



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:45 AM UTC

PDB ID : 4YBL / pdb_00004ybl
Title : Crystal structure of the stabilized inner domain of clade A/E HIV-1 gp120 in complex with the ADCC mediating ANTI-HIV-1 antibody A32
Authors : Tolbert, W.D.; Gohain, N.; Pazgier, M.
Deposited on : 2015-02-18
Resolution : 3.10 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

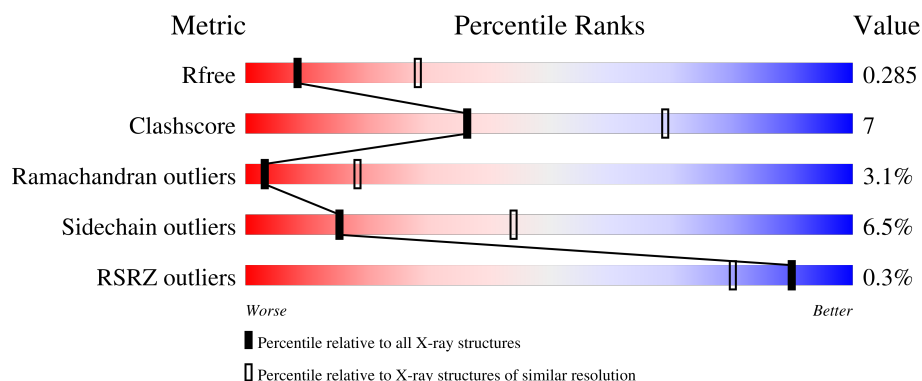
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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 of 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	A	154	
1	G	154	
2	B	224	
2	H	224	
3	C	210	

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Mol	Chain	Length	Quality of chain
3	L	210	90% 9% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8470 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 Stabilized inner domain of clade A/E gp120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	126	Total	C	N	O	S	0	0	0
			1007	633	171	192	11			
1	A	129	Total	C	N	O	S	0	0	0
			1036	652	176	197	11			

- Molecule 2 is a protein called A32 antibody Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	214	Total	C	N	O	S	0	0	0
			1627	1031	279	312	5			
2	B	222	Total	C	N	O	S	0	0	0
			1676	1057	287	327	5			

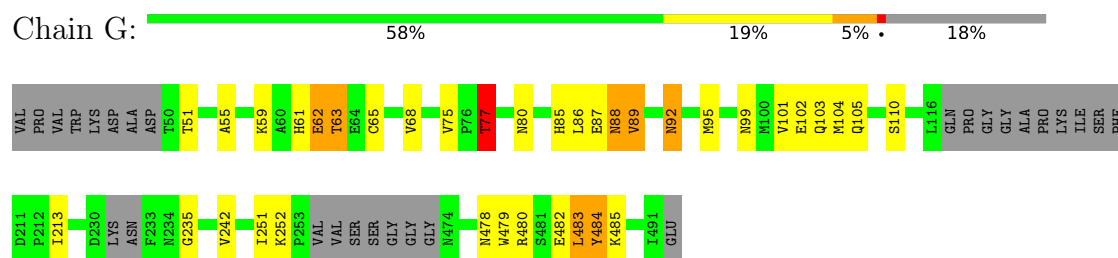
- Molecule 3 is a protein called A32 antibody light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	210	Total	C	N	O	S	0	0	0
			1562	977	261	320	4			
3	C	210	Total	C	N	O	S	0	0	0
			1562	977	261	320	4			

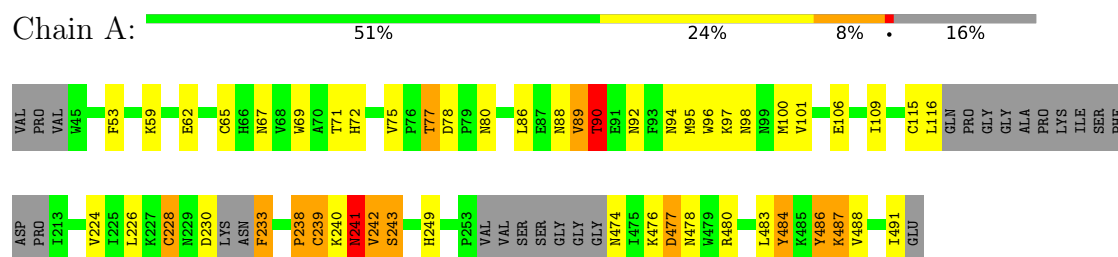
3 Residue-property plots

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.

