



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:11 PM GMT

PDB ID : 4R4H
Title : Crystal structure of non-neutralizing, A32-like antibody 2.2c in complex with HIV-1 Env gp120
Authors : McLellan, J.; Acharya, P.; Huang, C.-C.; Kwong, P.D.
Deposited on : 2014-08-19
Resolution : 4.28 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

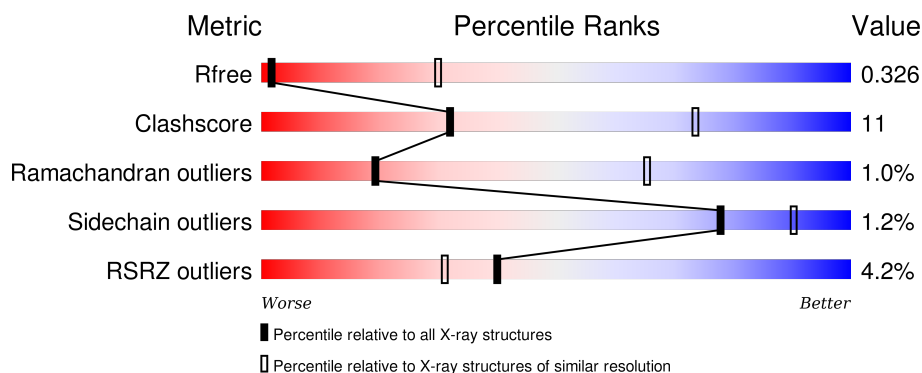
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.28 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	1043 (4.92-3.60)
Clashscore	102246	1145 (4.92-3.60)
Ramachandran outliers	100387	1088 (4.92-3.60)
Sidechain outliers	100360	1072 (4.92-3.60)
RSRZ outliers	91569	1047 (4.92-3.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	433	<div> <div>11%</div> <div>59%</div> <div>18%</div> <div>22%</div> </div>
2	B	178	<div> <div>11%</div> <div>83%</div> <div>16%</div> </div>
3	L	210	<div> <div>86%</div> <div>14%</div> </div>
4	H	220	<div> <div>7%</div> <div>75%</div> <div>24%</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
5	NAG	A	601	-	-	-	X
5	NAG	A	603	-	-	-	X
5	NAG	A	604	-	-	X	-
5	NAG	A	605	-	-	-	X
5	NAG	A	606	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7415 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 HIV-1 Env gp120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2629	1644	461	503	21			

- Molecule 2 is a protein called T-cell surface glycoprotein CD4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	178	Total	C	N	O	S	0	3	0
			1391	868	243	276	4			

- Molecule 3 is a protein called Antibody 2.2c, Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	210	Total	C	N	O	S	0	0	0
			1612	1012	272	323	5			

- Molecule 4 is a protein called Antibody 2.2c, Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	220	Total	C	N	O	S	0	0	0
			1671	1064	283	319	5			

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).

