



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:12 AM UTC

PDB ID : 7PHS / pdb_00007phs
Title : highly potent IL6 antagonistic antibody selected from a camelid immune phage display repertoire
Authors : Cambillau, C.; Pannecoucke, E.; De Haard, H.
Deposited on : 2021-08-18
Resolution : 2.77 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

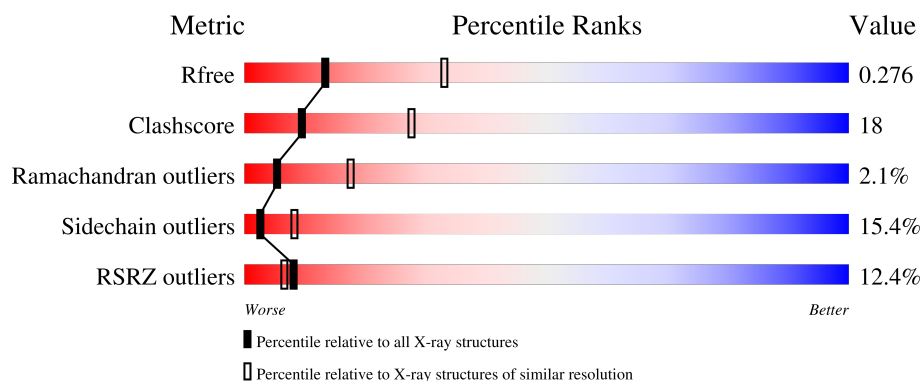
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	H	223	15% 61% 30% 7%
2	L	212	9% 71% 22% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3325 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 Heavy chain of Fab129D3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	223	Total	C	N	O	S	0	0	0
			1667	1059	269	334	5			

- Molecule 2 is a protein called Light chain of Fab129D3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	211	Total	C	N	O	S	0	0	0
			1569	984	261	319	5			

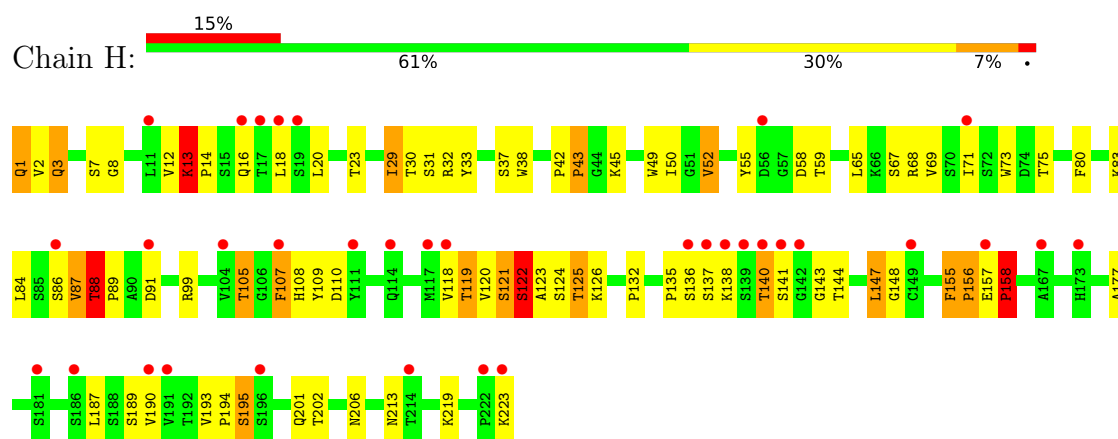
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	51	Total	O	0	0
			51	51		
3	L	38	Total	O	0	0
			38	38		

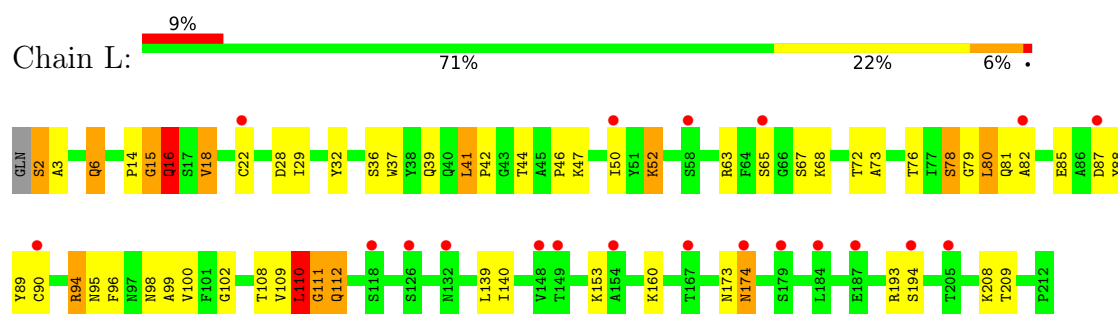
3 Residue-property plots [i](#)