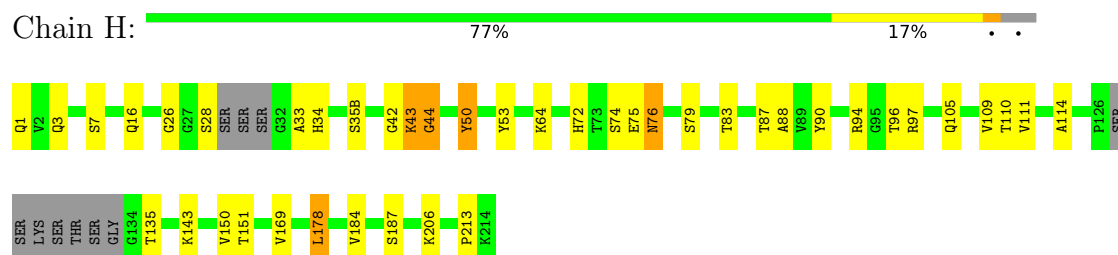
- Molecule 1: Stabilized inner domain of clade A/E gp120



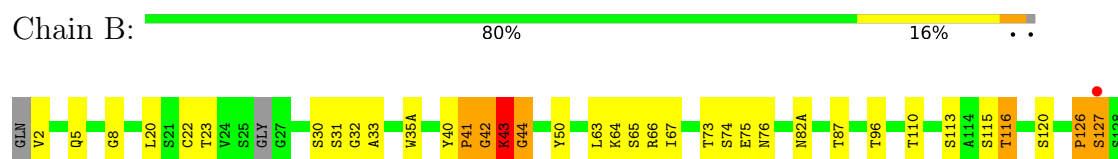
- Molecule 1: Stabilized inner domain of clade A/E gp120



- Molecule 2: A32 antibody Fab heavy chain

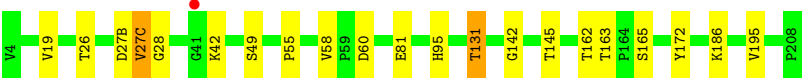
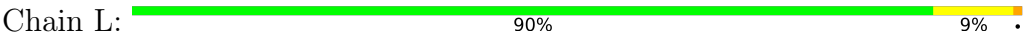


- Molecule 2: A32 antibody Fab heavy chain

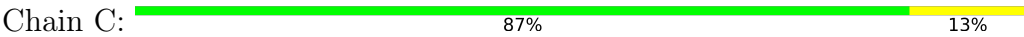




● Molecule 3: A32 antibody light chain



● Molecule 3: A32 antibody light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	75.74Å 208.21Å 73.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.10 50.00 – 3.10	Depositor EDS
% Data completeness (in resolution range)	91.5 (50.00-3.10) 91.5 (50.00-3.10)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 3.06Å)	Xtriage
Refinement program	REFMAC 5.8.0103, PHENIX, CNS	Depositor
R, R_{free}	0.232 , 0.289 0.231 , 0.285	Depositor DCC
R_{free} test set	1021 reflections (4.45%)	wwPDB-VP
Wilson B-factor (Å ²)	82.6	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 53.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.021 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8470	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.22% 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

