



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 04:09 AM EDT

PDB ID : 7L79 / pdb_00007L79
Title : Crystal structure of broadly HIV-1-neutralizing antibody VRC40.01
Authors : Zhou, T.; Kwong, P.D.
Deposited on : 2020-12-27
Resolution : 2.83 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

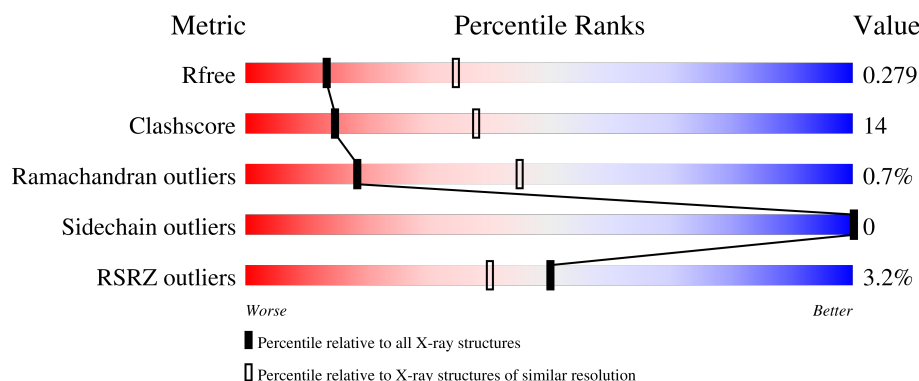
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.83 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	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	230	5% 73% 27%
1	H	230	5% 67% 31% .
2	B	214	% 75% 24%
2	L	214	% 77% 22% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9872 atoms, of which 2987 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 Heavy chain of VRC40.01.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	H	229	Total	C	H	N	O	S	0	0	0
			2444	1102	715	297	324	6			
1	A	229	Total	C	H	N	O	S	0	0	0
			2425	1102	696	297	324	6			

- Molecule 2 is a protein called Light chain of VRC40.01.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	L	214	Total	C	H	N	O	S	0	0	0
			2447	1036	788	286	330	7			
2	B	214	Total	C	H	N	O	S	0	0	0
			2447	1036	788	286	330	7			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	1	Total	Zn	0	0
			1	1		
3	L	5	Total	Zn	0	0
			5	5		
3	B	3	Total	Zn	0	0
			3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	22	Total	O	0	0
			22	22		
4	L	37	Total	O	0	0
			37	37		
4	A	20	Total	O	0	0
			20	20		

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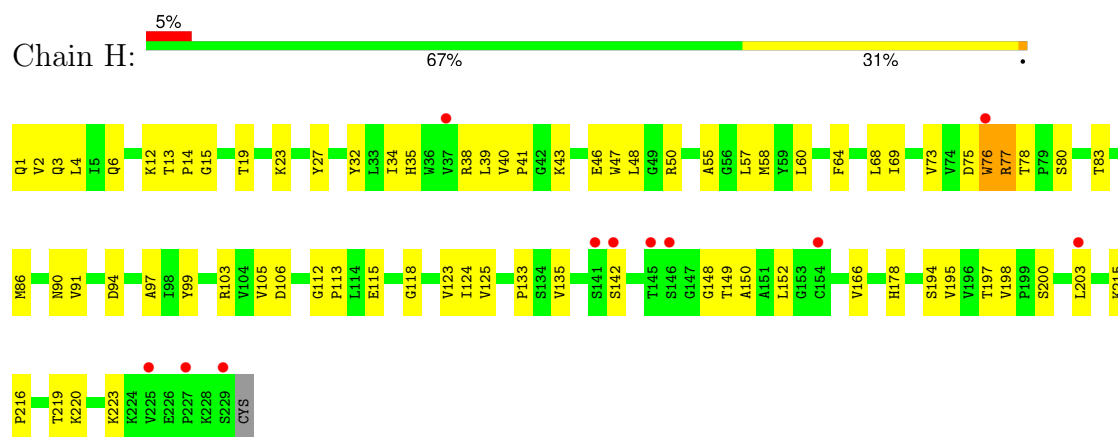
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	21	Total	O	0	0
			21	21		

