



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:58 AM UTC

PDB ID : 1FH5 / pdb_00001fh5
Title : CRYSTAL STRUCTURE OF THE FAB FRAGMENT OF THE MONO-CLONAL ANTIBODY MAK33
Authors : Augustine, J.G.; de la Calle, A.; Knarr, G.; Buchner, J.; Frederick, C.A.
Deposited on : 2000-07-31
Resolution : 2.90 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

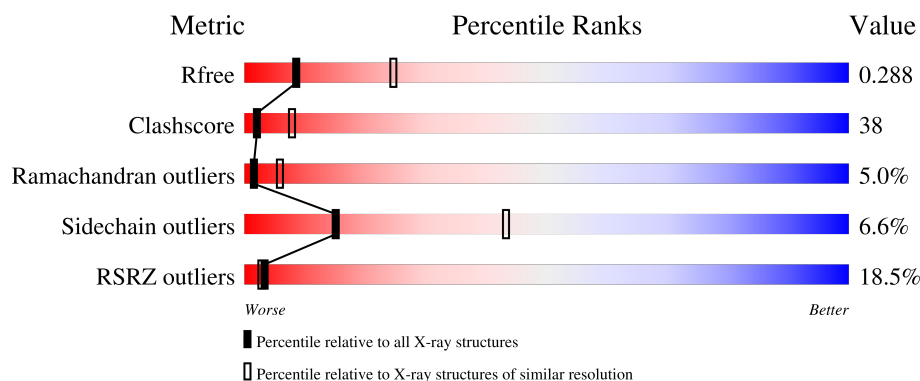
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	L	213	10% 48% 43% 8%
2	H	198	27% 57% 32% 8% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3159 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 MONOCLONAL ANTIBODY MAK33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	213	Total	C	N	O	S	0	0	0
			1650	1023	279	342	6			

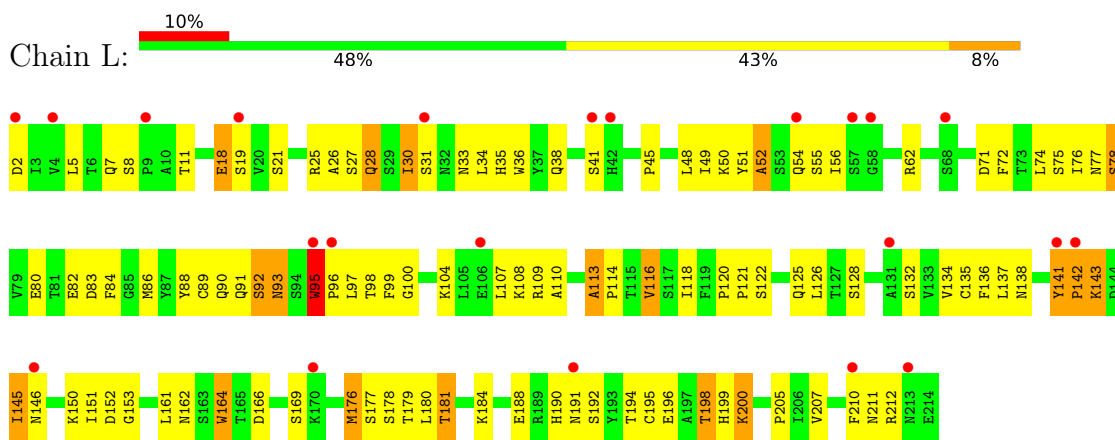
- Molecule 2 is a protein called MONOCLONAL ANTIBODY MAK33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	198	Total	C	N	O	S	0	0	0
			1509	960	245	296	8			