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: Heavy chain of Fab129D3



- Molecule 2: Light chain of Fab129D3



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	58.99Å 58.99Å 285.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.05 – 2.77 41.05 – 2.77	Depositor EDS
% Data completeness (in resolution range)	98.4 (41.05-2.77) 98.7 (41.05-2.77)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.29 (at 2.77Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.273 , 0.303 (Not available) , 0.276	Depositor DCC
R_{free} test set	1344 reflections (9.91%)	wwPDB-VP
Wilson B-factor (Å ²)	51.9	Xtriage
Anisotropy	0.761	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	3325	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% 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](#)

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	H	1.01	4/1714 (0.2%)	1.54	23/2349 (1.0%)
2	L	0.84	1/1607 (0.1%)	1.27	9/2197 (0.4%)
All	All	0.93	5/3321 (0.2%)	1.42	32/4546 (0.7%)

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	2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	88	THR	C-N	15.62	1.50	1.34
1	H	155	PHE	C-N	8.20	1.49	1.34
1	H	13	LYS	C-N	6.88	1.50	1.33
2	L	46	PRO	C-O	-5.49	1.17	1.23
1	H	132	PRO	C-O	-5.12	1.18	1.23

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	122	SER	CA-C-N	15.27	145.98	122.93
1	H	122	SER	C-N-CA	15.27	145.98	122.93
1	H	88	THR	CA-C-N	11.97	131.55	119.82
1	H	88	THR	C-N-CA	11.97	131.55	119.82
1	H	13	LYS	CA-C-N	10.94	133.52	119.84
1	H	13	LYS	C-N-CA	10.94	133.52	119.84
1	H	58	ASP	CB-CA-C	7.89	122.84	109.50
1	H	88	THR	O-C-N	-7.65	114.31	121.20
1	H	122	SER	N-CA-C	7.59	122.68	113.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	110	LEU	N-CA-C	7.26	124.92	114.39
1	H	107	PHE	CA-C-N	6.90	134.71	121.54
1	H	107	PHE	C-N-CA	6.90	134.71	121.54
1	H	135	PRO	N-CA-C	6.58	121.57	111.57
2	L	111	GLY	CA-C-N	6.48	129.40	120.39
2	L	111	GLY	C-N-CA	6.48	129.40	120.39
2	L	16	GLN	CA-C-N	6.41	133.23	121.70
2	L	16	GLN	C-N-CA	6.41	133.23	121.70
1	H	88	THR	CB-CA-C	-6.09	99.60	108.91
1	H	42	PRO	CA-C-N	6.03	127.37	119.84
1	H	42	PRO	C-N-CA	6.03	127.37	119.84
1	H	105	THR	CB-CA-C	5.92	117.16	109.28
2	L	42	PRO	N-CA-C	5.79	120.30	111.15
1	H	155	PHE	CA-C-O	5.76	124.96	119.72
2	L	82	ALA	CA-C-N	5.64	127.84	120.28
2	L	82	ALA	C-N-CA	5.64	127.84	120.28
1	H	201	GLN	N-CA-C	5.30	118.12	110.28
1	H	43	PRO	N-CA-CB	-5.29	97.69	103.25
1	H	158	PRO	N-CA-CB	-5.24	96.83	102.60
1	H	119	THR	CB-CA-C	5.21	119.12	109.70
1	H	31	SER	CA-C-N	5.12	131.32	121.54
1	H	31	SER	C-N-CA	5.12	131.32	121.54
2	L	44	THR	CA-CB-OG1	-5.05	102.03	109.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	122	SER	Peptide,Mainchain

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	H	1667	0	1606	66	0
2	L	1569	0	1524	60	0
3	H	51	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	38	0	0	1	0
All	All	3325	0	3130	114	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 18.