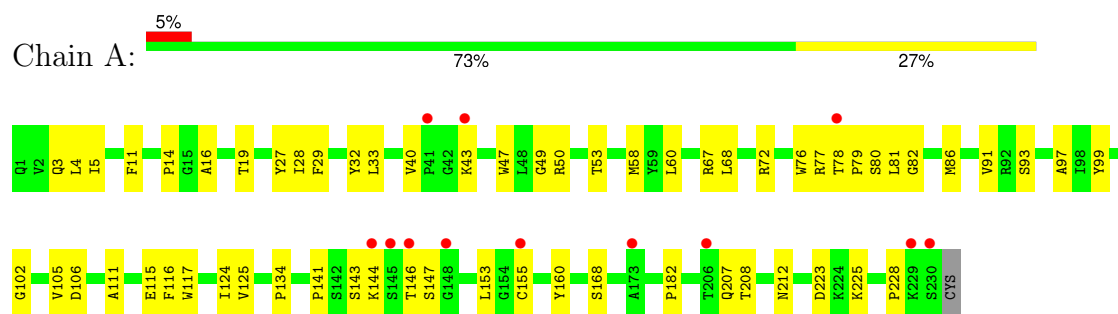
3 Residue-property plots

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.

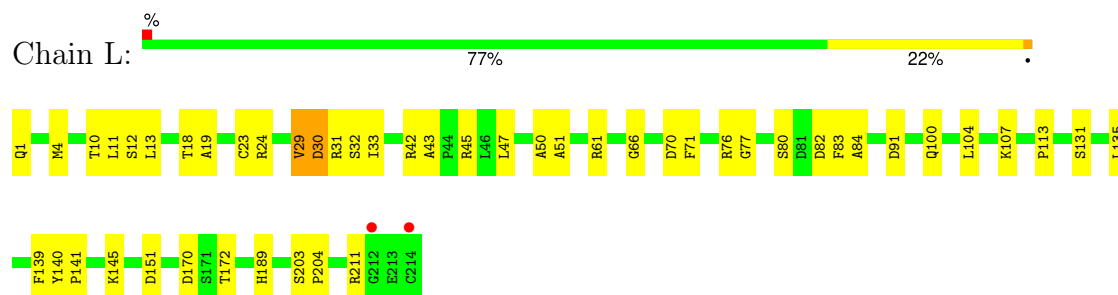
- Molecule 1: Heavy chain of VRC40.01



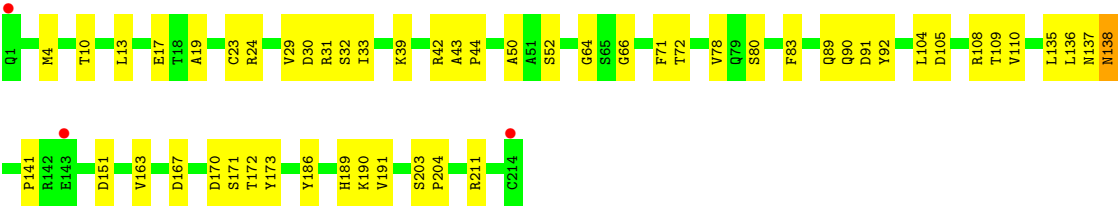
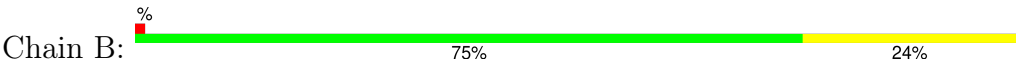
- Molecule 1: Heavy chain of VRC40.01



- Molecule 2: Light chain of VRC40.01



- Molecule 2: Light chain of VRC40.01



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.79Å 135.95Å 136.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.92 – 2.83 36.92 – 2.83	Depositor EDS
% Data completeness (in resolution range)	97.0 (36.92-2.83) 96.9 (36.92-2.83)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.81Å)	Xtriage
Refinement program	PHENIX V1.1	Depositor
R, R_{free}	0.220 , 0.280 0.222 , 0.279	Depositor DCC
R_{free} test set	1398 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	55.3	Xtriage
Anisotropy	0.119	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9872	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.07 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2064e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: ZN

