



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:40 PM UTC

PDB ID : 2AJ9 / pdb_00002aj9
Title : X-ray crystal structure of W42A,R161A double mutant of Mycobacterium tuberculosis beta-ketoacyl-ACP synthase III
Authors : Brown, A.K.; Sridharan, S.; Kremer, L.; Lindenberg, S.; Dover, L.G.; Sacchetti, J.C.; Besra, G.S.
Deposited on : 2005-08-01
Resolution : 2.50 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

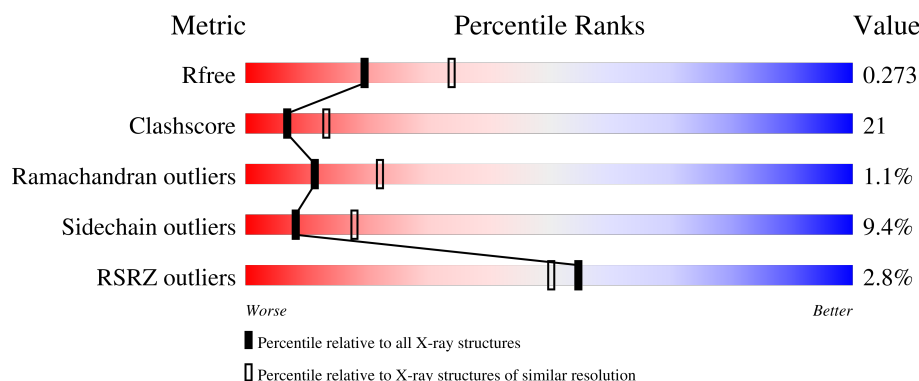
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

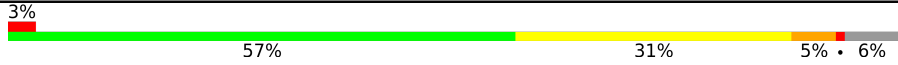
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	356	
1	B	356	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4955 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 3-oxoacyl-[acyl-carrier-protein] synthase III.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	0	0
			2427	1513	428	472	14			
1	B	334	Total	C	N	O	S	0	0	0
			2427	1513	428	472	14			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	cloning artifact	UNP P0A574
A	-19	GLY	-	cloning artifact	UNP P0A574
A	-18	SER	-	cloning artifact	UNP P0A574
A	-17	SER	-	cloning artifact	UNP P0A574
A	-16	HIS	-	expression tag	UNP P0A574
A	-15	HIS	-	expression tag	UNP P0A574
A	-14	HIS	-	expression tag	UNP P0A574
A	-13	HIS	-	expression tag	UNP P0A574
A	-12	HIS	-	expression tag	UNP P0A574
A	-11	HIS	-	expression tag	UNP P0A574
A	-10	SER	-	cloning artifact	UNP P0A574
A	-9	SER	-	cloning artifact	UNP P0A574
A	-8	GLY	-	cloning artifact	UNP P0A574
A	-7	LEU	-	cloning artifact	UNP P0A574
A	-6	VAL	-	cloning artifact	UNP P0A574
A	-5	PRO	-	cloning artifact	UNP P0A574
A	-4	ARG	-	cloning artifact	UNP P0A574
A	-3	GLY	-	cloning artifact	UNP P0A574
A	-2	SER	-	cloning artifact	UNP P0A574
A	-1	VAL	-	cloning artifact	UNP P0A574
A	0	CYS	-	cloning artifact	UNP P0A574
A	42	ALA	TRP	engineered mutation	UNP P0A574
A	161	ALA	ARG	engineered mutation	UNP P0A574
B	-20	MET	-	cloning artifact	UNP P0A574
B	-19	GLY	-	cloning artifact	UNP P0A574