All (114) 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:H:105:THR:HG22	2:L:52:LYS:NZ	1.60	1.15
1:H:138:LYS:NZ	2:L:209:THR:O	1.79	1.13
2:L:14:PRO:HG3	2:L:111:GLY:HA3	1.24	1.06
1:H:105:THR:HG22	2:L:52:LYS:HZ1	1.16	0.99
2:L:2:SER:N	2:L:94:ARG:HH22	1.69	0.91
1:H:143:GLY:O	1:H:195:SER:N	2.06	0.87
2:L:94:ARG:NH1	2:L:100:VAL:HG21	1.94	0.83
2:L:94:ARG:HH11	2:L:100:VAL:CB	1.93	0.82
1:H:59:THR:HG21	1:H:71:ILE:CG2	2.09	0.82
2:L:94:ARG:HH11	2:L:100:VAL:HB	1.44	0.82
2:L:16:GLN:HB3	2:L:79:GLY:CA	2.10	0.80
2:L:94:ARG:NH1	2:L:100:VAL:CG2	2.45	0.80
2:L:16:GLN:HB3	2:L:79:GLY:HA2	1.65	0.78
2:L:16:GLN:HB3	2:L:79:GLY:N	1.98	0.78
1:H:105:THR:CG2	2:L:52:LYS:NZ	2.44	0.77
2:L:80:LEU:N	2:L:80:LEU:HD23	2.01	0.76
1:H:1:GLN:HG2	1:H:2:VAL:H	1.51	0.74
1:H:105:THR:HG22	2:L:52:LYS:HZ3	1.49	0.74
1:H:49:TRP:HZ2	1:H:52:VAL:HG22	1.53	0.74
2:L:2:SER:N	2:L:94:ARG:NH2	2.38	0.71
1:H:33:TYR:HB2	3:H:311:HOH:O	1.91	0.70
2:L:18:VAL:HG13	2:L:80:LEU:HD11	1.74	0.70
1:H:105:THR:CG2	2:L:52:LYS:HZ3	2.04	0.69
1:H:143:GLY:O	1:H:194:PRO:HA	1.92	0.69
2:L:94:ARG:NH1	2:L:100:VAL:HB	2.08	0.68
1:H:137:SER:HA	1:H:140:THR:HG23	1.75	0.67
1:H:1:GLN:CG	1:H:2:VAL:H	2.08	0.67
1:H:14:PRO:HA	1:H:87:VAL:O	1.95	0.66
2:L:94:ARG:NH1	2:L:100:VAL:CB	2.60	0.65
1:H:121:SER:HB3	1:H:123:ALA:HB3	1.79	0.64
1:H:59:THR:HG21	1:H:71:ILE:HG23	1.79	0.64
2:L:14:PRO:CG	2:L:111:GLY:HA3	2.16	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:15:GLY:O	2:L:79:GLY:HA2	1.98	0.63
2:L:15:GLY:O	2:L:80:LEU:N	2.31	0.63
1:H:125:THR:HG23	1:H:155:PHE:O	1.99	0.63
2:L:14:PRO:HG3	2:L:111:GLY:CA	2.14	0.63
2:L:94:ARG:HH12	2:L:100:VAL:CG2	2.13	0.62
2:L:85:GLU:HG3	2:L:109:VAL:HG23	1.82	0.62
1:H:59:THR:CG2	1:H:71:ILE:CG2	2.78	0.62
2:L:63:ARG:HG2	2:L:63:ARG:HH11	1.64	0.61
2:L:68:LYS:HE3	3:L:311:HOH:O	2.01	0.59
1:H:125:THR:CG2	1:H:155:PHE:O	2.50	0.59
2:L:16:GLN:HB3	2:L:79:GLY:H	1.67	0.59
2:L:94:ARG:HH12	2:L:100:VAL:HG21	1.67	0.58
1:H:49:TRP:CD1	2:L:99:ALA:HB3	2.38	0.58
2:L:37:TRP:CZ3	2:L:90:CYS:HB3	2.39	0.58
2:L:16:GLN:CB	2:L:79:GLY:HA2	2.33	0.57
1:H:107:PHE:CE2	2:L:98:ASN:O	2.57	0.57
1:H:202:THR:HG23	1:H:219:LYS:HE3	1.88	0.55
1:H:18:LEU:HD11	1:H:118:VAL:HG11	1.89	0.55
1:H:107:PHE:HE2	2:L:98:ASN:O	1.90	0.54
1:H:87:VAL:HG12	1:H:91:ASP:OD2	2.06	0.54
1:H:37:SER:HB3	1:H:52:VAL:HG13	1.91	0.52
1:H:1:GLN:HG2	1:H:2:VAL:N	2.21	0.52
1:H:148:GLY:HA3	1:H:189:SER:O	2.09	0.52
2:L:63:ARG:HG2	2:L:63:ARG:NH1	2.25	0.52
1:H:59:THR:HG21	1:H:71:ILE:HG22	1.89	0.51
2:L:39:GLN:HG3	2:L:88:TYR:CE2	2.47	0.50