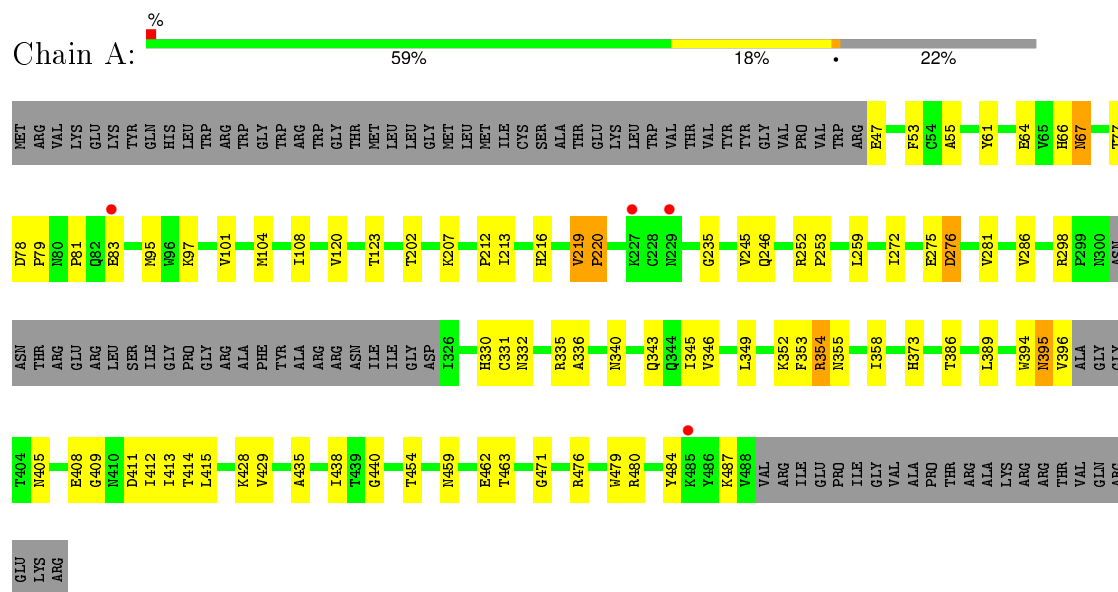


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	H	1	Total	C	N	O	0	0
			14	8	1	5		

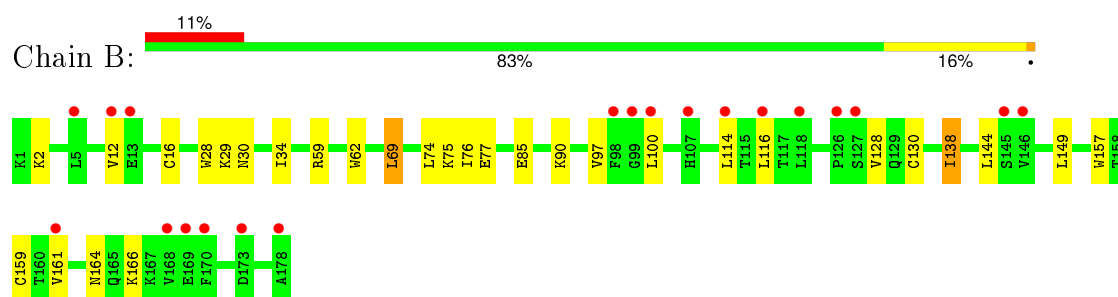
3 Residue-property plots [i](#)

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.

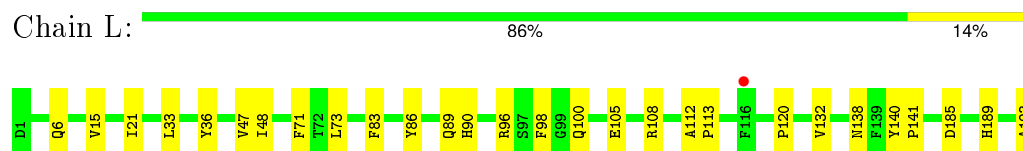
• Molecule 1: HIV-1 Env gp120



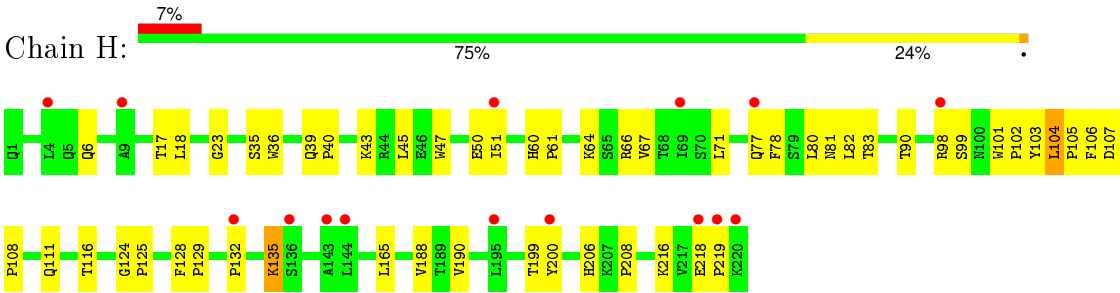
• Molecule 2: T-cell surface glycoprotein CD4



