



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2018 – 06:33 AM EST

PDB ID : 4L6W
Title : Carboxyltransferase subunit (AccD6) of Mycobacterium tuberculosis acetyl-CoA carboxylase
Authors : Anandhakrishnan, M.; Ehebauer, M.T.; Wilmanns, M.
Deposited on : 2013-06-13
Resolution : 1.95 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
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030736

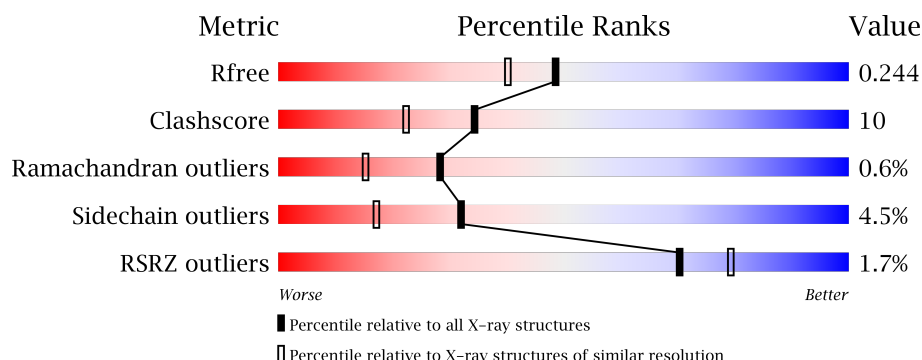
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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}	100719	2004 (1.96-1.96)
Clashscore	112137	2136 (1.96-1.96)
Ramachandran outliers	110173	2117 (1.96-1.96)
Sidechain outliers	110143	2117 (1.96-1.96)
RSRZ outliers	101464	2018 (1.96-1.96)

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	493	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 71%, yellow 12%, grey 16%);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> 1% 71% 12% • 16% </div> </div>
1	B	493	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, green 71%, yellow 12%, grey 15%);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> 2% 71% 12% • 15% </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6336 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 Probable propionyl-CoA carboxylase beta chain 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	412	Total	C	N	O	S	0	0	0
			3032	1901	541	575	15			
1	B	419	Total	C	N	O	S	0	0	0
			3109	1948	568	579	14			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP P63407
A	-18	GLY	-	EXPRESSION TAG	UNP P63407
A	-17	SER	-	EXPRESSION TAG	UNP P63407
A	-16	SER	-	EXPRESSION TAG	UNP P63407
A	-15	HIS	-	EXPRESSION TAG	UNP P63407
A	-14	HIS	-	EXPRESSION TAG	UNP P63407
A	-13	HIS	-	EXPRESSION TAG	UNP P63407
A	-12	HIS	-	EXPRESSION TAG	UNP P63407
A	-11	HIS	-	EXPRESSION TAG	UNP P63407
A	-10	HIS	-	EXPRESSION TAG	UNP P63407
A	-9	SER	-	EXPRESSION TAG	UNP P63407
A	-8	SER	-	EXPRESSION TAG	UNP P63407
A	-7	GLY	-	EXPRESSION TAG	UNP P63407
A	-6	LEU	-	EXPRESSION TAG	UNP P63407
A	-5	VAL	-	EXPRESSION TAG	UNP P63407
A	-4	PRO	-	EXPRESSION TAG	UNP P63407
A	-3	ARG	-	EXPRESSION TAG	UNP P63407
A	-2	GLY	-	EXPRESSION TAG	UNP P63407
A	-1	SER	-	EXPRESSION TAG	UNP P63407
A	0	HIS	-	EXPRESSION TAG	UNP P63407
B	-19	MET	-	EXPRESSION TAG	UNP P63407
B	-18	GLY	-	EXPRESSION TAG	UNP P63407
B	-17	SER	-	EXPRESSION TAG	UNP P63407
B	-16	SER	-	EXPRESSION TAG	UNP P63407
B	-15	HIS	-	EXPRESSION TAG	UNP P63407

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	EXPRESSION TAG	UNP P63407
B	-13	HIS	-	EXPRESSION TAG	UNP P63407
B	-12	HIS	-	EXPRESSION TAG	UNP P63407
B	-11	HIS	-	EXPRESSION TAG	UNP P63407
B	-10	HIS	-	EXPRESSION TAG	UNP P63407
B	-9	SER	-	EXPRESSION TAG	UNP P63407
B	-8	SER	-	EXPRESSION TAG	UNP P63407
B	-7	GLY	-	EXPRESSION TAG	UNP P63407
B	-6	LEU	-	EXPRESSION TAG	UNP P63407
B	-5	VAL	-	EXPRESSION TAG	UNP P63407
B	-4	PRO	-	EXPRESSION TAG	UNP P63407
B	-3	ARG	-	EXPRESSION TAG	UNP P63407
B	-2	GLY	-	EXPRESSION TAG	UNP P63407
B	-1	SER	-	EXPRESSION TAG	UNP P63407
B	0	HIS	-	EXPRESSION TAG	UNP P63407