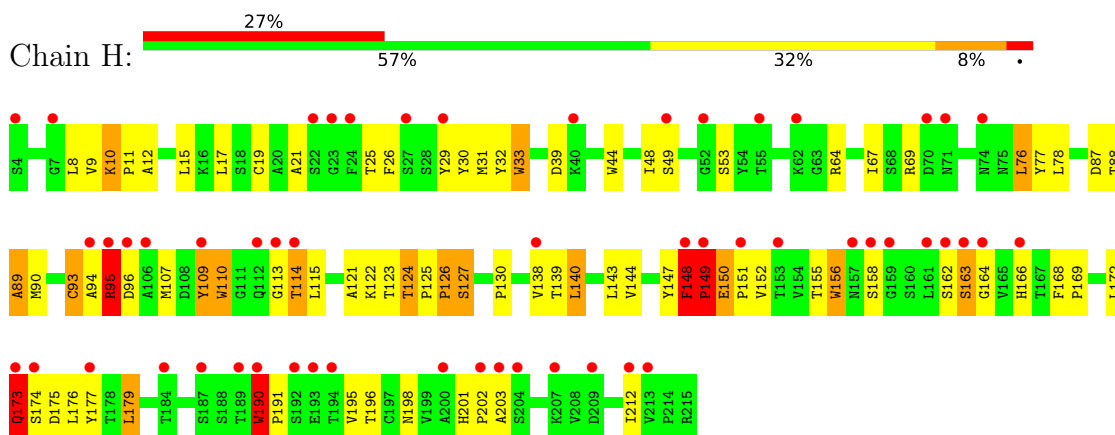
3 Residue-property plots

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: MONOCLONAL ANTIBODY MAK33



• Molecule 2: MONOCLONAL ANTIBODY MAK33



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.91Å 65.66Å 73.46Å 109.14° 106.94° 95.91°	Depositor
Resolution (Å)	8.00 – 2.90 8.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.90) 89.2 (8.00-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.88Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.198 , 0.243 0.276 , 0.288	Depositor DCC
R_{free} test set	1975 reflections (9.80%)	wwPDB-VP
Wilson B-factor (Å ²)	45.4	Xtriage
Anisotropy	0.188	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 60.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.62	EDS
Total number of atoms	3159	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.96% 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	L	0.83	5/1687 (0.3%)	1.25	8/2290 (0.3%)
2	H	0.93	4/1548 (0.3%)	1.26	9/2111 (0.4%)
All	All	0.88	9/3235 (0.3%)	1.26	17/4401 (0.4%)

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
2	H	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	126	PRO	C-N	-9.75	1.21	1.33
2	H	127	SER	C-N	-6.72	1.23	1.33
1	L	145	ILE	C-N	-5.67	1.26	1.33
2	H	190	TRP	NE1-CE2	-5.59	1.31	1.37
1	L	18	GLU	C-N	-5.47	1.25	1.33
1	L	95	TRP	NE1-CE2	-5.45	1.31	1.37
1	L	75	SER	C-N	-5.35	1.26	1.33
2	H	156	TRP	NE1-CE2	-5.26	1.31	1.37
1	L	164	TRP	NE1-CE2	-5.21	1.31	1.37

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	127	SER	CA-C-O	-8.11	110.31	120.20
1	L	113	ALA	CA-C-N	7.71	127.47	119.76
1	L	113	ALA	C-N-CA	7.71	127.47	119.76
2	H	148	PHE	CA-CB-CG	-6.96	106.84	113.80
1	L	145	ILE	CA-C-O	-6.79	113.41	121.28

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	150	GLU	O-C-N	-6.55	113.78	121.32
2	H	33	TRP	CA-C-O	-5.65	114.47	120.40
1	L	153	GLY	CA-C-N	-5.61	114.28	122.65
1	L	153	GLY	C-N-CA	-5.61	114.28	122.65
1	L	116	VAL	N-CA-C	5.59	117.64	108.97
2	H	166	HIS	CB-CA-C	-5.58	104.39	111.70
2	H	76	LEU	CA-C-O	-5.51	114.37	120.38
2	H	164	GLY	N-CA-C	-5.47	107.77	115.32
1	L	92	SER	N-CA-C	5.30	117.24	110.61
2	H	149	PRO	O-C-N	5.17	129.63	122.64
1	L	166	ASP	N-CA-C	-5.09	104.15	110.41
2	H	95	ARG	NE-CZ-NH2	5.05	123.74	119.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	127	SER	Mainchain
2	H	77	TYR	Sidechain

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	L	1650	0	1573	137	0
2	H	1509	0	1470	104	0
All	All	3159	0	3043	234	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 38.