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	0.36	1/1773 (0.1%)	0.57	0/2416
1	H	0.30	0/1773	0.51	0/2416
2	B	0.29	0/1695	0.50	0/2301
2	L	0.30	0/1695	0.51	0/2301
All	All	0.32	1/6936 (0.0%)	0.52	0/9434

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	H	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	155	CYS	CB-SG	-5.94	1.61	1.81

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	75	ASP	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	1729	696	1731	53	0
1	H	1729	715	1730	53	2
2	B	1659	788	1608	41	0
2	L	1659	788	1607	45	0
3	B	3	0	0	0	0
3	H	1	0	0	0	0
3	L	5	0	0	0	0
4	A	20	0	0	1	0
4	B	21	0	0	7	0
4	H	22	0	0	1	0
4	L	37	0	0	10	0
All	All	6885	2987	6676	187	2

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 14.

All (187) 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:B:151:ASP:OD2	4:B:401:HOH:O	1.84	0.95
2:L:18:THR:HG23	2:L:76:ARG:HA	1.47	0.93
2:L:151:ASP:OD2	4:L:401:HOH:O	1.91	0.89
2:L:30:ASP:HB3	4:L:413:HOH:O	1.74	0.87
2:L:12:SER:HB3	2:L:107:LYS:HD2	1.55	0.86
2:B:13:LEU:HD12	2:B:104:LEU:HD11	1.56	0.86
2:L:100:GLN:OE1	4:L:402:HOH:O	1.92	0.85
1:A:40:VAL:CG2	1:A:43:LYS:HB2	2.06	0.84
1:A:40:VAL:HG23	1:A:43:LYS:HB2	1.61	0.82
1:H:76:TRP:O	1:H:77:ARG:HB2	1.78	0.82
2:L:145:LYS:NZ	4:L:403:HOH:O	2.12	0.80
1:H:149:THR:HG21	1:H:197:THR:HG22	1.63	0.79
1:A:40:VAL:HG12	1:A:97:ALA:HB2	1.65	0.77
1:A:58:MET:HE3	1:A:60:LEU:HD21	1.66	0.77
2:B:50:ALA:O	4:B:402:HOH:O	2.04	0.76
2:B:17:GLU:O	2:B:78:VAL:HG23	1.85	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:52:SER:OG	4:B:402:HOH:O	2.03	0.74
1:H:47:TRP:HZ2	1:H:50:ARG:HG2	1.53	0.73
2:L:76:ARG:HG3	2:L:77:GLY:H	1.54	0.73
2:L:80:SER:HA	2:L:83:PHE:CE1	2.25	0.71
1:A:27:TYR:O	1:A:80:SER:HA	1.92	0.70
1:H:148:GLY:O	1:H:200:SER:OG	2.08	0.70
1:A:58:MET:CE	1:A:60:LEU:HD11	2.22	0.69
2:L:45:ARG:HG2	4:L:422:HOH:O	1.92	0.69
1:A:58:MET:HE1	1:A:60:LEU:HD11	1.75	0.68
2:L:47:LEU:HD23	4:L:422:HOH:O	1.93	0.67
2:B:190:LYS:HD2	2:B:191:VAL:HG23	1.76	0.67
2:L:12:SER:CB	2:L:107:LYS:HD2	2.25	0.67
1:A:58:MET:CE	1:A:60:LEU:HD21	2.25	0.66
2:B:10:THR:HG22	4:B:410:HOH:O	1.96	0.66
1:A:144:LYS:O	1:A:146:THR:N	2.26	0.66
2:L:131:SER:O	4:L:404:HOH:O	2.13	0.65