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	SER	-	cloning artifact	UNP P0A574
B	-17	SER	-	cloning artifact	UNP P0A574
B	-16	HIS	-	expression tag	UNP P0A574
B	-15	HIS	-	expression tag	UNP P0A574
B	-14	HIS	-	expression tag	UNP P0A574
B	-13	HIS	-	expression tag	UNP P0A574
B	-12	HIS	-	expression tag	UNP P0A574
B	-11	HIS	-	expression tag	UNP P0A574
B	-10	SER	-	cloning artifact	UNP P0A574
B	-9	SER	-	cloning artifact	UNP P0A574
B	-8	GLY	-	cloning artifact	UNP P0A574
B	-7	LEU	-	cloning artifact	UNP P0A574
B	-6	VAL	-	cloning artifact	UNP P0A574
B	-5	PRO	-	cloning artifact	UNP P0A574
B	-4	ARG	-	cloning artifact	UNP P0A574
B	-3	GLY	-	cloning artifact	UNP P0A574
B	-2	SER	-	cloning artifact	UNP P0A574
B	-1	VAL	-	cloning artifact	UNP P0A574
B	0	CYS	-	cloning artifact	UNP P0A574
B	42	ALA	TRP	engineered mutation	UNP P0A574
B	161	ALA	ARG	engineered mutation	UNP P0A574

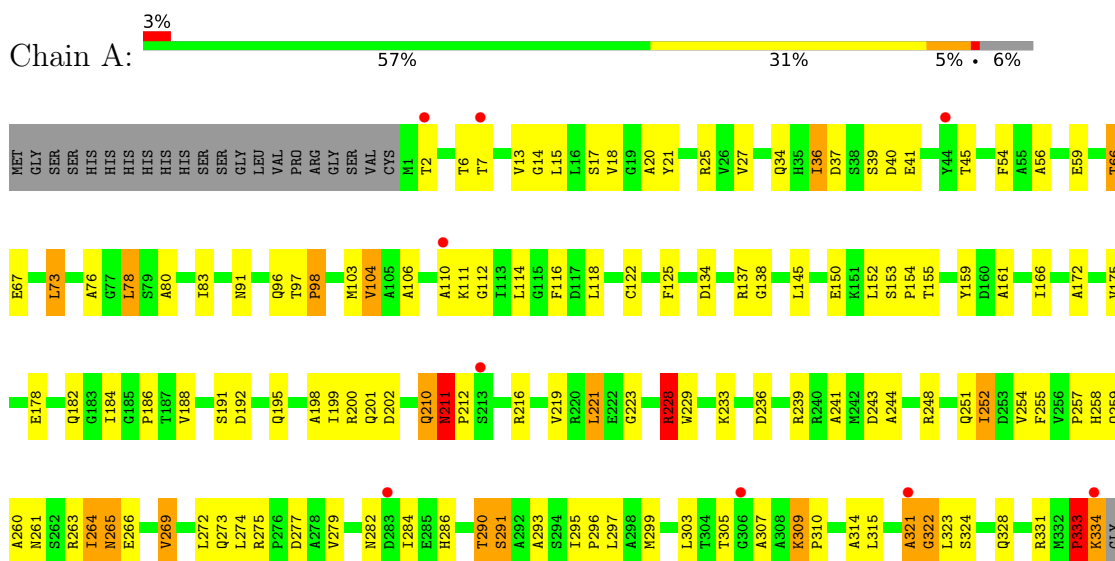
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	62	Total	O	0	0
			62	62		
2	B	39	Total	O	0	0
			39	39		

