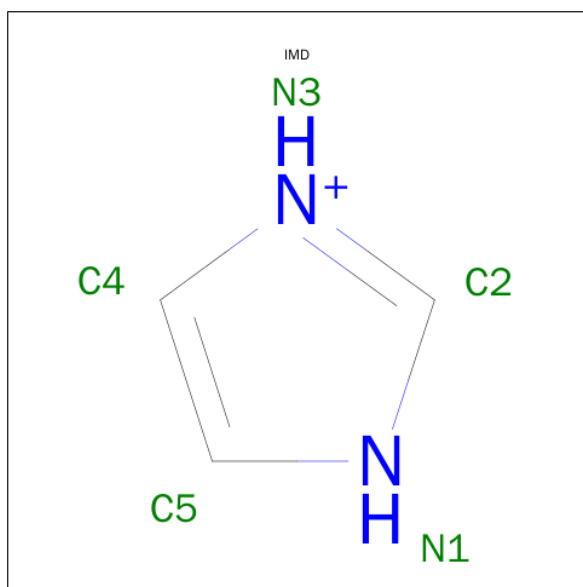
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	203	Total	C	N	O	S	0	0	0
			1599	1002	276	316	5			

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (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		
4	L	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is IMIDAZOLE (three-letter code: IMD) (formula: C₃H₅N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	L	1	Total	C	N	0	0
			5	3	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	G	30	Total 30	O 30	0	0
6	H	24	Total 24	O 24	0	0
6	L	23	Total 23	O 23	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	191.62Å 191.62Å 103.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.32 – 3.00 47.32 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.32-3.00) 99.0 (47.32-2.85)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.30 (at 2.86Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: DEV_998)	Depositor
R, R_{free}	0.260 , 0.299 0.263 , 0.303	Depositor DCC
R_{free} test set	998 reflections (5.12%)	DCC
Wilson B-factor (Å ²)	51.8	Xtriage
Anisotropy	0.984	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 70.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 22631 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6232	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.39% 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: NAG, IMD

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	0.21	0/2744	0.39	0/3722
2	H	0.21	0/1760	0.38	0/2404
3	L	0.21	0/1633	0.38	0/2217
All	All	0.21	0/6137	0.38	0/8343

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	2687	0	2593	53	0
2	H	1710	0	1643	15	0
3	L	1599	0	1561	21	0
4	G	140	0	130	10	1
4	L	14	0	13	1	0
5	L	5	0	5	1	0
6	G	30	0	0	2	0
6	H	24	0	0	0	0
6	L	23	0	0	0	0
All	All	6232	0	5945	89	1

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 7.