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	0.85	1/1062 (0.1%)	1.19	8/1445 (0.6%)
1	G	0.70	0/1032	1.04	7/1405 (0.5%)
2	B	0.66	0/1718	0.93	4/2343 (0.2%)
2	H	0.61	0/1668	0.90	1/2275 (0.0%)
3	C	0.67	0/1603	0.88	0/2191
3	L	0.63	0/1603	0.91	3/2191 (0.1%)
All	All	0.68	1/8686 (0.0%)	0.96	23/11850 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
2	B	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	77	THR	CB-CG2	-11.39	1.15	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	90	THR	N-CA-C	9.09	120.65	108.38
1	G	77	THR	CA-CB-OG1	-8.43	96.96	109.60
3	L	27(C)	VAL	N-CA-C	-8.00	101.33	110.05
1	A	77	THR	CA-CB-CG2	-7.64	97.51	110.50
1	G	77	THR	CB-CA-C	7.16	121.34	109.53
2	B	40	TYR	N-CA-C	7.05	125.39	109.81
1	A	477	ASP	N-CA-C	6.98	120.69	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	184	VAL	N-CA-C	6.82	114.78	109.19
1	A	241	ASN	N-CA-C	6.37	119.00	108.49
2	B	43	LYS	N-CA-C	6.34	124.31	110.80
2	B	40	TYR	C-N-CD	-6.21	106.93	120.60
1	G	88	ASN	N-CA-C	6.09	123.78	110.80
3	L	163	THR	CA-C-N	5.73	127.00	119.84
3	L	163	THR	C-N-CA	5.73	127.00	119.84
1	G	63	THR	N-CA-C	5.71	120.24	113.28
1	A	488	VAL	N-CA-C	5.62	116.57	108.48
1	A	77	THR	N-CA-C	5.57	117.88	110.53
1	G	62	GLU	N-CA-C	5.54	118.49	109.24
1	A	88	ASN	N-CA-C	5.37	122.24	110.80
1	G	251	ILE	N-CA-C	5.36	114.47	106.85
1	G	92	ASN	N-CA-C	5.28	117.33	109.25
1	A	228	CYS	CA-CB-SG	-5.24	102.35	114.40
2	B	129	LYS	N-CA-C	-5.12	106.81	113.16

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	233	PHE	Peptide
1	A	484	TYR	Peptide
1	A	89	VAL	Peptide
2	B	35(A)	TRP	Peptide