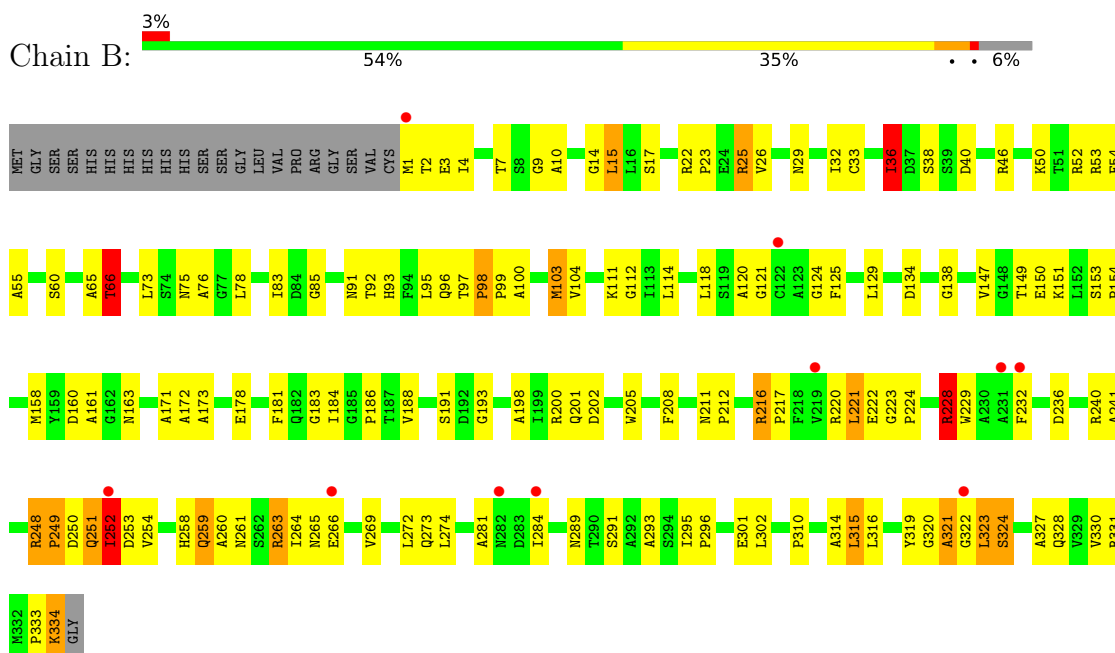
3 Residue-property plots

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: 3-oxoacyl-[acyl-carrier-protein] synthase III



• Molecule 1: 3-oxoacyl-[acyl-carrier-protein] synthase III



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.17Å 56.02Å 88.19Å 90.00° 90.57° 90.00°	Depositor
Resolution (Å)	28.01 – 2.50 28.01 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.0 (28.01-2.50) 92.2 (28.01-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.15 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.225 , 0.279 0.232 , 0.273	Depositor DCC
R_{free} test set	1980 reflections (9.39%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.473	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.038 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4955	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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	2.02	3/2469 (0.1%)	1.41	43/3358 (1.3%)
1	B	0.87	4/2469 (0.2%)	1.42	36/3358 (1.1%)
All	All	1.56	7/4938 (0.1%)	1.42	79/6716 (1.2%)

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	4
1	B	0	8
All	All	0	12

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	334	LYS	C-O	92.51	3.08	1.23
1	B	253	ASP	C-N	10.51	1.47	1.33
1	B	324	SER	C-O	9.84	1.36	1.23
1	A	309	LYS	C-N	8.03	1.43	1.33
1	B	10	ALA	N-CA	5.18	1.52	1.46
1	B	250	ASP	C-N	-5.11	1.26	1.33
1	A	310	PRO	N-CD	5.01	1.54	1.47

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	334	LYS	CA-C-O	11.40	140.18	120.80
1	A	309	LYS	CA-C-N	-10.84	108.34	119.93
1	A	309	LYS	C-N-CA	-10.84	108.34	119.93
1	B	248	ARG	CA-C-N	-9.86	107.52	119.84
1	B	248	ARG	C-N-CA	-9.86	107.52	119.84