All (234) 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:L:145:ILE:HG22	1:L:164:TRP:CH2	1.74	1.21

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:95:TRP:CG	1:L:96:PRO:HD3	1.76	1.19
1:L:95:TRP:HB3	1:L:96:PRO:CD	1.71	1.19
2:H:31:MET:HE1	2:H:110:TRP:HZ2	1.13	1.14
1:L:192:SER:HB3	1:L:211:ASN:OD1	1.50	1.10
2:H:190:TRP:HB3	2:H:191:PRO:HD3	1.32	1.10
1:L:49:ILE:HD12	1:L:74:LEU:CD1	1.81	1.09
1:L:95:TRP:HB3	1:L:96:PRO:HD2	1.35	1.08
1:L:95:TRP:CB	1:L:96:PRO:CD	2.34	1.06
1:L:196:GLU:HG3	1:L:207:VAL:HG12	1.37	1.06
1:L:145:ILE:HG22	1:L:164:TRP:HH2	0.96	1.06
2:H:148:PHE:HB3	2:H:149:PRO:HD3	1.39	1.01
2:H:190:TRP:HB3	2:H:191:PRO:CD	1.88	1.01
1:L:95:TRP:CB	1:L:96:PRO:HD3	1.92	0.95
2:H:31:MET:HE1	2:H:110:TRP:CZ2	2.02	0.93
2:H:148:PHE:HB3	2:H:149:PRO:CD	1.98	0.93
1:L:86:MET:SD	1:L:104:LYS:HD2	2.08	0.92
1:L:26:ALA:CB	1:L:30:ILE:HG21	2.00	0.92
2:H:9:VAL:HG21	2:H:15:LEU:HD22	1.54	0.90
2:H:190:TRP:CB	2:H:191:PRO:HD3	2.02	0.89
1:L:145:ILE:CG2	1:L:164:TRP:CH2	2.56	0.88
1:L:33:ASN:HD22	1:L:93:ASN:HB3	1.40	0.87
1:L:95:TRP:CD2	1:L:96:PRO:HD3	2.09	0.87
1:L:95:TRP:CG	1:L:96:PRO:CD	2.59	0.85
1:L:26:ALA:CB	1:L:30:ILE:CG2	2.55	0.84
1:L:50:LYS:O	1:L:54:GLN:HB2	1.78	0.84
1:L:99:PHE:HZ	2:H:107:MET:HE1	1.42	0.83
2:H:190:TRP:CB	2:H:191:PRO:CD	2.56	0.83
1:L:38:GLN:HB2	1:L:48:LEU:HD11	1.59	0.83
1:L:99:PHE:CZ	2:H:107:MET:HE1	2.15	0.82
2:H:31:MET:CE	2:H:110:TRP:HZ2	1.91	0.81
1:L:49:ILE:HD12	1:L:74:LEU:HD11	1.61	0.81
1:L:95:TRP:HB3	1:L:96:PRO:HD3	1.53	0.81
2:H:123:THR:HG23	2:H:149:PRO:HG2	1.62	0.81
1:L:84:PHE:CE1	1:L:107:LEU:HB2	2.17	0.79
1:L:11:THR:HG22	1:L:104:LYS:HB3	1.65	0.79
2:H:123:THR:HG23	2:H:149:PRO:CG	2.13	0.79
1:L:2:ASP:OD1	1:L:96:PRO:HG2	1.83	0.78
1:L:49:ILE:CD1	1:L:74:LEU:CD1	2.59	0.78
2:H:32:TYR:CD1	2:H:107:MET:HG2	2.18	0.78
1:L:108:LYS:HA	1:L:141:TYR:OH	1.84	0.77
2:H:19:CYS:HB3	2:H:76:LEU:HB3	1.64	0.77