2:L:4:MET:O	4:L:402:HOH:O	2.15	0.64
1:H:195:VAL:HG21	2:L:135:LEU:HD22	1.78	0.64
1:A:86:MET:HE1	1:A:99:TYR:CD2	2.32	0.64
1:H:27:TYR:O	1:H:80:SER:HA	1.98	0.63
1:A:3:GLN:O	1:A:4:LEU:HD23	1.99	0.63
2:L:66:GLY:HA3	2:L:71:PHE:HA	1.80	0.63
1:H:152:LEU:HD23	1:H:203:LEU:HD21	1.79	0.62
1:H:34:ILE:HD13	1:H:103:ARG:HB3	1.81	0.62
1:A:28:ILE:HD12	1:A:79:PRO:HB2	1.82	0.62
1:A:72:ARG:O	1:A:72:ARG:HD2	2.00	0.61
1:A:76:TRP:O	1:A:77:ARG:HB3	2.00	0.61
1:A:19:THR:HA	1:A:86:MET:O	2.00	0.61
2:L:189:HIS:O	2:L:211:ARG:NH1	2.34	0.61
2:L:170:ASP:HB2	2:L:172:THR:HG23	1.82	0.61
1:A:47:TRP:HZ2	1:A:50:ARG:HG2	1.66	0.61
2:B:108:ARG:HD2	4:B:405:HOH:O	1.99	0.60
1:H:68:LEU:HD11	1:H:86:MET:HG3	1.82	0.60
1:A:143:SER:HB3	1:A:147:SER:OG	2.02	0.59
1:H:133:PRO:HD2	1:H:219:THR:HG21	1.85	0.58
1:A:16:ALA:O	1:A:91:VAL:HG22	2.03	0.58
1:H:40:VAL:HG13	1:H:41:PRO:HD2	1.85	0.58
1:H:15:GLY:HA2	1:H:90:ASN:OD1	2.04	0.58
2:L:29:VAL:O	2:L:30:ASP:C	2.47	0.58
1:A:32:TYR:HE2	1:A:106:ASP:HA	1.70	0.57
2:B:170:ASP:OD1	2:B:172:THR:OG1	2.21	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:58:MET:HE1	1:H:69:ILE:HG23	1.86	0.57
2:L:82:ASP:O	2:L:104:LEU:HD23	2.05	0.57
2:L:4:MET:HE3	2:L:23:CYS:SG	2.44	0.56
1:A:40:VAL:HG12	1:A:97:ALA:CB	2.34	0.56
2:B:13:LEU:HD12	2:B:104:LEU:CD1	2.33	0.56
1:A:3:GLN:C	1:A:4:LEU:HD23	2.31	0.56
2:B:4:MET:HE2	2:B:23:CYS:SG	2.44	0.56
1:A:14:PRO:HA	1:A:91:VAL:O	2.06	0.56
1:A:153:LEU:C	1:A:153:LEU:HD12	2.31	0.55
1:A:72:ARG:HD2	1:A:72:ARG:C	2.32	0.55
1:A:117:TRP:CZ3	2:B:44:PRO:HG2	2.41	0.55
1:H:3:GLN:C	1:H:4:LEU:HD12	2.33	0.54
2:B:135:LEU:C	2:B:136:LEU:HD12	2.32	0.54
1:H:55:ALA:HB3	1:H:57:LEU:HD12	1.88	0.54
2:L:189:HIS:ND1	4:L:401:HOH:O	2.26	0.54
1:H:47:TRP:CZ2	1:H:50:ARG:HG2	2.41	0.53
2:L:24:ARG:HH21	2:L:24:ARG:HG3	1.72	0.53
2:B:137:ASN:ND2	2:B:138:ASN:OD1	2.41	0.53
1:H:55:ALA:HB3	1:H:57:LEU:CD1	2.37	0.53
1:H:105:VAL:HG22	1:H:115:GLU:HB2	1.91	0.53
1:A:11:PHE:CE2	1:A:124:ILE:HD13	2.44	0.53
1:A:29:PHE:CD1	1:A:82:GLY:HA3	2.43	0.53
1:A:68:LEU:HD11	1:A:86:MET:HG3	1.90	0.52
1:A:86:MET:HE1	1:A:99:TYR:CE2	2.45	0.52
1:H:19:THR:HA	1:H:86:MET:O	2.11	0.51
1:A:47:TRP:CZ2	1:A:49:GLY:HA2	2.45	0.51
1:H:38:ARG:HD3	1:H:99:TYR:CE2	2.46	0.50