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	211	ASN	CA-C-N	-8.65	109.03	119.84
1	A	211	ASN	C-N-CA	-8.65	109.03	119.84
1	A	211	ASN	CA-CB-CG	8.54	121.14	112.60
1	B	2	THR	CA-C-N	8.25	134.38	123.00
1	B	2	THR	C-N-CA	8.25	134.38	123.00
1	B	273	GLN	O-C-N	-7.95	112.01	122.59
1	B	248	ARG	CA-C-O	7.87	127.69	120.60
1	B	228	ARG	NE-CZ-NH2	7.71	126.14	119.20
1	B	40	ASP	CA-CB-CG	7.70	120.30	112.60
1	A	202	ASP	CA-CB-CG	7.69	120.29	112.60
1	A	37	ASP	O-C-N	-7.63	113.17	122.71
1	B	273	GLN	CA-C-O	7.62	131.41	120.51
1	B	228	ARG	NE-CZ-NH1	-7.62	113.89	121.50
1	A	104	VAL	N-CA-CB	7.58	118.92	110.51
1	A	264	ILE	N-CA-C	7.46	117.98	110.23
1	A	309	LYS	O-C-N	7.45	128.21	121.80
1	B	160	ASP	CA-CB-CG	7.43	120.03	112.60
1	A	212	PRO	CA-N-CD	-7.27	101.82	112.00
1	B	75	ASN	CA-CB-CG	7.16	119.76	112.60
1	A	265	ASN	CA-CB-CG	7.13	119.73	112.60
1	A	195	GLN	CA-C-O	6.97	129.23	120.65
1	B	66	THR	CA-CB-OG1	6.92	119.98	109.60
1	A	103	MET	CA-C-N	6.80	129.38	120.60
1	A	103	MET	C-N-CA	6.80	129.38	120.60
1	A	40	ASP	CA-CB-CG	6.79	119.39	112.60
1	A	66	THR	CA-CB-OG1	6.77	119.76	109.60
1	A	37	ASP	CA-CB-CG	6.67	119.27	112.60
1	B	249	PRO	CA-N-CD	-6.64	102.70	112.00
1	A	333	PRO	CA-C-N	6.56	133.51	121.70
1	A	333	PRO	C-N-CA	6.56	133.51	121.70
1	A	309	LYS	CA-C-O	6.51	127.82	120.46
1	A	282	ASN	CA-CB-CG	6.48	119.08	112.60
1	A	210	GLN	N-CA-C	6.46	118.12	111.14
1	B	220	ARG	O-C-N	6.44	130.89	123.29
1	B	252	ILE	N-CA-CB	6.44	120.95	110.13
1	B	208	PHE	CB-CA-C	6.42	121.05	110.90
1	B	66	THR	CA-CB-CG2	6.35	121.30	110.50
1	A	310	PRO	CA-N-CD	-6.33	103.13	112.00
1	A	195	GLN	O-C-N	-6.33	113.72	122.20
1	A	228	ARG	CD-NE-CZ	6.33	133.26	124.40
1	A	104	VAL	CA-CB-CG2	6.17	120.89	110.40
1	A	34	GLN	N-CA-C	-6.16	105.73	113.18

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	286	HIS	CA-C-O	6.09	126.66	119.48
1	A	104	VAL	CA-CB-CG1	6.07	120.72	110.40
1	B	252	ILE	CA-CB-CG1	6.07	120.72	110.40
1	B	134	ASP	CA-CB-CG	6.06	118.66	112.60
1	B	248	ARG	C-N-CD	6.00	149.58	125.00
1	A	66	THR	CA-CB-CG2	5.99	120.68	110.50
1	A	211	ASN	N-CA-C	5.98	123.02	109.81
1	A	265	ASN	OD1-CG-ND2	-5.97	116.63	122.60
1	B	253	ASP	CA-CB-CG	5.95	118.55	112.60
1	A	178	GLU	CB-CA-C	5.92	119.77	109.65
1	A	290	THR	CA-C-O	5.90	126.80	119.78
1	A	269	VAL	O-C-N	5.88	127.82	121.94
1	A	211	ASN	C-N-CD	5.83	148.91	125.00
1	B	1	MET	CA-C-N	5.78	132.94	122.02
1	B	1	MET	C-N-CA	5.78	132.94	122.02
1	A	134	ASP	CA-CB-CG	5.75	118.35	112.60
1	A	309	LYS	C-N-CD	5.72	148.46	125.00
1	B	253	ASP	O-C-N	5.67	130.63	122.13
1	B	252	ILE	CA-CB-CG2	5.61	120.04	110.50
1	B	228	ARG	CD-NE-CZ	5.61	132.25	124.40
1	B	200	ARG	N-CA-C	5.60	115.94	108.38
1	B	92	THR	CA-C-O	5.56	126.68	119.95
1	B	251	GLN	CA-C-N	5.55	127.94	120.50
1	B	251	GLN	C-N-CA	5.55	127.94	120.50
1	B	202	ASP	CA-CB-CG	5.48	118.08	112.60
1	B	3	GLU	O-C-N	-5.21	117.15	123.29
1	A	37	ASP	CA-C-O	5.16	129.29	122.63
1	A	104	VAL	O-C-N	5.13	127.07	121.94
1	A	265	ASN	CA-C-O	-5.07	115.15	120.63
1	B	98	PRO	CA-C-N	5.05	124.98	119.78
1	B	98	PRO	C-N-CA	5.05	124.98	119.78
1	B	75	ASN	OD1-CG-ND2	-5.03	117.57	122.60