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:109:ARG:O	1:L:141:TYR:CE2	2.38	0.77
1:L:51:TYR:O	1:L:52:ALA:HB3	1.84	0.76
2:H:149:PRO:C	2:H:151:PRO:HD2	2.08	0.76
1:L:109:ARG:HG2	1:L:141:TYR:CD1	2.21	0.76
1:L:164:TRP:O	2:H:169:PRO:HG2	1.86	0.76
2:H:158:SER:H	2:H:198:ASN:ND2	1.84	0.76
1:L:7:GLN:HE22	1:L:88:TYR:HA	1.51	0.76
2:H:122:LYS:O	2:H:124:THR:HG22	1.88	0.73
1:L:164:TRP:CZ2	1:L:176:MET:HE3	2.23	0.73
1:L:116:VAL:HG22	1:L:137:LEU:HD12	1.69	0.73
1:L:192:SER:CB	1:L:211:ASN:OD1	2.35	0.72
1:L:51:TYR:O	1:L:52:ALA:CB	2.37	0.72
2:H:148:PHE:CB	2:H:149:PRO:CD	2.68	0.72
1:L:77:ASN:O	1:L:78:SER:HB3	1.89	0.72
2:H:148:PHE:CB	2:H:149:PRO:HD3	2.18	0.71
2:H:10:LYS:HD3	2:H:10:LYS:H	1.56	0.71
1:L:152:ASP:OD2	1:L:190:HIS:HB3	1.91	0.71
1:L:33:ASN:ND2	1:L:93:ASN:HB3	2.06	0.70
1:L:176:MET:HG2	1:L:177:SER:N	2.06	0.70
1:L:90:GLN:NE2	1:L:97:LEU:HD11	2.05	0.70
2:H:138:VAL:HG12	2:H:140:LEU:HD23	1.74	0.69
2:H:144:VAL:HG12	2:H:147:TYR:CD2	2.29	0.68
2:H:150:GLU:O	2:H:152:VAL:HG23	1.93	0.68
2:H:158:SER:H	2:H:198:ASN:HD21	1.40	0.68
2:H:190:TRP:CD1	2:H:190:TRP:C	2.69	0.68
1:L:190:HIS:O	1:L:212:ARG:NH1	2.26	0.68
2:H:64:ARG:NH2	2:H:87:ASP:OD2	2.27	0.68
1:L:26:ALA:HB3	1:L:30:ILE:CG2	2.24	0.67
2:H:21:ALA:HA	2:H:110:TRP:HZ3	1.60	0.67
1:L:91:GLN:HE21	1:L:98:THR:H	1.41	0.67
1:L:141:TYR:HB3	1:L:142:PRO:HD3	1.76	0.67
2:H:190:TRP:CG	2:H:191:PRO:HD3	2.29	0.66
1:L:27:SER:O	1:L:28:GLN:HB2	1.96	0.66
1:L:120:PRO:HB3	1:L:210:PHE:CE1	2.30	0.66
1:L:30:ILE:HD11	1:L:34:LEU:HB2	1.78	0.66
2:H:32:TYR:CE1	2:H:96:ASP:HB2	2.31	0.66
1:L:126:LEU:HD12	1:L:184:LYS:HG3	1.78	0.66
1:L:116:VAL:HG22	1:L:137:LEU:CD1	2.26	0.65
2:H:114:THR:O	2:H:114:THR:HG23	1.95	0.65
1:L:109:ARG:HG2	1:L:141:TYR:CE1	2.31	0.65
1:L:50:LYS:O	1:L:50:LYS:HG3	1.96	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:198:THR:HG23	1:L:205:PRO:HB3	1.79	0.64
1:L:49:ILE:HD12	1:L:74:LEU:HD13	1.74	0.63
1:L:146:ASN:HB3	1:L:198:THR:OG1	1.99	0.62
2:H:9:VAL:CG2	2:H:15:LEU:HD22	2.25	0.62
2:H:138:VAL:CG1	2:H:139:THR:N	2.61	0.62
1:L:5:LEU:HD23	1:L:26:ALA:HA	1.83	0.61
1:L:121:PRO:HB2	1:L:126:LEU:HD21	1.82	0.61
2:H:140:LEU:HD12	2:H:195:VAL:HG11	1.82	0.61
2:H:21:ALA:HA	2:H:110:TRP:CZ3	2.36	0.60
2:H:138:VAL:HG12	2:H:139:THR:N	2.16	0.60
2:H:148:PHE:O	2:H:150:GLU:N	2.35	0.60
2:H:48:ILE:HB	2:H:67:ILE:HD13	1.82	0.59
1:L:26:ALA:HB1	1:L:30:ILE:CG2	2.33	0.59
1:L:142:PRO:HG2	1:L:199:HIS:CE1	2.37	0.59
1:L:109:ARG:NH1	1:L:110:ALA:O	2.36	0.59
1:L:84:PHE:CZ	1:L:107:LEU:HB2	2.38	0.58
1:L:118:ILE:CD1	1:L:210:PHE:CE2	2.86	0.58
2:H:10:LYS:H	2:H:10:LYS:CD	2.15	0.58
2:H:31:MET:HB2	2:H:48:ILE:HG22	1.86	0.58
2:H:88:THR:O	2:H:89:ALA:HB2	2.03	0.58
1:L:118:ILE:HG12	1:L:210:PHE:HE2	1.69	0.58
1:L:141:TYR:O	1:L:143:LYS:N	2.37	0.57
2:H:94:ALA:HA	2:H:110:TRP:NE1	2.19	0.57
1:L:34:LEU:HD13	1:L:72:PHE:CD2	2.39	0.57