2:L:113:PRO:HB3	2:L:139:PHE:HB3	1.92	0.50
2:L:203:SER:HB2	2:L:204:PRO:HD2	1.93	0.50
1:H:32:TYR:HE2	1:H:106:ASP:HA	1.76	0.50
2:L:12:SER:O	2:L:13:LEU:HG	2.11	0.50
2:B:110:VAL:HG22	2:B:141:PRO:HD3	1.93	0.50
1:H:149:THR:HG23	1:H:198:VAL:O	2.11	0.50
2:L:83:PHE:O	2:L:84:ALA:HB2	2.13	0.49
1:A:141:PRO:HD2	1:A:228:PRO:HA	1.95	0.49
1:A:58:MET:HE2	1:A:60:LEU:CG	2.43	0.49
1:H:35:HIS:CE1	1:H:50:ARG:HD3	2.48	0.49
2:B:108:ARG:NE	2:B:109:THR:O	2.45	0.49
1:A:29:PHE:CE2	1:A:53:THR:HB	2.48	0.48
1:H:13:THR:OG1	1:H:14:PRO:HD2	2.12	0.48
1:H:73:VAL:O	1:H:73:VAL:HG23	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:208:THR:HG23	1:A:225:LYS:HE2	1.95	0.48
1:H:166:VAL:HG11	1:H:194:SER:CB	2.43	0.48
2:L:170:ASP:OD1	2:L:170:ASP:N	2.34	0.48
1:H:6:GLN:OE1	1:H:118:GLY:HA3	2.14	0.48
1:A:58:MET:HE2	1:A:60:LEU:HG	1.96	0.48
1:H:38:ARG:HD3	1:H:99:TYR:CZ	2.49	0.48
1:H:38:ARG:HH21	1:H:46:GLU:CD	2.21	0.47
2:L:24:ARG:HG3	2:L:24:ARG:NH2	2.30	0.47
2:L:32:SER:HB3	2:L:91:ASP:OD1	2.14	0.47
1:H:124:ILE:HD12	1:H:124:ILE:N	2.29	0.47
1:A:5:ILE:HD12	1:A:5:ILE:N	2.30	0.46
1:A:143:SER:HB3	1:A:147:SER:CB	2.46	0.46
2:B:80:SER:HA	2:B:83:PHE:CE1	2.50	0.46
2:L:10:THR:HG22	2:L:11:LEU:N	2.30	0.46
2:B:13:LEU:CD1	2:B:19:ALA:HB2	2.44	0.46
2:B:189:HIS:ND1	4:B:401:HOH:O	2.28	0.46
1:H:91:VAL:HB	1:H:125:VAL:HG21	1.97	0.46
1:H:123:VAL:HG13	1:H:123:VAL:O	2.16	0.46
2:B:105:ASP:OD2	2:B:173:TYR:OH	2.32	0.46
2:L:24:ARG:NH2	2:L:70:ASP:OD1	2.49	0.46
2:B:30:ASP:HB2	2:B:92:TYR:CE1	2.51	0.46
1:H:112:GLY:N	1:H:113:PRO:CD	2.79	0.45
1:H:149:THR:CG2	1:H:150:ALA:N	2.79	0.45
1:H:12:LYS:O	1:H:125:VAL:HA	2.16	0.45
2:L:76:ARG:HG3	2:L:77:GLY:N	2.27	0.45
2:B:39:LYS:HD2	2:B:42:ARG:NH2	2.31	0.45
2:B:64:GLY:HA2	2:B:72:THR:O	2.17	0.45
2:B:186:TYR:CZ	2:B:211:ARG:HG3	2.51	0.45
2:L:80:SER:HA	2:L:83:PHE:CD1	2.51	0.45
1:H:1:GLN:HG3	1:H:2:VAL:H	1.81	0.45
2:B:30:ASP:HB3	2:B:31:ARG:H	1.63	0.45
1:H:94:ASP:OD1	4:H:401:HOH:O	2.21	0.45
1:H:48:LEU:HD23	1:H:64:PHE:CE2	2.52	0.45
2:L:1:GLN:HA	2:L:1:GLN:OE1	2.16	0.44
1:H:135:VAL:O	1:H:223:LYS:HE3	2.16	0.44
1:H:195:VAL:HG21	2:L:135:LEU:CD2	2.45	0.44
1:A:33:LEU:HD23	1:A:50:ARG:HD2	1.99	0.44
2:B:203:SER:HB2	2:B:204:PRO:HD2	1.99	0.44