• Molecule 3: Antibody 2.2c, Light chain



• Molecule 4: Antibody 2.2c, Heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.36Å 69.43Å 328.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.73 – 4.28 47.73 – 4.28	Depositor EDS
% Data completeness (in resolution range)	76.8 (47.73-4.28) 68.0 (47.73-4.28)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 4.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.309 , 0.320 0.311 , 0.326	Depositor DCC
R_{free} test set	382 reflections (5.44%)	DCC
Wilson B-factor (Å ²)	119.2	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 133.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Outliers	0 of 7945 reflections	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	7415	wwPDB-VP
Average B, all atoms (Å ²)	166.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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

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	0.29	0/2681	0.45	1/3642 (0.0%)
2	B	0.22	0/1424	0.38	0/1921
3	L	0.22	0/1646	0.37	0/2231
4	H	0.23	0/1720	0.39	0/2352
All	All	0.25	0/7471	0.41	1/10146 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	276	ASP	CB-CG-OD2	5.20	122.98	118.30

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	2629	0	2555	83	0
2	B	1391	0	1417	26	0
3	L	1612	0	1583	18	0
4	H	1671	0	1648	40	0
5	A	98	0	89	20	0
5	H	14	0	13	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7415	0	7305	155	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 11.

All (155) 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)
4:H:81:ASN:HD21	5:H:301:NAG:C1	1.32	1.43
4:H:81:ASN:ND2	5:H:301:NAG:O5	1.63	1.30
1:A:332:ASN:HD21	5:A:604:NAG:H2	1.07	1.10
1:A:332:ASN:OD1	5:A:604:NAG:C1	2.06	1.04
1:A:340:ASN:HA	1:A:343:GLN:HG3	1.40	1.03
1:A:330:HIS:O	5:A:604:NAG:H81	1.58	1.01
1:A:332:ASN:ND2	5:A:604:NAG:H2	1.76	1.00
1:A:336:ALA:HA	5:A:605:NAG:O7	1.67	0.95
1:A:275:GLU:O	1:A:276:ASP:OD1	1.86	0.91
1:A:330:HIS:O	5:A:604:NAG:C8	2.18	0.90
1:A:281:VAL:CG2	2:B:29:LYS:HE2	2.03	0.89
1:A:332:ASN:OD1	5:A:604:NAG:C2	2.22	0.87
1:A:330:HIS:C	5:A:604:NAG:C8	2.42	0.87
1:A:281:VAL:HG21	2:B:29:LYS:CE	2.04	0.86
1:A:330:HIS:C	5:A:604:NAG:H82	1.95	0.86
1:A:335:ARG:NE	1:A:409:GLY:O	2.09	0.85
4:H:39:GLN:HB2	4:H:45:LEU:HD23	1.65	0.79
2:B:100:LEU:HD21	2:B:116:LEU:HB3	1.64	0.78
1:A:79:PRO:HG2	1:A:246:GLN:HG2	1.67	0.76
1:A:281:VAL:HG21	2:B:29:LYS:HE2	1.68	0.74
1:A:332:ASN:OD1	5:A:604:NAG:N2	2.21	0.73
1:A:298:ARG:NH2	1:A:440:GLY:O	2.18	0.72
1:A:281:VAL:CG2	2:B:29:LYS:CE	2.65	0.72
4:H:18:LEU:HB3	4:H:82:LEU:HB3	1.73	0.69
4:H:17:THR:HG21	5:H:301:NAG:H62	1.74	0.69
1:A:77:THR:HB	1:A:78:ASP:HA	1.73	0.69
4:H:6:GLN:H	4:H:111:GLN:HE22	1.41	0.69
1:A:120:VAL:HG22	1:A:202:THR:HG22	1.74	0.68
1:A:281:VAL:HG12	1:A:281:VAL:O	1.93	0.66
3:L:89:GLN:HE22	4:H:106:PHE:HE2	1.45	0.64