2:H:140:LEU:HD23	2:H:140:LEU:N	2.20	0.57
1:L:49:ILE:CD1	1:L:74:LEU:HD11	2.28	0.57
2:H:148:PHE:CD1	2:H:148:PHE:C	2.80	0.56
1:L:38:GLN:HB2	1:L:48:LEU:CD1	2.33	0.56
1:L:164:TRP:CE2	1:L:176:MET:HE3	2.41	0.56
2:H:162:SER:O	2:H:163:SER:HB3	2.05	0.56
1:L:50:LYS:HG3	1:L:51:TYR:HD1	1.71	0.56
1:L:198:THR:CG2	1:L:205:PRO:HB3	2.34	0.56
1:L:122:SER:O	1:L:126:LEU:HD23	2.06	0.55
2:H:172:LEU:HD13	2:H:177:TYR:CE1	2.42	0.55
2:H:173:GLN:HG3	2:H:176:LEU:O	2.06	0.55
1:L:84:PHE:CE1	1:L:107:LEU:HD22	2.42	0.55
2:H:94:ALA:HA	2:H:110:TRP:CD1	2.41	0.55
2:H:201:HIS:CE1	2:H:203:ALA:HB3	2.43	0.54
1:L:118:ILE:HG12	1:L:210:PHE:CE2	2.43	0.54
1:L:26:ALA:HB2	1:L:30:ILE:HG21	1.85	0.54
2:H:162:SER:O	2:H:163:SER:CB	2.55	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:174:SER:OG	2:H:175:ASP:N	2.40	0.54
1:L:150:LYS:HB2	1:L:194:THR:HB	1.88	0.54
2:H:138:VAL:HG12	2:H:140:LEU:CD2	2.37	0.54
2:H:30:TYR:CZ	2:H:49:SER:HA	2.43	0.53
2:H:149:PRO:O	2:H:151:PRO:HD2	2.08	0.53
1:L:26:ALA:HB3	1:L:30:ILE:HG23	1.89	0.53
1:L:35:HIS:O	1:L:89:CYS:HA	2.08	0.53
1:L:90:GLN:HE22	1:L:97:LEU:HD11	1.73	0.53
2:H:148:PHE:O	2:H:177:TYR:CD2	2.62	0.53
1:L:98:THR:HG22	1:L:99:PHE:N	2.23	0.53
1:L:86:MET:SD	1:L:104:LYS:CD	2.91	0.52
2:H:139:THR:C	2:H:140:LEU:HD23	2.35	0.52
1:L:90:GLN:HB2	1:L:99:PHE:CD2	2.44	0.51
1:L:191:ASN:OD1	1:L:211:ASN:HB3	2.09	0.51
2:H:123:THR:HA	2:H:149:PRO:HD3	1.92	0.51
1:L:118:ILE:CD1	1:L:210:PHE:HE2	2.23	0.51
1:L:97:LEU:HD23	2:H:44:TRP:HB2	1.91	0.51
2:H:31:MET:HB2	2:H:48:ILE:CG2	2.41	0.51
2:H:179:LEU:C	2:H:179:LEU:HD23	2.36	0.51
2:H:109:TYR:C	2:H:110:TRP:CD1	2.89	0.51
1:L:109:ARG:HG3	1:L:109:ARG:HH11	1.75	0.51
1:L:191:ASN:O	1:L:211:ASN:HA	2.12	0.50
1:L:134:VAL:HG22	1:L:179:THR:HG23	1.92	0.50
1:L:49:ILE:CD1	1:L:74:LEU:HD12	2.39	0.50
1:L:90:GLN:HE21	1:L:97:LEU:CG	2.24	0.50
1:L:50:LYS:O	1:L:50:LYS:CG	2.60	0.50
1:L:98:THR:CG2	1:L:99:PHE:N	2.75	0.49
1:L:116:VAL:HG13	1:L:137:LEU:HD13	1.94	0.49
1:L:118:ILE:HD13	1:L:210:PHE:CD2	2.46	0.49
2:H:190:TRP:CG	2:H:191:PRO:CD	2.93	0.49
1:L:164:TRP:O	2:H:169:PRO:CG	2.59	0.49
2:H:17:LEU:HD12	2:H:78:LEU:HD23	1.95	0.49
1:L:191:ASN:O	1:L:212:ARG:N	2.43	0.49
2:H:10:LYS:HD3	2:H:10:LYS:N	2.25	0.49
2:H:147:TYR:CE1	2:H:177:TYR:HB2	2.48	0.48
2:H:19:CYS:HB2	2:H:33:TRP:CH2	2.47	0.48
1:L:132:SER:OG	1:L:181:THR:HG23	2.12	0.48
1:L:151:ILE:HD11	1:L:180:LEU:HD21	1.96	0.48
1:L:35:HIS:CD2	1:L:51:TYR:H	2.32	0.48
1:L:26:ALA:HB1	1:L:30:ILE:HG21	1.91	0.47
1:L:162:ASN:HD22	1:L:178:SER:HA	1.78	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:200:LYS:NZ	1:L:200:LYS:H	2.12	0.47
2:H:109:TYR:C	2:H:110:TRP:HD1	2.23	0.47
2:H:21:ALA:HB2	2:H:110:TRP:CH2	2.48	0.47
2:H:144:VAL:HG12	2:H:147:TYR:HD2	1.79	0.47
2:H:156:TRP:HA	2:H:196:THR:O	2.14	0.47
2:H:190:TRP:CG	2:H:191:PRO:N	2.80	0.47
2:H:149:PRO:C	2:H:151:PRO:CD	2.84	0.47