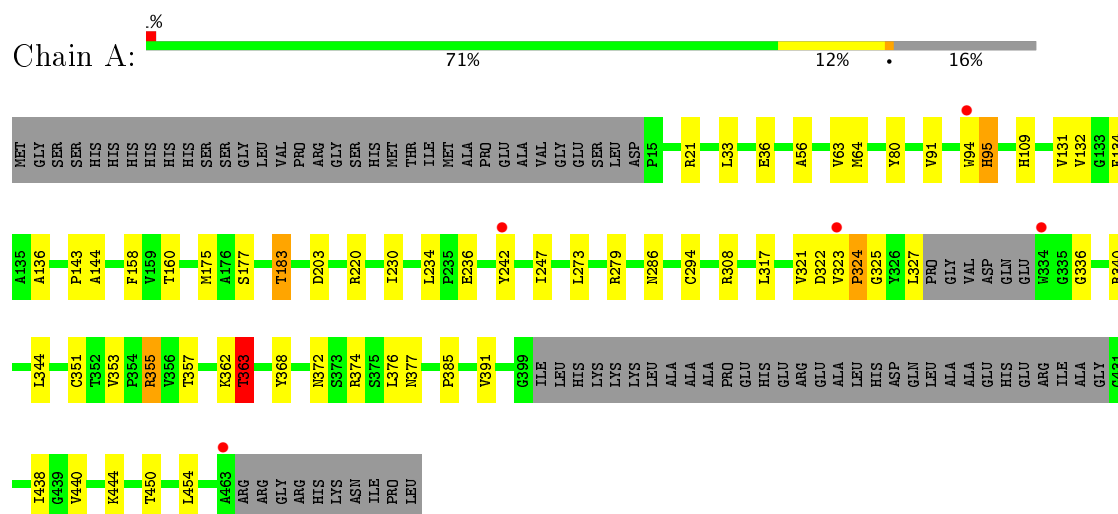
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	91	Total O 91 91	0	0
2	B	104	Total O 104 104	0	0

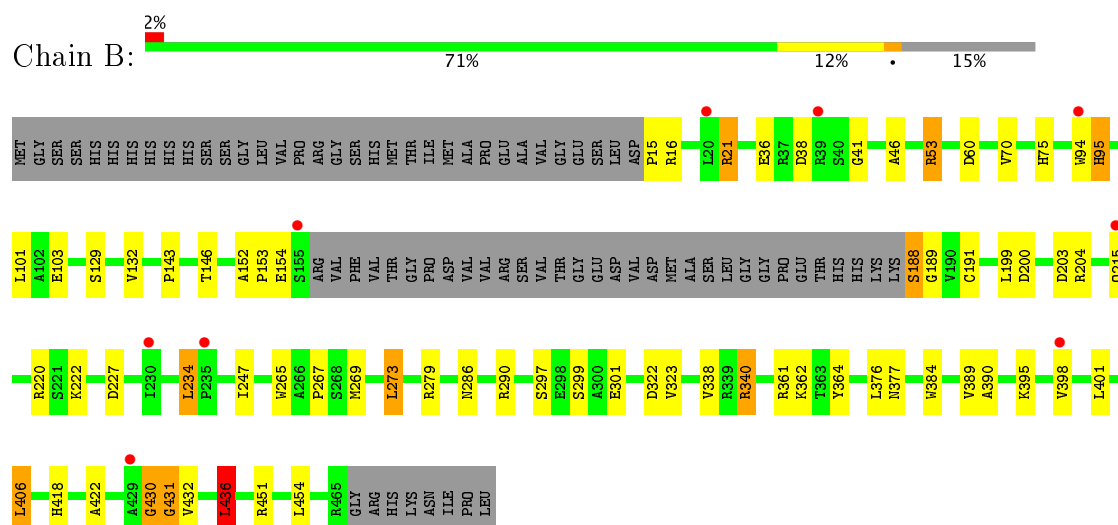
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Probable propionyl-CoA carboxylase beta chain 6