1:A:335:ARG:HD3	1:A:408:GLU:O	1.98	0.64
1:A:340:ASN:CA	1:A:343:GLN:HG3	2.22	0.64
4:H:71:LEU:HD23	4:H:78:PHE:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:85:GLU:HG2	2:B:90:LYS:HG2	1.81	0.62
1:A:343:GLN:HG2	1:A:394:TRP:HH2	1.65	0.61
4:H:17:THR:HG21	5:H:301:NAG:C6	2.31	0.61
1:A:343:GLN:HG2	1:A:394:TRP:CH2	2.36	0.61
1:A:438:ILE:HG22	1:A:440:GLY:H	1.66	0.60
1:A:108:ILE:HD12	1:A:253:PRO:HB3	1.84	0.60
1:A:281:VAL:HG23	2:B:29:LYS:HE2	1.81	0.59
1:A:55:ALA:HB3	1:A:216:HIS:HB2	1.84	0.59
1:A:272:ILE:HG22	1:A:286:VAL:HG22	1.83	0.59
1:A:281:VAL:HG21	2:B:29:LYS:HE3	1.83	0.59
1:A:83:GLU:HG2	1:A:245:VAL:HG12	1.85	0.58
4:H:35:SER:HB2	4:H:50:GLU:HB3	1.85	0.58
3:L:120:PRO:HD3	3:L:132:VAL:HG22	1.85	0.57
5:A:601:NAG:H62	5:A:607:NAG:H81	1.87	0.57
3:L:47:VAL:HG12	3:L:48:ILE:HG13	1.85	0.57
4:H:51:ILE:HD13	4:H:71:LEU:HG	1.86	0.57
1:A:354:ARG:HG3	1:A:355:ASN:H	1.70	0.57
1:A:394:TRP:O	1:A:395:ASN:HB2	2.05	0.57
1:A:67:ASN:ND2	1:A:213:ILE:HG13	2.21	0.56
1:A:340:ASN:HA	1:A:343:GLN:CG	2.26	0.56
1:A:332:ASN:CG	5:A:604:NAG:C2	2.74	0.56
1:A:332:ASN:CG	5:A:604:NAG:H2	2.27	0.55
1:A:346:VAL:HG21	1:A:394:TRP:CD1	2.41	0.55
1:A:331:CYS:N	5:A:604:NAG:H82	2.22	0.55
1:A:332:ASN:ND2	5:A:604:NAG:C2	2.61	0.55
3:L:21:ILE:HD11	3:L:73:LEU:HD23	1.89	0.54
2:B:59:ARG:HA	2:B:62:TRP:CD1	2.42	0.54
1:A:335:ARG:CZ	1:A:413:ILE:HG12	2.38	0.54
1:A:77:THR:HB	1:A:78:ASP:CA	2.37	0.54
2:B:138:ILE:HD12	2:B:144:LEU:HB3	1.90	0.53
2:B:128:VAL:HG22	2:B:161:VAL:HG22	1.90	0.53
4:H:23:GLY:HA2	4:H:77:GLN:HG2	1.90	0.53
4:H:199:THR:HG23	4:H:216:LYS:HE3	1.91	0.53
3:L:90:HIS:O	3:L:96:ARG:HA	2.09	0.53
1:A:462:GLU:HG2	1:A:463:THR:H	1.74	0.52
1:A:414:THR:OG1	5:A:604:NAG:C1	2.57	0.52
1:A:476:ARG:HA	1:A:479:TRP:CD1	2.44	0.51
1:A:395:ASN:O	1:A:396:VAL:HB	2.09	0.51
1:A:281:VAL:CG1	1:A:281:VAL:O	2.59	0.51
4:H:101:TRP:CD1	4:H:102:PRO:HA	2.45	0.51
1:A:101:VAL:HG13	1:A:479:TRP:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:ASN:HD21	1:A:213:ILE:HG13	1.75	0.50
1:A:335:ARG:HG3	1:A:335:ARG:HH11	1.76	0.49
1:A:53:PHE:CD2	4:H:101:TRP:HZ2	2.30	0.49
2:B:130:CYS:HA	2:B:159:CYS:HA	1.95	0.49
3:L:193:ALA:HB2	3:L:208:SER:HB3	1.95	0.49
1:A:123:THR:HG21	1:A:428:LYS:HE3	1.93	0.49
3:L:6:GLN:HB2	3:L:100:GLN:HE22	1.78	0.49
1:A:332:ASN:CG	5:A:604:NAG:C1	2.80	0.49
1:A:332:ASN:OD1	5:A:604:NAG:C7	2.60	0.49
4:H:190:VAL:HG11	4:H:200:TYR:CE1	2.48	0.49
1:A:207:LYS:HE2	1:A:435:ALA:HB3	1.94	0.48
1:A:354:ARG:HG3	1:A:355:ASN:N	2.29	0.48
4:H:165:LEU:HD21	4:H:188:VAL:HG21	1.96	0.48
1:A:346:VAL:HG13	1:A:358:ILE:HB	1.95	0.48
1:A:66:HIS:CD2	1:A:212:PRO:HA	2.49	0.47
1:A:335:ARG:HG3	1:A:335:ARG:NH1	2.30	0.47
3:L:89:GLN:HB2	3:L:98:PHE:CE2	2.49	0.47
1:A:346:VAL:HG21	1:A:394:TRP:CE2	2.50	0.47
2:B:76:ILE:HA	2:B:97:VAL:HG23	1.96	0.47
4:H:206:HIS:CE1	4:H:208:PRO:HB2	2.50	0.47
2:B:12:VAL:HG12	2:B:74:LEU:HD21	1.96	0.47