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	GLN	Mainchain,Peptide
1	A	211	ASN	Mainchain
1	A	305	THR	Mainchain
1	B	171	ALA	Mainchain
1	B	228	ARG	Sidechain

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	B	232	PHE	Mainchain
1	B	323	LEU	Mainchain
1	B	36	ILE	Mainchain
1	B	46	ARG	Mainchain
1	B	78	LEU	Mainchain
1	B	9	GLY	Mainchain

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	2427	0	2402	109	0
1	B	2427	0	2402	125	0
2	A	62	0	0	3	0
2	B	39	0	0	2	0
All	All	4955	0	4804	206	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 21.

All (206) 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:B:261:ASN:HD22	1:B:264:ILE:HG13	1.26	0.99
1:A:228:ARG:HG3	1:A:228:ARG:HH11	1.29	0.95
1:B:258:HIS:CD2	1:B:260:ALA:HB2	2.02	0.94
1:A:259:GLN:HE21	1:A:284:ILE:H	0.94	0.90
1:A:152:LEU:O	1:A:155:THR:HG22	1.74	0.87
1:B:66:THR:HB	1:B:104:VAL:HG13	1.57	0.86
1:A:96:GLN:HE21	1:B:205:TRP:HE1	1.20	0.86
1:B:125:PHE:CE1	1:B:296:PRO:HG3	2.11	0.86
1:A:36:ILE:HD11	1:A:159:TYR:O	1.77	0.85
1:A:122:CYS:HB3	1:A:258:HIS:CE1	2.12	0.85
1:B:252:ILE:HD12	1:B:315:LEU:HB2	1.58	0.84
1:B:153:SER:HB2	1:B:154:PRO:HD3	1.60	0.83
1:A:322:GLY:HA3	1:B:99:PRO:HD2	1.62	0.81