2:L:94:ARG:HH11	2:L:100:VAL:CG2	2.19	0.50
1:H:99:ARG:O	1:H:109:TYR:HA	2.11	0.50
2:L:29:ILE:HG22	2:L:29:ILE:O	2.12	0.49
1:H:110:ASP:OD2	1:H:110:ASP:N	2.46	0.49
1:H:143:GLY:O	1:H:194:PRO:CA	2.61	0.49
1:H:49:TRP:CG	2:L:99:ALA:HB3	2.48	0.48
1:H:1:GLN:CG	1:H:2:VAL:N	2.76	0.48
1:H:37:SER:HB3	1:H:52:VAL:CG1	2.44	0.47
2:L:22:CYS:O	2:L:73:ALA:N	2.38	0.47
1:H:190:VAL:HG21	2:L:139:LEU:HD13	1.96	0.47
2:L:63:ARG:HB2	2:L:78:SER:O	2.15	0.47
1:H:30:THR:HG23	1:H:75:THR:HG22	1.96	0.46
2:L:95:ASN:HA	2:L:96:PHE:HA	1.69	0.46
1:H:89:PRO:HA	1:H:120:VAL:HB	1.98	0.46
2:L:173:ASN:O	2:L:174:ASN:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:55:TYR:H	1:H:55:TYR:HD1	1.63	0.45
1:H:147:LEU:HD21	1:H:193:VAL:CG2	2.47	0.44
1:H:88:THR:HG23	1:H:88:THR:O	2.16	0.44
1:H:3:GLN:HE21	1:H:3:GLN:HB3	1.58	0.44
1:H:137:SER:OG	1:H:137:SER:O	2.31	0.44
1:H:38:TRP:HB3	1:H:50:ILE:HD12	1.98	0.44
1:H:144:THR:HA	1:H:193:VAL:O	2.18	0.44
1:H:30:THR:CG2	1:H:75:THR:CG2	2.95	0.44
1:H:13:LYS:HB2	1:H:13:LYS:HE3	1.57	0.44
2:L:65:SER:HB3	2:L:76:THR:HB	2.00	0.43
1:H:49:TRP:CD1	1:H:109:TYR:HH	2.37	0.43
2:L:6:GLN:HE22	2:L:89:TYR:HA	1.83	0.43
2:L:80:LEU:N	2:L:80:LEU:CD2	2.71	0.43
2:L:110:LEU:C	2:L:112:GLN:H	2.25	0.43
1:H:49:TRP:CZ2	1:H:52:VAL:HG22	2.43	0.43
1:H:157:GLU:HA	1:H:158:PRO:HA	1.67	0.43
2:L:22:CYS:HB2	2:L:37:TRP:CH2	2.54	0.43
1:H:2:VAL:HG11	1:H:99:ARG:NH1	2.34	0.43
2:L:18:VAL:CG1	2:L:80:LEU:HD11	2.46	0.43
1:H:1:GLN:HG2	1:H:2:VAL:HG13	2.01	0.42
1:H:177:ALA:HA	1:H:187:LEU:HB3	2.02	0.42
2:L:112:GLN:HE21	2:L:112:GLN:HB3	1.55	0.42
1:H:121:SER:HB3	1:H:123:ALA:CB	2.48	0.42
1:H:190:VAL:CG2	2:L:139:LEU:HD13	2.50	0.42
2:L:6:GLN:HE21	2:L:102:GLY:HA3	1.84	0.42
2:L:32:TYR:OH	2:L:95:ASN:HB2	2.19	0.42
2:L:39:GLN:HG3	2:L:88:TYR:CZ	2.54	0.42
2:L:208:LYS:HD3	2:L:208:LYS:HA	1.79	0.42
1:H:73:TRP:HA	1:H:80:PHE:HA	2.02	0.41
1:H:30:THR:CG2	1:H:75:THR:HG22	2.50	0.41
1:H:55:TYR:N	1:H:55:TYR:CD1	2.88	0.41
1:H:155:PHE:HA	1:H:156:PRO:HA	1.85	0.41
2:L:41:LEU:H	2:L:41:LEU:HG	1.70	0.41
1:H:147:LEU:HD21	1:H:193:VAL:HG22	2.03	0.41
1:H:13:LYS:HE3	1:H:16:GLN:NE2	2.35	0.41
1:H:29:ILE:O	1:H:29:ILE:HG12	2.20	0.41
1:H:147:LEU:H	1:H:147:LEU:HG	1.74	0.41
2:L:2:SER:HB2	2:L:3:ALA:H	1.68	0.41
2:L:95:ASN:ND2	2:L:96:PHE:HB2	2.36	0.41
1:H:8:GLY:HA3	1:H:20:LEU:HD23	2.02	0.40
1:H:177:ALA:HB2	1:H:187:LEU:HD23	2.03	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	H	221/223 (99%)	198 (90%)	17 (8%)	6 (3%)	4	12
2	L	209/212 (99%)	194 (93%)	12 (6%)	3 (1%)	9	26
All	All	430/435 (99%)	392 (91%)	29 (7%)	9 (2%)	5	17