All (89) 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:G:289:ASN:CG	4:G:505:NAG:H82	1.56	1.25
1:G:460:GLN:OE1	1:G:460:GLN:HA	1.55	1.00
4:G:505:NAG:O6	6:G:625:HOH:O	1.93	0.87
1:G:290:GLU:HB2	4:G:505:NAG:H83	1.60	0.80
1:G:463:ASN:O	1:G:465:THR:CG2	2.30	0.79
1:G:463:ASN:O	1:G:465:THR:HG22	1.83	0.79
1:G:289:ASN:CG	4:G:505:NAG:C8	2.47	0.76
3:L:33:LEU:HD11	3:L:88:CYS:HB2	1.67	0.76
1:G:289:ASN:ND2	4:G:505:NAG:H82	2.02	0.74
1:G:460:GLN:OE1	1:G:460:GLN:CA	2.36	0.73
1:G:289:ASN:CB	4:G:505:NAG:H82	2.19	0.71
1:G:229:ASN:HB2	1:G:241:ASN:HB3	1.75	0.67
1:G:55:ALA:HB3	1:G:216:HIS:HB2	1.78	0.65
1:G:270:ILE:HB	1:G:348:GLU:HG3	1.77	0.65
2:H:87:THR:HG23	2:H:110:THR:HA	1.79	0.64
2:H:193:THR:HG23	2:H:210:LYS:HE2	1.79	0.64
1:G:113:ASP:HB2	1:G:428:GLN:HE22	1.63	0.64
1:G:254:VAL:HG21	1:G:262:ASN:HB2	1.78	0.64
2:H:197:ASN:ND2	2:H:208:ASP:OD2	2.36	0.59
1:G:289:ASN:OD1	4:G:505:NAG:H82	2.00	0.59
1:G:463:ASN:O	1:G:465:THR:HG23	2.03	0.58
3:L:123:GLU:HA	3:L:126:LYS:HD3	1.86	0.58
1:G:460:GLN:C	1:G:463:ASN:N	2.57	0.58
2:H:151:THR:O	2:H:199:ASN:ND2	2.35	0.58
3:L:21:ILE:HD11	3:L:73:LEU:HD23	1.85	0.57
1:G:290:GLU:HB2	4:G:505:NAG:C8	2.31	0.57
1:G:294:ILE:HG12	1:G:333:ILE:HG12	1.87	0.57
2:H:45:LEU:HB2	3:L:98:VAL:HG11	1.87	0.56
3:L:6:GLN:HB3	3:L:99:VAL:HG11	1.88	0.55
1:G:362:GLU:OE1	1:G:469:ARG:NH2	2.40	0.55
1:G:463:ASN:OD1	1:G:465:THR:CG2	2.55	0.55
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.87	0.55
2:H:7:SER:O	2:H:107:THR:OG1	2.26	0.54
1:G:385:CYS:HA	1:G:418:CYS:HA	1.90	0.53
1:G:254:VAL:HG11	1:G:261:LEU:HB2	1.90	0.53
3:L:187:GLU:HA	3:L:211:ARG:HH12	1.75	0.51
1:G:423:ILE:HG13	1:G:434:ILE:HD12	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:420:ILE:HG21	1:G:438:PRO:HG3	1.91	0.51
1:G:64:GLU:OE2	1:G:67:ASN:ND2	2.45	0.50
3:L:116:PHE:HB2	3:L:135:LEU:HB3	1.92	0.50
3:L:136:LEU:HB2	3:L:175:LEU:HB3	1.93	0.50
1:G:92:ASN:O	1:G:487:LYS:NZ	2.36	0.50
1:G:109:ILE:HG12	1:G:428:GLN:HG3	1.93	0.50
2:H:71(C):SER:OG	2:H:71(D):TRP:N	2.44	0.49