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:LEU:HD11	1:A:83:ILE:HD11	1.61	0.81
1:A:259:GLN:NE2	1:A:284:ILE:H	1.79	0.80
1:B:228:ARG:HH21	1:B:228:ARG:HB3	1.48	0.78
1:A:172:ALA:HB3	1:A:296:PRO:HB2	1.65	0.76
1:A:259:GLN:HB3	1:A:284:ILE:HB	1.69	0.75
1:A:236:ASP:OD2	2:A:395:HOH:O	2.04	0.75
1:B:32:ILE:HD11	1:B:52:ARG:HD2	1.69	0.74
1:A:96:GLN:HE21	1:B:205:TRP:NE1	1.86	0.74
1:B:228:ARG:HB3	1:B:228:ARG:NH2	2.03	0.74
1:B:259:GLN:HB3	1:B:284:ILE:HB	1.71	0.73
1:A:259:GLN:HE21	1:A:284:ILE:N	1.80	0.73
1:B:281:ALA:HA	1:B:302:LEU:HD11	1.72	0.72
1:A:199:ILE:HG12	1:A:221:LEU:HD23	1.70	0.72
1:B:36:ILE:HG13	1:B:161:ALA:HB2	1.69	0.71
1:B:261:ASN:HB3	1:B:264:ILE:HD12	1.70	0.71
1:A:137:ARG:HH21	1:A:182:GLN:NE2	1.90	0.70
1:A:201:GLN:NE2	1:B:96:GLN:H	1.90	0.70
1:A:137:ARG:HH21	1:A:182:GLN:HE22	1.38	0.70
1:A:198:ALA:HB3	1:A:323:LEU:HD11	1.74	0.70
1:A:96:GLN:H	1:B:201:GLN:NE2	1.91	0.69
1:B:321:ALA:C	1:B:323:LEU:H	2.02	0.67
1:A:321:ALA:C	1:A:323:LEU:H	2.03	0.67
1:A:188:VAL:HG21	1:A:241:ALA:HA	1.76	0.66
1:A:76:ALA:HB1	1:A:78:LEU:HD22	1.76	0.66
1:B:258:HIS:HD2	1:B:260:ALA:HB2	1.59	0.65
1:A:15:LEU:CB	1:A:333:PRO:HB3	2.27	0.65
1:A:7:THR:HG23	1:A:138:GLY:O	1.97	0.65
1:B:125:PHE:CD1	1:B:296:PRO:HG3	2.31	0.65
1:B:36:ILE:CG1	1:B:161:ALA:HB2	2.26	0.64
1:A:7:THR:HG22	1:B:186:PRO:HB3	1.80	0.64
1:B:150:GLU:HG3	1:B:291:SER:HB3	1.82	0.62
1:B:29:ASN:OD1	1:B:50:LYS:O	2.17	0.62
1:A:161:ALA:HB3	2:A:379:HOH:O	1.97	0.62
1:A:76:ALA:CB	1:A:78:LEU:HD22	2.29	0.62
1:B:118:LEU:C	1:B:118:LEU:HD23	2.25	0.62
1:A:258:HIS:CD2	1:A:260:ALA:HB2	2.35	0.61
1:B:321:ALA:C	1:B:323:LEU:N	2.57	0.61
1:B:7:THR:HG23	1:B:138:GLY:O	2.01	0.61
1:B:252:ILE:CD1	1:B:315:LEU:HB2	2.30	0.61
1:B:172:ALA:HB3	1:B:296:PRO:HB2	1.83	0.61
1:B:221:LEU:HD22	1:B:222:GLU:N	2.15	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:GLN:HG2	1:B:201:GLN:HE22	1.66	0.61
1:A:17:SER:OG	1:A:76:ALA:HB2	2.00	0.60
1:A:125:PHE:CD1	1:A:296:PRO:HG3	2.36	0.60
1:A:192:ASP:OD2	1:B:111:LYS:HD2	2.01	0.60
1:A:36:ILE:CD1	1:A:159:TYR:O	2.49	0.60
1:A:322:GLY:CA	1:B:98:PRO:HB2	2.31	0.59
1:A:122:CYS:SG	1:A:291:SER:OG	2.43	0.59
1:B:114:LEU:HD12	1:B:114:LEU:C	2.29	0.58
1:B:252:ILE:HD13	1:B:274:LEU:HD22	1.86	0.57
1:A:56:ALA:HB3	1:A:59:GLU:HG3	1.87	0.57
1:B:188:VAL:O	1:B:188:VAL:HG13	2.04	0.57
1:B:125:PHE:CD1	1:B:296:PRO:CG	2.88	0.56
1:B:188:VAL:HG13	1:B:327:ALA:HB3	1.88	0.56
1:A:191:SER:O	1:B:112:GLY:HA2	2.05	0.56