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	A	1036	0	968	52	0
1	G	1007	0	943	26	0
2	B	1676	0	1648	17	0
2	H	1627	0	1603	17	0
3	C	1562	0	1499	8	1
3	L	1562	0	1499	6	0
All	All	8470	0	8160	119	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 (119) 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:249:HIS:ND1	1:A:486:TYR:OH	2.02	0.92
1:A:230:ASP:CG	1:A:233:PHE:HB2	2.02	0.84
1:A:230:ASP:OD2	1:A:233:PHE:HB2	1.80	0.81
1:A:77:THR:HG22	1:A:78:ASP:N	1.95	0.79
1:A:86:LEU:HD23	1:A:242:VAL:C	2.12	0.74
2:B:63:LEU:HB2	2:B:67:ILE:HD11	1.72	0.71
1:A:90:THR:HG21	1:A:240:LYS:HA	1.72	0.70
1:A:86:LEU:HD23	1:A:243:SER:N	2.06	0.70
1:G:55:ALA:HB1	1:G:77:THR:OG1	1.92	0.70
1:A:480:ARG:HA	1:A:484:TYR:CE2	2.27	0.69
1:A:92:ASN:HA	1:A:238:PRO:HA	1.76	0.68
1:G:75:VAL:HG11	2:H:33:ALA:HB1	1.76	0.67
1:G:101:VAL:HG11	1:G:480:ARG:HE	1.59	0.67
1:A:75:VAL:HG11	2:B:33:ALA:HB1	1.77	0.67
1:A:480:ARG:HD3	1:A:484:TYR:HE2	1.60	0.66
1:G:101:VAL:CG2	1:G:483:LEU:HD22	2.27	0.64
1:A:101:VAL:HG21	1:A:484:TYR:OH	2.00	0.61
3:C:92:THR:O	3:C:93:ASP:O	2.18	0.61
1:A:86:LEU:HB2	1:A:89:VAL:HG22	1.82	0.60
1:A:106:GLU:HA	1:A:109:ILE:HD12	1.82	0.60
1:A:90:THR:CG2	1:A:240:LYS:HA	2.32	0.60
1:A:86:LEU:HD12	1:A:86:LEU:O	2.02	0.60
3:C:115:VAL:O	3:C:204:LYS:HD2	2.01	0.59
1:A:477:ASP:O	1:A:480:ARG:HB3	2.04	0.58
1:G:86:LEU:HD22	1:G:242:VAL:HB	1.87	0.57
1:A:69:TRP:HA	1:A:72:HIS:CE1	2.40	0.57
1:A:90:THR:HG21	1:A:239:CYS:C	2.30	0.57
1:A:477:ASP:OD1	1:A:478:ASN:N	2.38	0.56
1:G:86:LEU:HD11	1:G:89:VAL:HB	1.88	0.56
2:B:66:ARG:NH1	2:B:82(A):ASN:O	2.39	0.55
1:G:104:MET:HG2	1:G:479:TRP:CD1	2.41	0.55
2:H:43:LYS:HB2	2:H:44:GLY:HA3	1.89	0.55
2:B:63:LEU:O	2:B:65:SER:N	2.41	0.54
2:B:41:PRO:O	2:B:42:GLY:C	2.51	0.54
1:A:480:ARG:HD3	1:A:484:TYR:CE2	2.43	0.54
1:A:101:VAL:HG22	1:A:483:LEU:CD1	2.38	0.53
2:H:79:SER:OG	1:A:80:ASN:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:65:CYS:O	1:G:68:VAL:HB	2.11	0.51
2:B:8:GLY:HA3	2:B:20:LEU:HD23	1.94	0.50
1:A:224:VAL:HG12	1:A:491:ILE:HD11	1.93	0.50
1:G:483:LEU:O	1:G:485:LYS:N	2.45	0.50
3:C:93:ASP:O	3:C:94:ILE:C	2.55	0.49
1:A:67:ASN:O	1:A:71:THR:HG23	2.12	0.49
1:A:92:ASN:HA	1:A:238:PRO:CA	2.40	0.49
2:H:88:ALA:HB3	2:H:90:TYR:CE1	2.48	0.49
2:B:43:LYS:HB2	2:B:44:GLY:HA2	1.95	0.48
1:A:90:THR:HG21	1:A:240:LYS:CA	2.43	0.48
1:G:62:GLU:CD	1:G:63:THR:H	2.22	0.47
1:A:86:LEU:O	1:A:86:LEU:CG	2.62	0.47
1:A:59:LYS:HD2	1:A:62:GLU:HB2	1.97	0.47
1:A:228:CYS:HA	1:A:242:VAL:HA	1.95	0.47
1:A:53:PHE:CE1	2:B:33:ALA:HB2	2.50	0.47
1:G:51:THR:HA	1:G:103:GLN:HE22	1.80	0.47
2:H:43:LYS:CB	2:H:44:GLY:CA	2.93	0.46
2:H:87:THR:HG23	2:H:110:THR:HA	1.97	0.46
3:C:34:SER:OG	3:C:89:SER:OG	2.31	0.46
1:A:249:HIS:CE1	1:A:486:TYR:OH	2.64	0.46