1:G:463:ASN:OD1	1:G:465:THR:HG23	2.12	0.49
1:G:289:ASN:OD1	1:G:290:GLU:N	2.46	0.49
1:G:257:THR:O	1:G:259:LEU:N	2.41	0.48
1:G:343:GLN:HG2	4:G:506:NAG:H62	1.95	0.48
1:G:280:ASN:HD22	1:G:458:GLY:HA3	1.78	0.48
3:L:145:LYS:HB3	3:L:197:THR:HB	1.96	0.48
2:H:155:ASN:HA	2:H:195:ILE:HG23	1.97	0.47
2:H:27:TYR:CE2	2:H:29:ILE:HA	2.49	0.47
1:G:464:ASP:OD1	1:G:464:ASP:N	2.48	0.47
2:H:144:ASP:HA	2:H:175:LEU:HB3	1.96	0.46
1:G:258:GLN:NE2	1:G:372:THR:O	2.47	0.46
3:L:37:GLN:OE1	5:L:302:IMD:H2	2.16	0.45
1:G:343:GLN:NE2	1:G:398:GLY:O	2.34	0.45
1:G:289:ASN:OD1	4:G:505:NAG:C8	2.60	0.45
3:L:108:ARG:HG2	3:L:109:THR:H	1.80	0.45
3:L:113:PRO:HD3	3:L:198:HIS:ND1	2.31	0.45
1:G:82:GLN:NE2	6:G:618:HOH:O	2.48	0.45
1:G:381:GLU:HG3	1:G:443:ILE:HD13	1.99	0.45
2:H:159:LEU:HD21	2:H:182:VAL:HG11	1.99	0.44
3:L:61:ARG:NH2	3:L:81:GLU:OE2	2.45	0.44
1:G:273:ARG:HB2	1:G:285:ILE:HB	1.99	0.44
1:G:389:ASP:HB3	1:G:407:GLY:O	2.17	0.44
3:L:65:ARG:NE	4:L:301:NAG:O7	2.51	0.44
2:H:146:PHE:HA	2:H:147:PRO:HA	1.76	0.43
1:G:117:LYS:HE3	1:G:117:LYS:HB2	1.88	0.42
2:H:136:ALA:O	2:H:184:VAL:N	2.44	0.42
1:G:293:ASN:OD1	1:G:446:ASN:ND2	2.50	0.42
2:H:201:LYS:N	2:H:202:PRO:HD2	2.35	0.42
1:G:216:HIS:ND1	1:G:248:THR:O	2.45	0.42
1:G:286:VAL:HB	1:G:452:LEU:HB2	2.02	0.42
3:L:198:HIS:CD2	3:L:199:GLN:H	2.38	0.41
2:H:82:LEU:HA	2:H:82:LEU:HD12	1.93	0.41
3:L:99:VAL:HG12	3:L:101:GLY:H	1.85	0.41
3:L:35:TRP:CE2	3:L:73:LEU:HB2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:61:ARG:HB2	3:L:76:ASN:O	2.20	0.41
1:G:406:THR:HB	1:G:407:GLY:H	1.59	0.41
1:G:219:ALA:HA	1:G:220:PRO:HD3	1.88	0.41
3:L:150:VAL:HB	3:L:155:GLN:HE21	1.85	0.41
1:G:403:TYR:CE2	1:G:405:HIS:HB2	2.56	0.41
1:G:298:ARG:HD3	1:G:443:ILE:HD12	2.04	0.40
1:G:300:ASN:HD22	1:G:441:GLY:HA2	1.86	0.40
1:G:123:THR:OG1	1:G:199:SER:O	2.33	0.40
3:L:17:ASP:OD1	3:L:18:THR:N	2.55	0.40
1:G:207:LYS:HD3	1:G:439:ILE:HG22	2.04	0.40
1:G:390:LEU:HD22	1:G:416:LEU:HD11	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:502:NAG:O4	4:G:502:NAG:O4[7_554]	2.16	0.04