4:H:135:LYS:H	4:H:135:LYS:HD3	1.80	0.46
2:B:30:ASN:HD21	2:B:34:ILE:HD12	1.81	0.46
1:A:104:MET:O	1:A:108:ILE:HG12	2.16	0.46
4:H:132:PRO:HG2	4:H:219:PRO:HG3	1.98	0.46
4:H:40:PRO:HB2	4:H:43:LYS:HD2	1.98	0.46
2:B:16:CYS:HB2	2:B:28:TRP:CZ2	2.51	0.45
3:L:185:ASP:O	3:L:189:HIS:HD2	1.99	0.45
1:A:353:PHE:O	1:A:355:ASN:N	2.50	0.45
3:L:33:LEU:HD22	3:L:71:PHE:CG	2.52	0.45
1:A:77:THR:HG22	4:H:101:TRP:HB3	1.99	0.45
1:A:272:ILE:HD11	1:A:352:LYS:HG3	1.98	0.45
4:H:40:PRO:HB2	4:H:43:LYS:HB2	1.98	0.45
3:L:98:PHE:CG	4:H:45:LEU:HB2	2.52	0.45
2:B:114:LEU:HB3	2:B:149:LEU:HD11	1.98	0.45
4:H:128:PHE:HA	4:H:129:PRO:HD3	1.85	0.45
2:B:130:CYS:SG	2:B:157:TRP:HB3	2.57	0.44
4:H:90:THR:HG23	4:H:116:THR:HA	2.00	0.44
4:H:99:SER:HB2	4:H:103:TYR:HD2	1.82	0.44
1:A:454:THR:HG23	1:A:471:GLY:HA3	2.00	0.44
4:H:101:TRP:CG	4:H:102:PRO:HA	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:601:NAG:H62	5:A:607:NAG:C8	2.47	0.44
1:A:429:VAL:HG11	2:B:59:ARG:O	2.18	0.44
2:B:28:TRP:CD2	2:B:69:LEU:HD12	2.52	0.44
1:A:95:MET:SD	1:A:235:GLY:HA3	2.57	0.44
2:B:77:GLU:CD	2:B:77:GLU:H	2.21	0.44
1:A:64:GLU:OE2	1:A:66:HIS:HB2	2.18	0.43
3:L:15:VAL:HG22	3:L:83:PHE:HE2	1.83	0.43
1:A:252:ARG:HA	1:A:253:PRO:HD3	1.84	0.43
4:H:216:LYS:HE2	4:H:218:GLU:HG2	1.99	0.43
1:A:386:THR:HG22	1:A:415:LEU:HD13	1.99	0.43
1:A:219:VAL:HA	1:A:220:PRO:HD3	1.87	0.43
2:B:12:VAL:HB	2:B:74:LEU:HD11	1.99	0.43
3:L:98:PHE:CD1	4:H:45:LEU:HB2	2.54	0.43
4:H:47:TRP:HZ2	4:H:50:GLU:HG2	1.84	0.42
1:A:389:LEU:HD11	1:A:415:LEU:HD11	2.00	0.42
3:L:112:ALA:HA	3:L:113:PRO:HD3	1.87	0.42
2:B:28:TRP:CE3	2:B:69:LEU:HD12	2.54	0.42
1:A:480:ARG:O	1:A:484:TYR:HB3	2.20	0.42
3:L:36:TYR:O	3:L:86:TYR:HA	2.20	0.42
1:A:61:TYR:CD1	4:H:64:LYS:HD2	2.55	0.42
4:H:104:LEU:H	4:H:105:PRO:CD	2.32	0.42
1:A:259:LEU:HB2	1:A:373:HIS:CE1	2.54	0.42
2:B:75:LYS:O	2:B:97:VAL:HG21	2.20	0.42
1:A:386:THR:HB	1:A:389:LEU:HD12	2.02	0.42
1:A:281:VAL:HG23	2:B:29:LYS:CE	2.46	0.41
1:A:346:VAL:HG21	1:A:394:TRP:NE1	2.35	0.41
4:H:104:LEU:N	4:H:105:PRO:CD	2.83	0.41
3:L:15:VAL:HG22	3:L:83:PHE:CE2	2.56	0.41
4:H:67:VAL:HA	4:H:81:ASN:O	2.20	0.41
4:H:107:ASP:OD1	4:H:108:PRO:HA	2.21	0.41
3:L:6:GLN:HB2	3:L:100:GLN:NE2	2.36	0.41
3:L:140:TYR:CG	3:L:141:PRO:HA	2.56	0.41
4:H:36:TRP:CE2	4:H:80:LEU:HB2	2.56	0.41
4:H:66:ARG:HD2	4:H:83:THR:O	2.21	0.41
4:H:124:GLY:HA2	4:H:125:PRO:HD3	1.84	0.41
2:B:164:ASN:C	2:B:166:LYS:H	2.25	0.40
1:A:412:ILE:HD11	5:A:604:NAG:O6	2.21	0.40
1:A:345:ILE:O	1:A:349:LEU:HG	2.21	0.40
1:A:47:GLU:HG2	1:A:487:LYS:HE3	2.03	0.40
4:H:60:HIS:CD2	4:H:61:PRO:HD2	2.57	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	A	331/433 (76%)	288 (87%)	36 (11%)	7 (2%)	9	52
2	B	179/178 (101%)	170 (95%)	9 (5%)	0	100	100
3	L	208/210 (99%)	196 (94%)	11 (5%)	1 (0%)	34	77
4	H	218/220 (99%)	201 (92%)	16 (7%)	1 (0%)	34	77
All	All	936/1041 (90%)	855 (91%)	72 (8%)	9 (1%)	19	65