1:L:30:ILE:CD1	1:L:34:LEU:HD12	2.45	0.47
1:L:118:ILE:CG1	1:L:210:PHE:HE2	2.27	0.47
2:H:31:MET:HG3	2:H:76:LEU:HD13	1.96	0.47
2:H:130:PRO:CB	2:H:212:ILE:HD13	2.44	0.47
1:L:36:TRP:CE2	1:L:74:LEU:HB2	2.50	0.47
2:H:144:VAL:HG12	2:H:147:TYR:CE2	2.50	0.47
2:H:148:PHE:O	2:H:149:PRO:C	2.58	0.46
1:L:84:PHE:HE1	1:L:107:LEU:HD22	1.80	0.46
2:H:148:PHE:O	2:H:177:TYR:HD2	1.97	0.46
1:L:141:TYR:HB3	1:L:142:PRO:CD	2.43	0.46
2:H:11:PRO:O	2:H:12:ALA:HB3	2.15	0.46
1:L:80:GLU:O	1:L:83:ASP:HB2	2.16	0.46
1:L:18:GLU:HA	1:L:18:GLU:OE1	2.16	0.45
2:H:32:TYR:O	2:H:93:CYS:HA	2.16	0.45
1:L:125:GLN:O	1:L:128:SER:HB2	2.16	0.45
2:H:113:GLY:O	2:H:114:THR:HB	2.17	0.45
1:L:90:GLN:HE21	1:L:97:LEU:CD1	2.30	0.45
1:L:99:PHE:HZ	2:H:107:MET:CE	2.22	0.45
2:H:90:MET:HG2	2:H:115:LEU:HD13	1.99	0.45
1:L:62:ARG:O	1:L:76:ILE:HA	2.17	0.45
2:H:190:TRP:HB3	2:H:191:PRO:HD2	1.92	0.45
1:L:90:GLN:HE21	1:L:97:LEU:HG	1.81	0.44
2:H:125:PRO:HA	2:H:126:PRO:HD3	1.76	0.44
1:L:50:LYS:O	1:L:51:TYR:HB2	2.17	0.44
1:L:77:ASN:O	1:L:78:SER:CB	2.57	0.44
1:L:35:HIS:HE1	1:L:92:SER:OG	2.01	0.44
2:H:114:THR:O	2:H:114:THR:CG2	2.63	0.43
2:H:155:THR:O	2:H:198:ASN:ND2	2.51	0.43
1:L:19:SER:HB2	1:L:77:ASN:HA	1.99	0.43
1:L:135:CYS:CB	1:L:195:CYS:SG	3.06	0.43
1:L:162:ASN:ND2	1:L:178:SER:OG	2.51	0.43
1:L:25:ARG:HE	1:L:25:ARG:HB2	1.69	0.43
2:H:168:PHE:O	2:H:179:LEU:HG	2.19	0.43
1:L:142:PRO:HD2	1:L:199:HIS:CE1	2.54	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:30:TYR:CE1	2:H:49:SER:HA	2.54	0.43
1:L:188:GLU:O	1:L:212:ARG:NH1	2.51	0.43
1:L:45:PRO:HD2	2:H:109:TYR:CE1	2.54	0.43
1:L:121:PRO:HB2	1:L:126:LEU:CD2	2.48	0.43
1:L:62:ARG:NH1	1:L:83:ASP:OD1	2.52	0.42
1:L:118:ILE:HD11	1:L:210:PHE:CE2	2.54	0.42
1:L:136:PHE:HB3	1:L:138:ASN:OD1	2.19	0.42
1:L:90:GLN:NE2	1:L:97:LEU:CD1	2.78	0.42
2:H:29:TYR:O	2:H:69:ARG:NH2	2.43	0.42
2:H:126:PRO:HB3	2:H:144:VAL:HG13	2.02	0.42
1:L:113:ALA:HA	1:L:114:PRO:HD3	1.75	0.41
2:H:30:TYR:CE2	2:H:49:SER:HA	2.56	0.41
2:H:121:ALA:HB3	2:H:148:PHE:CE2	2.55	0.41
1:L:126:LEU:C	1:L:128:SER:N	2.77	0.41
2:H:29:TYR:HD1	2:H:95:ARG:HB2	1.85	0.41
2:H:88:THR:O	2:H:89:ALA:CB	2.69	0.41
2:H:173:GLN:CG	2:H:176:LEU:O	2.66	0.41
2:H:201:HIS:O	2:H:202:PRO:C	2.64	0.41
1:L:192:SER:HA	1:L:211:ASN:HA	2.02	0.41
2:H:94:ALA:HA	2:H:110:TRP:HE1	1.84	0.41
2:H:190:TRP:CD2	2:H:191:PRO:HD3	2.56	0.41
1:L:5:LEU:CD2	1:L:26:ALA:HB2	2.50	0.41
2:H:150:GLU:O	2:H:151:PRO:C	2.61	0.41
1:L:7:GLN:HE21	1:L:100:GLY:HA3	1.85	0.40
1:L:56:ILE:N	1:L:56:ILE:HD13	2.35	0.40
1:L:188:GLU:C	1:L:212:ARG:NH1	2.79	0.40
2:H:11:PRO:O	2:H:12:ALA:CB	2.70	0.40
1:L:2:ASP:CG	1:L:96:PRO:HG2	2.45	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	L	211/213 (99%)	187 (89%)	14 (7%)	10 (5%)	2	7
2	H	192/198 (97%)	166 (86%)	16 (8%)	10 (5%)	1	5
All	All	403/411 (98%)	353 (88%)	30 (7%)	20 (5%)	1	6