3:C:66:LYS:O	3:C:66:LYS:HG3	2.16	0.46
1:A:94:ASN:CB	1:A:97:LYS:HB3	2.46	0.46
1:G:101:VAL:HG21	1:G:483:LEU:HD22	1.96	0.46
2:H:143:LYS:HE2	3:L:131:THR:HG21	1.96	0.46
1:A:95:MET:HG3	1:A:484:TYR:CD2	2.51	0.46
1:A:480:ARG:CD	1:A:484:TYR:HE2	2.25	0.46
3:C:134:CYS:HB2	3:C:148:TRP:CH2	2.51	0.46
1:A:100:MET:SD	1:A:487:LYS:HA	2.56	0.46
2:H:72:HIS:CE1	1:A:80:ASN:ND2	2.84	0.46
1:A:86:LEU:O	1:A:86:LEU:HG	2.16	0.46
1:G:99:ASN:HA	1:G:102:GLU:HG2	1.98	0.46
1:A:101:VAL:HG22	1:A:483:LEU:HD13	1.97	0.46
2:B:115:SER:O	2:B:116:THR:C	2.59	0.45
1:G:59:LYS:HD3	1:G:61:HIS:CE1	2.50	0.45
2:B:87:THR:HG23	2:B:110:THR:HA	1.97	0.45
2:B:2:VAL:HG21	2:B:30:SER:HB3	1.97	0.45
1:A:86:LEU:HD21	1:A:242:VAL:N	2.32	0.45
2:B:126:PRO:O	2:B:127:SER:CB	2.65	0.45
1:G:55:ALA:CB	1:G:77:THR:OG1	2.62	0.45
1:G:85:HIS:ND1	1:G:86:LEU:O	2.49	0.45
2:H:43:LYS:HB2	2:H:44:GLY:CA	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:95:MET:SD	1:G:235:GLY:HA3	2.58	0.44
1:A:474:ASN:O	1:A:477:ASP:CG	2.60	0.44
1:A:78:ASP:OD1	1:A:78:ASP:O	2.35	0.44
1:A:86:LEU:HD21	1:A:241:ASN:HA	2.00	0.44
3:C:162:THR:HG22	3:C:163:THR:O	2.18	0.44
1:G:85:HIS:CE1	1:G:86:LEU:O	2.71	0.44
1:G:101:VAL:HG23	1:G:483:LEU:HD22	1.99	0.44
1:A:239:CYS:SG	1:A:240:LYS:N	2.91	0.44
3:L:142:GLY:HA3	3:L:172:TYR:CD1	2.53	0.43
1:G:101:VAL:O	1:G:101:VAL:HG12	2.18	0.43
2:H:178:LEU:C	2:H:178:LEU:HD12	2.44	0.43
2:H:1:GLN:O	2:H:3:GLN:HG3	2.18	0.43
1:G:80:ASN:O	1:G:80:ASN:ND2	2.52	0.43
1:A:94:ASN:HB2	1:A:97:LYS:HB3	2.01	0.43
2:H:75:GLU:O	2:H:76:ASN:C	2.62	0.42
1:A:86:LEU:O	1:A:86:LEU:CD1	2.67	0.42
1:G:480:ARG:HA	1:G:483:LEU:HB3	2.01	0.42
1:A:96:TRP:HB3	1:A:480:ARG:NH1	2.34	0.42
1:A:101:VAL:HG22	1:A:483:LEU:HD12	2.00	0.42
3:C:27(B):ASP:HB2	3:C:27(C):VAL:H	1.64	0.42
2:B:75:GLU:O	2:B:76:ASN:C	2.61	0.42
1:G:86:LEU:CD2	1:G:242:VAL:HB	2.50	0.42
3:L:26:THR:O	3:L:27(C):VAL:O	2.38	0.41
2:B:5:GLN:O	2:B:22:CYS:HA	2.20	0.41
2:B:127:SER:HB3	2:B:129:LYS:HG2	2.02	0.41
3:L:55:PRO:HD2	3:L:58:VAL:HG21	2.02	0.41
2:H:42:GLY:O	2:H:44:GLY:HA3	2.21	0.41
2:B:41:PRO:O	2:B:43:LYS:HB2	2.20	0.41
2:B:2:VAL:HG13	2:B:2:VAL:O	2.20	0.41
2:H:83:THR:C	2:H:111:VAL:HG11	2.46	0.41
2:H:97:ARG:HH21	3:L:95:HIS:CE1	2.39	0.41
3:L:27(B):ASP:HB2	3:L:27(C):VAL:H	1.73	0.41
1:G:478:ASN:O	1:G:482:GLU:OE1	2.38	0.41
2:H:34:HIS:CE1	2:H:94:ARG:CZ	3.04	0.41
1:A:86:LEU:HD22	1:A:242:VAL:HB	2.03	0.41
1:A:95:MET:O	1:A:487:LYS:NZ	2.50	0.41
1:A:226:LEU:HB3	1:A:242:VAL:CG1	2.51	0.41
2:H:35(B):SER:HB3	2:H:50:TYR:HB3	2.03	0.41
1:G:99:ASN:O	1:G:99:ASN:ND2	2.54	0.40
1:A:101:VAL:HG21	1:A:484:TYR:CZ	2.56	0.40
1:G:483:LEU:O	1:G:484:TYR:C	2.64	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)
3:C:93:ASP:OD1	3:C:151:ASP:OD2[2_455]	2.06	0.14