- Molecule 1: Probable propionyl-CoA carboxylase beta chain 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.94Å 90.31Å 154.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.97 – 1.95 68.88 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (68.97-1.95) 99.7 (68.88-1.95)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.11 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.174 , 0.244 0.183 , 0.244	Depositor DCC
R_{free} test set	3967 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.374	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.6	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	0.96	EDS
Total number of atoms	6336	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% 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	A	0.88	2/3089 (0.1%)	0.90	2/4201 (0.0%)
1	B	0.86	1/3168 (0.0%)	0.88	6/4303 (0.1%)
All	All	0.87	3/6257 (0.0%)	0.89	8/8504 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	153	PRO	N-CD	5.17	1.55	1.47
1	A	80	TYR	CE1-CZ	5.16	1.45	1.38
1	A	324	PRO	N-CD	5.07	1.54	1.47

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	355	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	B	436	LEU	CA-CB-CG	5.80	128.65	115.30
1	B	21	ARG	NE-CZ-NH2	-5.71	117.45	120.30
1	A	363	THR	CB-CA-C	-5.65	96.33	111.60
1	B	273	LEU	CB-CG-CD1	5.61	120.53	111.00
1	B	152	ALA	C-N-CD	5.51	139.97	128.40
1	B	273	LEU	CA-CB-CG	5.46	127.85	115.30
1	B	406	LEU	CB-CG-CD1	5.45	120.27	111.00

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	3032	0	2987	53	0
1	B	3109	0	3085	65	0
2	A	91	0	0	0	0
2	B	104	0	0	2	0
All	All	6336	0	6072	116	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 10.