1:A:78:THR:HG21	1:A:81:LEU:HD12	1.99	0.44
2:B:13:LEU:HD13	2:B:19:ALA:HB2	1.98	0.44
1:A:111:ALA:HB1	2:B:91:ASP:OD2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:29:VAL:O	2:B:30:ASP:C	2.61	0.44
1:H:39:LEU:O	1:H:97:ALA:HB1	2.18	0.44
1:H:32:TYR:CD2	1:H:103:ARG:NH1	2.82	0.44
1:H:77:ARG:HB3	1:H:78:THR:HG23	1.99	0.44
1:A:80:SER:O	1:A:81:LEU:HD23	2.18	0.44
2:L:13:LEU:HD13	2:L:19:ALA:HB2	1.98	0.44
2:B:32:SER:HB3	2:B:91:ASP:OD2	2.18	0.43
1:A:134:PRO:HB3	1:A:160:TYR:HB3	2.00	0.43
2:B:24:ARG:HG3	2:B:24:ARG:HH21	1.83	0.43
1:A:58:MET:HE2	1:A:60:LEU:HD11	2.00	0.43
2:B:33:ILE:HG13	2:B:71:PHE:CE1	2.54	0.43
2:B:42:ARG:HG2	2:B:43:ALA:N	2.33	0.43
2:B:66:GLY:HA3	2:B:71:PHE:HA	2.01	0.43
1:H:32:TYR:CE2	1:H:106:ASP:HA	2.54	0.43
1:A:67:ARG:HD3	4:A:310:HOH:O	2.19	0.43
1:H:40:VAL:CG1	1:H:41:PRO:HD2	2.49	0.42
1:A:168:SER:HB2	1:A:212:ASN:HB2	2.01	0.42
1:A:207:GLN:HG3	1:A:208:THR:N	2.35	0.42
2:L:61:ARG:HH21	2:L:82:ASP:CG	2.27	0.42
2:B:108:ARG:CD	4:B:405:HOH:O	2.63	0.42
2:B:167:ASP:O	2:B:171:SER:HA	2.20	0.42
1:H:58:MET:HG3	1:H:60:LEU:HD21	2.01	0.42
1:A:102:GLY:HA2	1:A:116:PHE:O	2.20	0.42
1:H:40:VAL:HG22	1:H:97:ALA:HB2	2.00	0.42
1:H:178:HIS:CE1	4:L:410:HOH:O	2.73	0.42
2:L:13:LEU:CD1	2:L:19:ALA:HB2	2.50	0.42
2:L:140:TYR:CG	2:L:141:PRO:HA	2.55	0.42
1:A:105:VAL:HG22	1:A:115:GLU:HB2	2.01	0.42
2:B:4:MET:HE3	2:B:4:MET:HA	2.01	0.42
2:L:30:ASP:HB3	2:L:31:ARG:H	1.68	0.41
2:L:203:SER:HB2	2:L:204:PRO:CD	2.50	0.41
2:B:80:SER:HA	2:B:83:PHE:CD1	2.55	0.41
2:B:89:GLN:HG2	2:B:90:GLN:N	2.34	0.41
1:H:149:THR:HG21	1:H:197:THR:CG2	2.41	0.41
1:A:182:PRO:HG2	2:B:163:VAL:O	2.20	0.41
2:B:108:ARG:NH2	2:B:170:ASP:O	2.53	0.41
2:L:33:ILE:HG13	2:L:71:PHE:CE1	2.56	0.41
1:A:93:SER:HA	1:A:125:VAL:HB	2.02	0.41
1:H:86:MET:HE1	1:H:99:TYR:CD2	2.55	0.41
2:L:13:LEU:HD12	2:L:104:LEU:HD11	2.02	0.41
1:A:212:ASN:HA	1:A:223:ASP:OD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:215:LYS:N	1:H:216:PRO:CD	2.83	0.40
1:A:28:ILE:CD1	1:A:79:PRO:HB2	2.50	0.40
1:A:144:LYS:HA	1:A:144:LYS:HD3	1.87	0.40
1:H:99:TYR:CE1	1:H:123:VAL:CG1	3.05	0.40
2:L:42:ARG:HG2	2:L:43:ALA:N	2.36	0.40
2:L:50:ALA:O	2:L:51:ALA:HB3	2.20	0.40
1:H:23:LYS:HG2	1:H:83:THR:HG23	2.04	0.40