1:B:229:TRP:CZ3	1:B:319:TYR:HB2	2.40	0.56
1:A:265:ASN:O	1:A:269:VAL:HG23	2.06	0.56
1:B:216:ARG:HD3	1:B:217:PRO:O	2.06	0.56
1:B:17:SER:HA	1:B:333:PRO:HB2	1.88	0.56
1:A:15:LEU:HB2	1:A:333:PRO:HB3	1.86	0.56
1:A:258:HIS:HD2	1:A:260:ALA:HB2	1.72	0.55
1:B:258:HIS:HE1	1:B:289:ASN:OD1	1.88	0.55
1:A:114:LEU:C	1:A:114:LEU:HD12	2.31	0.55
1:A:125:PHE:CE1	1:A:296:PRO:HG3	2.41	0.54
1:A:73:LEU:CD1	1:A:83:ILE:HD11	2.35	0.54
1:A:96:GLN:NE2	1:B:205:TRP:HE1	1.98	0.54
1:A:221:LEU:HD13	1:A:223:GLY:N	2.23	0.54
1:B:236:ASP:O	1:B:240:ARG:HG3	2.07	0.54
1:A:153:SER:OG	1:A:154:PRO:HD3	2.06	0.54
1:A:293:ALA:O	1:A:296:PRO:HD2	2.08	0.54
1:A:228:ARG:HG3	1:A:228:ARG:NH1	2.02	0.54
1:B:153:SER:CB	1:B:154:PRO:HD3	2.35	0.53
1:B:14:GLY:HA2	1:B:184:ILE:HG13	1.89	0.53
1:A:21:TYR:CD1	1:A:67:GLU:HG2	2.43	0.53
1:B:188:VAL:HG11	1:B:241:ALA:HB2	1.90	0.53
1:A:322:GLY:HA2	1:B:98:PRO:HB2	1.90	0.53
1:A:321:ALA:C	1:A:323:LEU:N	2.66	0.53
1:B:32:ILE:CD1	1:B:52:ARG:HD2	2.38	0.52
1:B:261:ASN:HB3	1:B:264:ILE:CD1	2.39	0.52
1:B:120:ALA:HB3	1:B:124:GLY:HA2	1.90	0.52
1:A:15:LEU:HB3	1:A:333:PRO:HB3	1.92	0.52
1:A:198:ALA:HB3	1:A:323:LEU:CD1	2.39	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:PRO:O	1:B:53:ARG:HD2	2.10	0.52
1:B:36:ILE:HG12	1:B:161:ALA:HA	1.92	0.52
1:A:275:ARG:HB3	1:A:277:ASP:OD1	2.10	0.52
1:B:125:PHE:CE1	1:B:296:PRO:CG	2.90	0.52
1:B:198:ALA:HB3	1:B:323:LEU:HD13	1.92	0.52
1:A:201:GLN:HE22	1:B:96:GLN:H	1.57	0.51
1:B:66:THR:CA	1:B:104:VAL:HG13	2.41	0.51
1:A:106:ALA:HB2	1:B:193:GLY:HA3	1.93	0.51
1:B:52:ARG:HH11	1:B:52:ARG:HG2	1.74	0.51
1:A:15:LEU:CD1	1:A:299:MET:HE1	2.41	0.51
1:B:259:GLN:O	1:B:284:ILE:HD12	2.11	0.51
1:A:248:ARG:O	1:A:251:GLN:HB2	2.11	0.50
1:A:45:THR:O	1:A:263:ARG:HD3	2.11	0.50
1:B:17:SER:OG	1:B:76:ALA:HB2	2.11	0.50
1:A:98:PRO:HB2	1:B:322:GLY:CA	2.42	0.50
1:A:221:LEU:HD13	1:A:223:GLY:H	1.77	0.50
1:A:331:ARG:HG3	1:B:4:ILE:HD12	1.93	0.50
1:B:97:THR:OG1	1:B:98:PRO:HA	2.11	0.49
1:A:321:ALA:HB1	1:B:97:THR:OG1	2.13	0.49
1:A:254:VAL:O	1:A:314:ALA:HA	2.13	0.49
1:B:248:ARG:O	1:B:249:PRO:C	2.53	0.49
1:B:258:HIS:CE1	1:B:289:ASN:OD1	2.66	0.49
1:A:96:GLN:HB3	1:B:91:ASN:O	2.13	0.49
1:A:261:ASN:HB3	1:A:264:ILE:HD12	1.94	0.49
1:B:66:THR:HA	1:B:104:VAL:HG13	1.95	0.49
1:B:15:LEU:HB3	1:B:333:PRO:HB3	1.95	0.48
1:B:29:ASN:O	1:B:33:CYS:SG	2.70	0.48
1:B:222:GLU:O	1:B:223:GLY:C	2.54	0.48
1:A:201:GLN:HE22	1:B:96:GLN:HG2	1.78	0.48
1:B:36:ILE:HG23	1:B:38:SER:HB2	1.96	0.48