All (116) 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:94:TRP:NE1	1:A:143:PRO:HG3	1.43	1.31
1:B:94:TRP:NE1	1:B:143:PRO:HG3	1.80	0.97
1:A:94:TRP:HE1	1:A:143:PRO:HG3	1.26	0.96
1:B:430:GLY:O	1:B:431:GLY:O	1.85	0.95
1:B:21:ARG:HH12	1:B:95:HIS:CD2	1.85	0.94
1:A:94:TRP:NE1	1:A:143:PRO:CG	2.32	0.93
1:B:188:SER:N	1:B:189:GLY:HA2	1.84	0.92
1:A:94:TRP:CD1	1:A:143:PRO:HG3	2.11	0.84
1:A:234:LEU:CD2	1:A:242:TYR:CZ	2.64	0.81
1:A:177:SER:O	1:A:183:THR:HG21	1.82	0.79
1:B:21:ARG:HH12	1:B:95:HIS:HD2	1.26	0.79
1:B:398:VAL:HG23	1:B:418:HIS:CE1	2.21	0.76
1:A:323:VAL:HG22	1:A:325:GLY:H	1.54	0.73
1:A:21:ARG:HH12	1:A:95:HIS:CD2	2.07	0.73
1:A:94:TRP:CZ2	1:A:143:PRO:HD3	2.24	0.72
1:A:234:LEU:HD22	1:A:242:TYR:CZ	2.25	0.71
1:B:301:GLU:HG3	1:B:340:ARG:HG2	1.70	0.71
1:A:21:ARG:HH12	1:A:95:HIS:HD2	1.39	0.70
1:B:215:GLN:OE1	1:B:279:ARG:NH1	2.26	0.69
1:B:297:SER:CB	1:B:340:ARG:HD3	2.25	0.66
1:B:94:TRP:O	1:B:95:HIS:HB2	1.96	0.65
1:B:430:GLY:O	1:B:431:GLY:C	2.34	0.65
1:A:94:TRP:O	1:A:95:HIS:HB2	1.97	0.65
1:B:15:PRO:HG2	1:B:16:ARG:H	1.61	0.65
1:B:322:ASP:CG	1:B:361:ARG:HB2	2.17	0.64
1:B:322:ASP:OD1	1:B:361:ARG:HB2	1.98	0.63
1:B:199:LEU:H	1:B:199:LEU:HD12	1.63	0.63
1:A:363:THR:HG22	1:A:368:TYR:HA	1.81	0.62
1:A:286:ASN:HD22	1:A:322:ASP:H	1.47	0.61
1:B:398:VAL:HG11	1:B:422:ALA:HA	1.81	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:376:LEU:O	1:B:377:ASN:HB2	2.01	0.60
1:A:94:TRP:CE2	1:A:143:PRO:HD3	2.37	0.60
1:B:398:VAL:CG2	1:B:418:HIS:CE1	2.84	0.60
1:B:21:ARG:NH1	1:B:95:HIS:HD2	1.98	0.59
1:B:398:VAL:HG23	1:B:418:HIS:HE1	1.67	0.59
1:A:323:VAL:HG23	1:A:324:PRO:HD2	1.84	0.59
1:B:129:SER:OG	1:B:146:THR:HG21	2.03	0.58
1:A:94:TRP:O	1:A:132:VAL:HG22	2.03	0.58
1:B:297:SER:OG	1:B:340:ARG:HD3	2.02	0.58
1:A:234:LEU:HD23	1:A:242:TYR:CE2	2.39	0.58
1:A:294:CYS:HA	1:A:323:VAL:HG23	1.86	0.57
1:A:158:PHE:HZ	1:A:175:MET:HE1	1.69	0.57
1:B:395:LYS:O	1:B:398:VAL:HG22	2.04	0.57
1:A:363:THR:CG2	1:A:368:TYR:HA	2.34	0.57
1:B:70:VAL:HG23	1:B:103:GLU:OE1	2.03	0.57
1:B:94:TRP:CE2	1:B:143:PRO:HG3	2.38	0.57
1:B:297:SER:HB3	1:B:340:ARG:HD3	1.86	0.57
1:B:389:VAL:H	1:B:431:GLY:HA3	1.71	0.56
1:B:199:LEU:HD12	1:B:199:LEU:N	2.21	0.56
1:A:234:LEU:CD2	1:A:242:TYR:OH	2.55	0.55
1:A:385:PRO:HA	1:A:444:LYS:HE3	1.87	0.55
1:A:336:GLY:O	1:A:340:ARG:HG2	2.07	0.55
1:A:321:VAL:CG1	1:A:363:THR:HG23	2.37	0.54
1:B:430:GLY:C	1:B:431:GLY:O	2.41	0.54
1:A:308:ARG:HH11	1:A:308:ARG:HG2	1.73	0.54
1:A:351:CYS:SG	1:A:353:VAL:HG22	2.47	0.54
1:B:234:LEU:HD22	1:B:384:TRP:HZ3	1.73	0.53
1:A:376:LEU:O	1:A:377:ASN:HB2	2.09	0.53
1:B:188:SER:N	1:B:189:GLY:CA	2.67	0.52
1:B:395:LYS:O	1:B:418:HIS:HE1	1.93	0.52
1:B:301:GLU:CG	1:B:340:ARG:HG2	2.38	0.52
1:B:38:ASP:OD2	1:B:75:HIS:HE1	1.93	0.51
1:B:297:SER:HB3	1:B:340:ARG:NE	2.26	0.51
1:B:297:SER:HB3	1:B:340:ARG:CZ	2.41	0.50
1:A:391:VAL:HG11	1:B:101:LEU:HD12	1.94	0.50
1:B:36:GLU:HG2	2:B:528:HOH:O	2.12	0.49
1:A:323:VAL:HG22	1:A:324:PRO:N	2.28	0.49
1:A:94:TRP:CD1	1:A:143:PRO:CG	2.87	0.49
1:B:297:SER:HB3	1:B:340:ARG:CD	2.43	0.48
1:B:199:LEU:H	1:B:199:LEU:CD1	2.26	0.48
1:A:94:TRP:HB2	1:A:131:VAL:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:94:TRP:O	1:B:132:VAL:HG22	2.12	0.48
1:B:267:PRO:HB2	1:B:290:ARG:HG3	1.94	0.48
1:B:227:ASP:OD1	1:B:451:ARG:HB3	2.14	0.48
1:A:158:PHE:HZ	1:A:175:MET:CE	2.26	0.47
1:B:398:VAL:HG11	1:B:422:ALA:CA	2.45	0.47
1:A:308:ARG:NH1	1:A:308:ARG:HG2	2.29	0.47
1:B:398:VAL:CG2	1:B:418:HIS:HE1	2.25	0.47
1:B:395:LYS:O	1:B:398:VAL:CG2	2.62	0.47
1:A:234:LEU:HD21	1:A:242:TYR:OH	2.16	0.46
1:B:94:TRP:CZ2	1:B:143:PRO:HD3	2.51	0.46
1:A:63:VAL:HG12	1:A:64:MET:HG3	1.98	0.46
1:B:220:ARG:HB3	1:B:220:ARG:CZ	2.45	0.45
1:A:234:LEU:CD2	1:A:242:TYR:CE2	2.98	0.45
1:A:317:LEU:HB2	1:A:355:ARG:HB2	1.98	0.44
1:B:323:VAL:O	1:B:323:VAL:HG13	2.16	0.44
1:A:323:VAL:HG23	1:A:324:PRO:CD	2.46	0.44
1:B:364:TYR:HA	1:B:390:ALA:O	2.17	0.44
1:B:432:VAL:HG23	1:B:436:LEU:HD22	2.00	0.44
1:B:389:VAL:HG23	1:B:432:VAL:HG12	1.99	0.44
1:A:94:TRP:CE2	1:A:143:PRO:CG	2.98	0.44
1:B:362:LYS:HB2	1:B:362:LYS:HE2	1.80	0.44
1:A:109:HIS:HD2	2:B:509:HOH:O	2.00	0.43
1:B:143:PRO:O	1:B:146:THR:OG1	2.26	0.43
1:B:269:MET:HB2	1:B:299:SER:HB2	2.01	0.43
1:A:134:PHE:HE1	1:A:175:MET:SD	2.41	0.43
1:B:94:TRP:HE1	1:B:143:PRO:HG3	1.76	0.43
1:A:136:ALA:HA	1:A:158:PHE:O	2.18	0.43
1:B:234:LEU:HD13	1:B:247:ILE:HD11	2.00	0.43
1:B:189:GLY:HA3	1:B:265:TRP:CE3	2.53	0.43
1:B:286:ASN:HD22	1:B:322:ASP:H	1.66	0.43
1:B:94:TRP:CD1	1:B:143:PRO:HG3	2.49	0.42
1:B:41:GLY:HA2	1:B:60:ASP:HB3	2.01	0.42
1:A:438:ILE:HG22	1:A:440:VAL:HG13	2.00	0.42
1:B:15:PRO:CG	1:B:16:ARG:H	2.30	0.42
1:B:200:ASP:O	1:B:204:ARG:HG3	2.19	0.42
1:A:321:VAL:HG11	1:A:363:THR:HG23	2.01	0.42
1:A:323:VAL:CG2	1:A:324:PRO:CD	2.97	0.42
1:A:323:VAL:CG2	1:A:324:PRO:N	2.83	0.41
1:A:357:THR:OG1	1:A:372:ASN:HB3	2.20	0.41
1:A:144:ALA:HB2	1:B:338:VAL:HG12	2.01	0.41
1:B:46:ALA:HB1	1:B:53:ARG:NH2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:VAL:CG2	1:A:324:PRO:HD2	2.48	0.41
1:A:230:ILE:HG13	1:A:450:THR:HB	2.03	0.40
1:A:56:ALA:HA	1:A:91:VAL:O	2.20	0.40
1:A:94:TRP:CE2	1:A:143:PRO:CD	3.02	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	A	406/493 (82%)	392 (97%)	13 (3%)	1 (0%)	51	41
1	B	415/493 (84%)	402 (97%)	9 (2%)	4 (1%)	18	7
All	All	821/986 (83%)	794 (97%)	22 (3%)	5 (1%)	28	15