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	52	ALA
1	L	95	TRP
2	H	110	TRP
2	H	114	THR
2	H	163	SER
2	H	173	GLN
2	H	190	TRP
1	L	28	GLN
1	L	143	LYS
2	H	26	PHE
2	H	148	PHE
1	L	31	SER
1	L	169	SER
2	H	109	TYR
1	L	78	SER
1	L	141	TYR
2	H	89	ALA
1	L	82	GLU
2	H	149	PRO
1	L	142	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	L	192/192 (100%)	180 (94%)	12 (6%)	16	45
2	H	170/170 (100%)	158 (93%)	12 (7%)	13	40
All	All	362/362 (100%)	338 (93%)	24 (7%)	15	43

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	8	SER
1	L	21	SER
1	L	30	ILE
1	L	41	SER
1	L	55	SER
1	L	71	ASP
1	L	93	ASN
1	L	161	LEU
1	L	176	MET
1	L	181	THR
1	L	198	THR
1	L	200	LYS
2	H	8	LEU
2	H	10	LYS
2	H	25	THR
2	H	39	ASP
2	H	53	SER
2	H	93	CYS
2	H	95	ARG
2	H	124	THR
2	H	140	LEU
2	H	143	LEU
2	H	173	GLN
2	H	179	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	7	GLN
1	L	28	GLN
1	L	33	ASN
1	L	35	HIS
1	L	38	GLN
1	L	90	GLN
1	L	93	ASN
1	L	162	ASN
2	H	198	ASN