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	354	ARG
1	A	395	ASN
1	A	405	ASN
3	L	138	ASN
1	A	97	LYS
1	A	411	ASP
1	A	220	PRO
4	H	104	LEU
1	A	81	PRO

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	A	298/379 (79%)	295 (99%)	3 (1%)	82	92
2	B	164/161 (102%)	161 (98%)	3 (2%)	66	88

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	L	183/183 (100%)	181 (99%)	2 (1%)	80	91
4	H	191/191 (100%)	189 (99%)	2 (1%)	82	92
All	All	836/914 (92%)	826 (99%)	10 (1%)	78	90

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	A	219	VAL
1	A	459	ASN
2	B	2	LYS
2	B	69	LEU
2	B	138	ILE
3	L	105	GLU
3	L	108	ARG
4	H	98	ARG
4	H	135	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	A	82	GLN
1	A	330	HIS
4	H	81	ASN

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

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$
5	NAG	A	601	1	14,14,15	0.56	0	15,19,21	0.72	1 (6%)
5	NAG	A	602	1	14,14,15	0.56	0	15,19,21	3.70	2 (13%)
5	NAG	A	603	1	14,14,15	0.52	0	15,19,21	0.40	0
5	NAG	A	604	-	14,14,15	1.06	1 (7%)	15,19,21	0.64	0
5	NAG	A	605	1	14,14,15	1.80	2 (14%)	15,19,21	2.36	3 (20%)
5	NAG	A	606	1	14,14,15	0.44	0	15,19,21	0.75	0
5	NAG	A	607	1	14,14,15	0.67	0	15,19,21	1.19	2 (13%)
5	NAG	H	301	4	14,14,15	0.42	0	15,19,21	0.96	1 (6%)