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	95	HIS
1	B	95	HIS
1	B	431	GLY
1	B	154	GLU
1	B	430	GLY

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	A	311/379 (82%)	295 (95%)	16 (5%)	28	13
1	B	314/379 (83%)	302 (96%)	12 (4%)	38	24
All	All	625/758 (82%)	597 (96%)	28 (4%)	32	17

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	36	GLU
1	A	160	THR
1	A	183	THR
1	A	203	ASP
1	A	220	ARG
1	A	236	GLU
1	A	247	ILE
1	A	273	LEU
1	A	279	ARG
1	A	327	LEU
1	A	344	LEU
1	A	362	LYS
1	A	363	THR
1	A	374	ARG
1	A	454	LEU
1	B	53	ARG
1	B	188	SER
1	B	191	CYS
1	B	203	ASP
1	B	222	LYS
1	B	234	LEU
1	B	273	LEU
1	B	340	ARG
1	B	401	LEU
1	B	406	LEU
1	B	436	LEU
1	B	454	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	HIS
1	A	109	HIS

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Mol	Chain	Res	Type
1	A	185	HIS
1	A	264	ASN
1	A	286	ASN
1	A	377	ASN
1	B	75	HIS
1	B	95	HIS
1	B	286	ASN
1	B	377	ASN
1	B	418	HIS

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 carbohydrates 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	412/493 (83%)	0.18	5 (1%) 79 85	28, 41, 60, 78	0
1	B	419/493 (84%)	0.27	9 (2%) 64 73	28, 43, 67, 92	0
All	All	831/986 (84%)	0.23	14 (1%) 70 79	28, 42, 63, 92	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	94	TRP	7.2
1	B	94	TRP	4.3
1	A	334	TRP	4.1
1	A	242	TYR	3.2
1	A	323	VAL	3.2
1	B	429	ALA	3.0
1	B	39	ARG	2.6
1	B	155	SER	2.4
1	B	235	PRO	2.4
1	B	398	VAL	2.2
1	A	463	ALA	2.1
1	B	230	ILE	2.1
1	B	215	GLN	2.0
1	B	20	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.