All (2) 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)
1:H:142:SER:O	1:H:220:LYS:HZ2[4_555]	1.25	0.35
1:H:142:SER:O	1:H:220:LYS:NZ[4_555]	1.86	0.34

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	227/230 (99%)	210 (92%)	17 (8%)	0	100	100
1	H	227/230 (99%)	212 (93%)	12 (5%)	3 (1%)	9	29
2	B	212/214 (99%)	194 (92%)	17 (8%)	1 (0%)	24	53
2	L	212/214 (99%)	195 (92%)	15 (7%)	2 (1%)	14	38
All	All	878/888 (99%)	811 (92%)	61 (7%)	6 (1%)	18	45

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	76	TRP
1	H	77	ARG

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Mol	Chain	Res	Type
2	B	138	ASN
1	H	43	LYS
2	L	30	ASP
2	L	29	VAL

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	195/196 (100%)	195 (100%)	0	100	100
1	H	195/196 (100%)	195 (100%)	0	100	100
2	B	186/186 (100%)	186 (100%)	0	100	100
2	L	186/186 (100%)	186 (100%)	0	100	100
All	All	762/764 (100%)	762 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	1	GLN
1	H	213	ASN
2	L	89	GLN
1	A	214	ASN
2	B	89	GLN
2	B	160	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](#)

Of 9 ligands modelled in this entry, 9 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	229/230 (99%)	0.40	12 (5%) 33 24	29, 51, 91, 132	0
1	H	229/230 (99%)	0.36	11 (4%) 35 27	28, 50, 102, 143	0
2	B	214/214 (100%)	0.23	3 (1%) 73 64	28, 45, 80, 130	0
2	L	214/214 (100%)	0.24	2 (0%) 81 74	25, 45, 89, 147	0
All	All	886/888 (99%)	0.31	28 (3%) 50 40	25, 48, 92, 147	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	142	SER	5.4
1	H	76	TRP	4.0
1	A	41	PRO	3.8
1	A	144	LYS	3.5
2	B	1	GLN	3.0
1	H	229	SER	2.8
2	B	214	CYS	2.7
1	A	145	SER	2.7
2	L	214	CYS	2.6
1	H	225	VAL	2.6
1	A	43	LYS	2.6
1	H	203	LEU	2.6
1	A	173	ALA	2.5
1	A	229	LYS	2.4
1	A	155	CYS	2.4
1	H	146	SER	2.4
1	H	145	THR	2.3
1	A	148	GLY	2.3
2	B	143	GLU	2.2
2	L	212	GLY	2.2
1	A	78	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	206	THR	2.1
1	H	227	PRO	2.1
1	H	37	VAL	2.1
1	A	230	SER	2.1
1	H	154	CYS	2.0
1	A	146	THR	2.0
1	H	141	SER	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 oligosaccharides 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
3	ZN	L	304	1/1	0.67	0.24	110,110,110,110	0
3	ZN	L	303	1/1	0.84	0.08	53,53,53,53	0
3	ZN	L	305	1/1	0.94	0.21	89,89,89,89	0
3	ZN	B	301	1/1	0.95	0.12	59,59,59,59	0
3	ZN	B	303	1/1	0.95	0.09	47,47,47,47	0
3	ZN	L	302	1/1	0.96	0.13	61,61,61,61	0
3	ZN	H	301	1/1	0.96	0.15	54,54,54,54	0
3	ZN	B	302	1/1	0.97	0.04	40,40,40,40	0
3	ZN	L	301	1/1	0.99	0.04	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.