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	G	337/358 (94%)	307 (91%)	28 (8%)	2 (1%)	30	72
2	H	217/226 (96%)	202 (93%)	14 (6%)	1 (0%)	34	76
3	L	201/206 (98%)	190 (94%)	10 (5%)	1 (0%)	34	76
All	All	755/790 (96%)	699 (93%)	52 (7%)	4 (0%)	34	76

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	116	LEU

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Mol	Chain	Res	Type
1	G	406	THR
2	H	82(A)	LYS
3	L	138	ASN

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	G	302/312 (97%)	299 (99%)	3 (1%)	82	95
2	H	190/193 (98%)	187 (98%)	3 (2%)	70	92
3	L	181/183 (99%)	180 (99%)	1 (1%)	90	97
All	All	673/688 (98%)	666 (99%)	7 (1%)	82	95

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

Mol	Chain	Res	Type
1	G	460	GLN
1	G	464	ASP
1	G	465	THR
2	H	71	ARG
2	H	71(D)	TRP
2	H	199	ASN
3	L	96	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	428	GLN
3	L	147	GLN
3	L	155	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

12 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	501	1	14,14,15	0.48	0	15,19,21	0.82	1 (6%)
4	NAG	G	502	1	14,14,15	0.47	0	15,19,21	0.89	0
4	NAG	G	503	1	14,14,15	0.47	0	15,19,21	0.63	0
4	NAG	G	504	1	14,14,15	0.46	0	15,19,21	0.81	0
4	NAG	G	505	1	14,14,15	0.28	0	15,19,21	0.55	0
4	NAG	G	506	1	14,14,15	0.50	0	15,19,21	0.88	1 (6%)
4	NAG	G	507	1	14,14,15	0.50	0	15,19,21	0.68	0
4	NAG	G	508	1	14,14,15	0.45	0	15,19,21	0.84	0
4	NAG	G	509	1	14,14,15	0.49	0	15,19,21	0.64	0
4	NAG	G	510	1	14,14,15	0.47	0	15,19,21	0.74	0
4	NAG	L	301	3	14,14,15	0.51	0	15,19,21	0.65	0
5	IMD	L	302	-	3,5,5	0.52	0	4,5,5	0.58	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	501	1	-	0/6/23/26	0/1/1/1
4	NAG	G	502	1	-	0/6/23/26	0/1/1/1
4	NAG	G	503	1	-	0/6/23/26	0/1/1/1
4	NAG	G	504	1	-	0/6/23/26	0/1/1/1
4	NAG	G	505	1	-	1/6/23/26	0/1/1/1
4	NAG	G	506	1	-	0/6/23/26	0/1/1/1
4	NAG	G	507	1	-	0/6/23/26	0/1/1/1
4	NAG	G	508	1	-	0/6/23/26	0/1/1/1
4	NAG	G	509	1	-	0/6/23/26	0/1/1/1
4	NAG	G	510	1	-	0/6/23/26	0/1/1/1
4	NAG	L	301	3	-	0/6/23/26	0/1/1/1
5	IMD	L	302	-	-	0/0/0/0	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	501	NAG	C1-O5-C5	2.06	114.87	112.25
4	G	506	NAG	C1-O5-C5	2.24	115.09	112.25

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	505	NAG	O7-C7-N2-C2

There are no ring outliers.

5 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	502	NAG	0	1
4	G	505	NAG	9	0
4	G	506	NAG	1	0
4	L	301	NAG	1	0
5	L	302	IMD	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	343/358 (95%)	0.60	23 (6%) 21 7	71, 113, 144, 181	0
2	H	221/226 (97%)	-0.07	3 (1%) 78 51	41, 68, 126, 178	0
3	L	203/206 (98%)	-0.14	0 100 100	33, 62, 103, 122	0
All	All	767/790 (97%)	0.21	26 (3%) 49 21	33, 91, 139, 181	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	127	SER	5.4
1	G	200	ALA	5.1
1	G	88	ASN	4.4
1	G	299	PRO	3.9
1	G	383	PHE	3.2
1	G	44	VAL	3.1
1	G	433	ALA	3.1
1	G	300	ASN	2.9
1	G	82	GLN	2.7
2	H	128	SER	2.7
1	G	436	ALA	2.7
1	G	198	GLY	2.6
1	G	353	PHE	2.6
1	G	410	SER	2.5
1	G	210	PHE	2.5
1	G	492	LYS	2.5
1	G	64	GLU	2.4
1	G	56	SER	2.4
1	G	201	ILE	2.4
1	G	224	ALA	2.4
1	G	66	HIS	2.3
1	G	226	LEU	2.3
1	G	122	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	130	SER	2.3
1	G	228	CYS	2.3
1	G	71	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](#)

There are no carbohydrates 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. 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
4	NAG	G	506	14/15	0.67	0.33	3.12	84,88,89,89	0
4	NAG	G	505	14/15	0.74	0.36	0.80	80,81,87,89	0
4	NAG	G	502	14/15	0.74	0.31	-0.10	81,83,91,92	0
4	NAG	L	301	14/15	0.91	0.23	-0.27	46,52,66,69	0
4	NAG	G	503	14/15	0.82	0.23	-0.81	61,69,72,75	0
4	NAG	G	504	14/15	0.95	0.15	-0.83	37,46,55,56	0
4	NAG	G	508	14/15	0.89	0.17	-0.84	61,63,64,64	0
4	NAG	G	507	14/15	0.87	0.20	-1.01	52,56,61,61	0
4	NAG	G	501	14/15	0.85	0.20	-1.66	57,65,70,71	0
5	IMD	L	302	5/5	0.96	0.14	-1.67	63,63,65,66	0
4	NAG	G	509	14/15	0.72	0.33	-	83,91,94,94	0
4	NAG	G	510	14/15	0.67	0.31	-	81,88,91,92	0

6.5 Other polymers [i](#)

There are no such residues in this entry.