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	121/154 (79%)	101 (84%)	16 (13%)	4 (3%)	3	17
1	G	118/154 (77%)	108 (92%)	7 (6%)	3 (2%)	4	21
2	B	218/224 (97%)	187 (86%)	19 (9%)	12 (6%)	1	9
2	H	208/224 (93%)	178 (86%)	20 (10%)	10 (5%)	2	11
3	C	208/210 (99%)	187 (90%)	19 (9%)	2 (1%)	12	41
3	L	208/210 (99%)	185 (89%)	21 (10%)	2 (1%)	12	41
All	All	1081/1176 (92%)	946 (88%)	102 (9%)	33 (3%)	3	18

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	88	ASN
1	G	484	TYR
2	H	114	ALA
3	L	60	ASP
1	A	238	PRO
1	A	242	VAL
2	B	41	PRO
2	B	64	LYS
2	B	126	PRO
2	B	127	SER
3	C	93	ASP
1	G	105	GLN

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Mol	Chain	Res	Type
2	H	43	LYS
2	H	64	LYS
1	A	239	CYS
1	A	243	SER
2	B	32	GLY
2	B	42	GLY
2	H	16	GLN
2	H	44	GLY
2	H	26	GLY
2	H	74	SER
2	H	187	SER
2	B	31	SER
2	B	43	LYS
2	B	116	THR
2	H	53	TYR
2	B	44	GLY
2	H	213	PRO
2	B	213	PRO
3	C	9	PRO
2	B	133	GLY
3	L	28	GLY

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	118/137 (86%)	109 (92%)	9 (8%)	12	39
1	G	116/137 (85%)	108 (93%)	8 (7%)	14	41
2	B	192/193 (100%)	181 (94%)	11 (6%)	18	49
2	H	184/193 (95%)	171 (93%)	13 (7%)	13	41
3	C	177/177 (100%)	165 (93%)	12 (7%)	14	42
3	L	177/177 (100%)	167 (94%)	10 (6%)	19	49
All	All	964/1014 (95%)	901 (94%)	63 (6%)	15	44

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	77	THR
1	G	87	GLU
1	G	89	VAL
1	G	92	ASN
1	G	110	SER
1	G	213	ILE
1	G	252	LYS
1	G	483	LEU
2	H	7	SER
2	H	28	SER
2	H	50	TYR
2	H	76	ASN
2	H	96	THR
2	H	105	GLN
2	H	109	VAL
2	H	135	THR
2	H	150	VAL
2	H	151	THR
2	H	169	VAL
2	H	178	LEU
2	H	206	LYS
3	L	19	VAL
3	L	42	LYS
3	L	49	SER
3	L	81	GLU
3	L	131	THR
3	L	145	THR
3	L	162	THR
3	L	165	SER
3	L	186	LYS
3	L	195	VAL
1	A	65	CYS
1	A	90	THR
1	A	98	ASN
1	A	115	CYS
1	A	116	LEU
1	A	241	ASN
1	A	476	LYS
1	A	486	TYR
1	A	487	LYS
2	B	23	THR
2	B	50	TYR

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Mol	Chain	Res	Type
2	B	73	THR
2	B	74	SER
2	B	96	THR
2	B	113	SER
2	B	120	SER
2	B	150	VAL
2	B	151	THR
2	B	169	VAL
2	B	178	LEU
3	C	4	VAL
3	C	61	ARG
3	C	73	LEU
3	C	81	GLU
3	C	97	VAL
3	C	121	SER
3	C	135	LEU
3	C	156	LYS
3	C	194	GLN
3	C	195	VAL
3	C	200	SER
3	C	201	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	61	HIS
1	G	66	HIS
1	G	92	ASN
1	G	103	GLN
3	L	95	HIS
3	L	169	ASN
1	A	67	ASN
1	A	72	HIS
2	B	39	GLN
2	B	76	ASN
3	C	53	ASN
3	C	79	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	A	129/154 (83%)	0.05	0 100 100	98, 136, 176, 196	0
1	G	126/154 (81%)	0.06	0 100 100	95, 133, 172, 201	0
2	B	222/224 (99%)	-0.38	1 (0%) 87 73	57, 84, 129, 157	0
2	H	214/224 (95%)	-0.43	0 100 100	60, 79, 111, 120	0
3	C	210/210 (100%)	-0.45	1 (0%) 87 73	53, 73, 100, 115	0
3	L	210/210 (100%)	-0.44	1 (0%) 87 73	64, 85, 105, 116	0
All	All	1111/1176 (94%)	-0.31	3 (0%) 90 80	53, 86, 158, 201	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	41	GLY	2.4
3	C	95	HIS	2.4
2	B	127	SER	2.3

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.