5.3.3 RNA [i](#)

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	H	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	132:ALA	C	138:VAL	N	10.61
1	H	96:ASP	C	106:ALA	N	6.39

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	L	213/213 (100%)	0.89	22 (10%)	12 10	8, 32, 56, 73	0
2	H	198/198 (100%)	1.27	54 (27%)	1 1	13, 39, 69, 91	0
All	All	411/411 (100%)	1.07	76 (18%)	3 3	8, 35, 62, 91	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	22	SER	8.6
1	L	213	ASN	7.1
2	H	193	GLU	4.9
2	H	159	GLY	4.7
2	H	189	THR	4.6
2	H	74	ASN	4.5
2	H	173	GLN	4.3
2	H	96	ASP	4.1
1	L	141	TYR	4.0
2	H	158	SER	3.7
2	H	113	GLY	3.7
1	L	95	TRP	3.5
2	H	157	ASN	3.5
2	H	209	ASP	3.4
2	H	151	PRO	3.4
2	H	29	TYR	3.4
2	H	187	SER	3.3
2	H	162	SER	3.2
2	H	27	SER	3.2
2	H	24	PHE	3.1
2	H	194	THR	3.1
2	H	200	ALA	3.0
2	H	23	GLY	3.0
2	H	138	VAL	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	H	71	ASN	2.9
1	L	106	GLU	2.8
2	H	203	ALA	2.8
1	L	42	HIS	2.8
2	H	204	SER	2.8
2	H	7	GLY	2.8
2	H	177	TYR	2.8
1	L	19	SER	2.8
1	L	58	GLY	2.8
1	L	146	ASN	2.8
2	H	95	ARG	2.7
2	H	153	THR	2.7
2	H	52	GLY	2.7
1	L	9	PRO	2.7
2	H	207	LYS	2.7
1	L	31	SER	2.7
1	L	68	SER	2.7
2	H	49	SER	2.6
2	H	163	SER	2.6
1	L	142	PRO	2.6
2	H	149	PRO	2.5
2	H	94	ALA	2.5
2	H	40	LYS	2.5
2	H	161	LEU	2.4
1	L	191	ASN	2.4
2	H	213	VAL	2.4
2	H	174	SER	2.4
2	H	106	ALA	2.4
1	L	2	ASP	2.4
2	H	190	TRP	2.4
2	H	55	THR	2.4
1	L	210	PHE	2.3
1	L	41	SER	2.3
2	H	70	ASP	2.3
1	L	57	SER	2.3
1	L	96	PRO	2.2
2	H	184	THR	2.2
2	H	202	PRO	2.2
2	H	62	LYS	2.2
2	H	112	GLN	2.2
2	H	4	SER	2.2
2	H	164	GLY	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	L	170	LYS	2.2
2	H	212	ILE	2.2
2	H	114	THR	2.2
2	H	192	SER	2.2
1	L	131	ALA	2.2
2	H	166	HIS	2.1
1	L	4	VAL	2.1
1	L	54	GLN	2.0
2	H	109	TYR	2.0
2	H	148	PHE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.