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	43	PRO
1	H	108	HIS
1	H	32	ARG
1	H	195	SER
2	L	174	ASN
2	L	28	ASP
2	L	15	GLY
1	H	156	PRO
1	H	158	PRO

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	H	190/194 (98%)	158 (83%)	32 (17%)	2	6
2	L	174/175 (99%)	150 (86%)	24 (14%)	3	11

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	364/369 (99%)	308 (85%)	56 (15%)	2 8

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1	GLN
1	H	3	GLN
1	H	7	SER
1	H	12	VAL
1	H	13	LYS
1	H	23	THR
1	H	29	ILE
1	H	45	LYS
1	H	52	VAL
1	H	65	LEU
1	H	67	SER
1	H	68	ARG
1	H	69	VAL
1	H	83	LYS
1	H	84	LEU
1	H	86	SER
1	H	87	VAL
1	H	88	THR
1	H	119	THR
1	H	121	SER
1	H	122	SER
1	H	124	SER
1	H	125	THR
1	H	126	LYS
1	H	136	SER
1	H	140	THR
1	H	141	SER
1	H	147	LEU
1	H	158	PRO
1	H	206	ASN
1	H	213	ASN
1	H	223	LYS
2	L	2	SER
2	L	6	GLN
2	L	16	GLN
2	L	18	VAL
2	L	36	SER

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Mol	Chain	Res	Type
2	L	41	LEU
2	L	47	LYS
2	L	50	ILE
2	L	52	LYS
2	L	67	SER
2	L	72	THR
2	L	78	SER
2	L	80	LEU
2	L	81	GLN
2	L	87	ASP
2	L	94	ARG
2	L	108	THR
2	L	110	LEU
2	L	112	GLN
2	L	140	ILE
2	L	153	LYS
2	L	160	LYS
2	L	193	ARG
2	L	194	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	3	GLN
1	H	5	GLN
1	H	114	GLN
1	H	180	GLN
1	H	201	GLN
1	H	208	ASN
2	L	6	GLN
2	L	71	ASN
2	L	95	ASN
2	L	112	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 ⓘ

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	H	223/223 (100%)	1.36	34 (15%) 5 4	10, 56, 76, 99	10 (4%)
2	L	211/212 (99%)	1.12	20 (9%) 14 11	22, 51, 69, 80	10 (4%)
All	All	434/435 (99%)	1.24	54 (12%) 8 6	10, 53, 74, 99	20 (4%)

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	181	SER	18.4
1	H	11	LEU	9.4
1	H	190	VAL	9.2
1	H	137	SER	9.2
2	L	167	THR	8.8
2	L	22	CYS	8.7
1	H	223	LYS	8.4
2	L	179	SER	8.0
2	L	90	CYS	7.9
2	L	126	SER	7.8
2	L	65	SER	7.2
1	H	173	HIS	7.1
1	H	186	SER	6.7
2	L	205	THR	6.3
2	L	149	THR	6.3
1	H	140	THR	6.1
1	H	141	SER	6.1
1	H	157	GLU	6.1
2	L	118	SER	6.0
2	L	58	SER	4.9
1	H	142	GLY	4.6
2	L	132	ASN	4.4
1	H	19	SER	4.4
1	H	138	LYS	4.1

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Mol	Chain	Res	Type	RSRZ
2	L	187	GLU	3.9
2	L	184	LEU	3.8
1	H	139	SER	3.4
1	H	222	PRO	3.2
1	H	17	THR	3.2
1	H	167	ALA	3.1
1	H	111	TYR	3.0
1	H	136	SER	2.9
1	H	107	PHE	2.9
1	H	56	ASP	2.9
2	L	82	ALA	2.9
1	H	196	SER	2.7
1	H	117	MET	2.5
1	H	114	GLN	2.5
1	H	86	SER	2.4
2	L	50	ILE	2.4
1	H	104	VAL	2.4
2	L	154	ALA	2.4
1	H	149	CYS	2.4
1	H	191	VAL	2.3
2	L	87	ASP	2.3
1	H	16	GLN	2.3
1	H	71	ILE	2.2
1	H	18	LEU	2.1
2	L	174	ASN	2.1
1	H	91	ASP	2.0
1	H	214	THR	2.0
2	L	194	SER	2.0
1	H	118	VAL	2.0
2	L	148	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 oligosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.