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
5	NAG	A	601	1	-	0/6/23/26	0/1/1/1
5	NAG	A	602	1	-	0/6/23/26	0/1/1/1
5	NAG	A	603	1	-	0/6/23/26	0/1/1/1
5	NAG	A	604	-	-	0/6/23/26	0/1/1/1
5	NAG	A	605	1	-	0/6/23/26	0/1/1/1
5	NAG	A	606	1	-	0/6/23/26	0/1/1/1
5	NAG	A	607	1	-	0/6/23/26	0/1/1/1
5	NAG	H	301	4	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	605	NAG	O5-C1	-5.66	1.34	1.43
5	A	604	NAG	O5-C1	-3.29	1.38	1.43
5	A	605	NAG	C1-C2	-2.92	1.48	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	605	NAG	C4-C3-C2	-3.48	105.82	111.23
5	A	607	NAG	C4-C3-C2	-2.89	106.73	111.23
5	A	602	NAG	C1-O5-C5	-2.03	109.67	112.25
5	H	301	NAG	O4-C4-C3	-2.01	105.81	110.34
5	A	607	NAG	C1-O5-C5	2.09	114.90	112.25
5	A	601	NAG	C1-O5-C5	2.19	115.02	112.25
5	A	605	NAG	O3-C3-C2	2.75	114.57	109.11
5	A	605	NAG	C2-N2-C7	7.41	132.56	123.04
5	A	602	NAG	C2-N2-C7	13.89	140.89	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	601	NAG	2	0
5	A	604	NAG	17	0
5	A	605	NAG	1	0
5	A	607	NAG	2	0
5	H	301	NAG	4	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	A	337/433 (77%)	0.17	4 (1%) 81 73	89, 150, 183, 214	0
2	B	178/178 (100%)	0.58	20 (11%) 7 6	135, 180, 273, 288	0
3	L	210/210 (100%)	0.07	1 (0%) 91 88	138, 171, 213, 232	0
4	H	220/220 (100%)	0.41	15 (6%) 20 14	132, 157, 208, 226	0
All	All	945/1041 (90%)	0.28	40 (4%) 40 31	89, 160, 227, 288	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	118	LEU	4.4
2	B	169	GLU	3.9
4	H	219	PRO	3.8
2	B	100	LEU	3.7
4	H	132	PRO	3.7
4	H	143	ALA	3.7
2	B	98	PHE	3.5
2	B	5	LEU	3.2
4	H	144	LEU	3.0
4	H	220	LYS	3.0
2	B	168	VAL	3.0
2	B	170	PHE	2.9
4	H	136	SER	2.9
4	H	9	ALA	2.8
2	B	12	VAL	2.8
2	B	173	ASP	2.7
4	H	69	ILE	2.7
1	A	227	LYS	2.7
1	A	83	GLU	2.7
2	B	127	SER	2.6
2	B	161	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	145	SER	2.5
4	H	4	LEU	2.5
2	B	116	LEU	2.5
2	B	126	PRO	2.5
2	B	114	LEU	2.4
2	B	178	ALA	2.3
1	A	485	LYS	2.3
4	H	218	GLU	2.3
4	H	195	LEU	2.3
2	B	99	GLY	2.3
4	H	77	GLN	2.2
4	H	200	TYR	2.2
1	A	229	ASN	2.2
4	H	51	ILE	2.2
2	B	146	VAL	2.1
4	H	98	ARG	2.1
2	B	107	HIS	2.1
3	L	116	PHE	2.0
2	B	13	GLU	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
5	NAG	A	605	14/15	0.68	0.36	3.19	175,181,185,187	0
5	NAG	A	603	14/15	0.52	0.60	3.13	154,158,162,163	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	NAG	A	601	14/15	0.75	0.45	1.78	135,138,141,141	0
5	NAG	A	606	14/15	0.71	0.46	1.03	161,164,168,168	0
5	NAG	A	604	14/15	0.86	0.21	-0.89	153,163,170,170	0
5	NAG	H	301	14/15	0.64	0.44	-	166,171,178,179	0
5	NAG	A	607	14/15	0.69	0.45	-	144,151,156,156	0
5	NAG	A	602	14/15	0.81	0.32	-	167,172,178,178	0

6.5 Other polymers [i](#)

There are no such residues in this entry.