1:A:293:ALA:C	1:A:296:PRO:HD2	2.38	0.48
1:A:2:THR:HG21	1:B:251:GLN:HE22	1.79	0.48
1:A:229:TRP:O	1:A:233:LYS:HB2	2.14	0.47
1:A:80:ALA:HA	1:A:83:ILE:HD12	1.96	0.47
1:B:23:PRO:HB3	1:B:55:ALA:HA	1.97	0.47
1:A:73:LEU:HD11	1:A:83:ILE:CD1	2.38	0.47
1:A:13:VAL:HG23	1:A:184:ILE:HD12	1.97	0.47
1:B:121:GLY:HA3	1:B:321:ALA:HB3	1.97	0.47
1:A:154:PRO:O	1:A:216:ARG:NH2	2.48	0.47
1:B:53:ARG:O	1:B:54:PHE:HD1	1.97	0.47
1:A:219:VAL:O	1:A:219:VAL:HG13	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:VAL:O	1:B:314:ALA:HA	2.15	0.46
1:B:320:GLY:O	1:B:323:LEU:HA	2.16	0.46
1:B:221:LEU:HD13	1:B:223:GLY:H	1.80	0.46
1:A:191:SER:O	1:B:111:LYS:O	2.34	0.46
1:B:60:SER:HA	1:B:151:LYS:HD2	1.95	0.46
1:B:151:LYS:O	1:B:154:PRO:HD2	2.16	0.46
1:B:198:ALA:HB3	1:B:323:LEU:CD1	2.46	0.46
1:A:96:GLN:H	1:B:201:GLN:HE22	1.63	0.46
1:B:26:VAL:HG22	1:B:53:ARG:NE	2.32	0.46
1:B:29:ASN:OD1	1:B:50:LYS:HA	2.15	0.46
1:B:103:MET:HE2	1:B:103:MET:HB2	1.73	0.46
1:A:252:ILE:HG12	1:A:274:LEU:HD22	1.97	0.45
1:B:295:ILE:HB	1:B:296:PRO:HD3	1.97	0.45
1:B:316:LEU:HD11	1:B:330:VAL:HG21	1.97	0.45
1:B:221:LEU:HD13	1:B:221:LEU:C	2.40	0.45
1:A:150:GLU:HG3	1:A:290:THR:O	2.16	0.45
1:A:7:THR:CG2	1:B:186:PRO:HB3	2.45	0.45
1:B:125:PHE:CZ	1:B:296:PRO:HG3	2.51	0.45
1:A:186:PRO:HG3	1:A:244:ALA:HB1	1.98	0.45
1:B:121:GLY:CA	1:B:321:ALA:HB3	2.47	0.44
1:A:15:LEU:HD12	1:A:333:PRO:HD3	2.00	0.44
1:A:118:LEU:C	1:A:118:LEU:HD23	2.42	0.44
1:A:272:LEU:O	1:A:273:GLN:HB2	2.17	0.44
1:A:27:VAL:HG21	1:A:54:PHE:CD2	2.52	0.44
1:A:73:LEU:HD11	1:A:145:LEU:HD11	1.98	0.44
1:A:116:PHE:HA	2:A:341:HOH:O	2.17	0.43
1:A:331:ARG:HG3	1:B:4:ILE:CD1	2.48	0.43
1:B:65:ALA:HB2	1:B:149:THR:HG21	1.99	0.43
1:A:188:VAL:CG2	1:A:241:ALA:HA	2.46	0.43
1:A:188:VAL:HG21	1:A:241:ALA:CA	2.46	0.43
1:A:295:ILE:HB	1:A:296:PRO:HD3	2.00	0.43
1:B:228:ARG:HB3	1:B:228:ARG:CZ	2.46	0.43
1:B:183:GLY:HA2	1:B:331:ARG:HG3	1.99	0.43
1:B:52:ARG:HG2	1:B:52:ARG:NH1	2.33	0.43
1:B:293:ALA:O	1:B:296:PRO:HD2	2.19	0.43
1:A:112:GLY:HA2	1:B:191:SER:O	2.18	0.42
1:B:100:ALA:O	1:B:103:MET:HB3	2.18	0.42
1:A:279:VAL:HG11	1:A:307:ALA:HB1	2.01	0.42
1:A:137:ARG:NH2	1:A:182:GLN:NE2	2.62	0.42
1:A:255:PHE:CZ	1:A:257:PRO:HB3	2.54	0.42
1:B:188:VAL:HG12	1:B:327:ALA:O	2.20	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:LEU:HD23	1:A:175:VAL:HG11	2.01	0.42
1:B:15:LEU:HD11	1:B:129:LEU:HD11	2.02	0.42
1:B:66:THR:HA	1:B:104:VAL:CG1	2.50	0.42
1:B:265:ASN:O	1:B:269:VAL:HG23	2.20	0.42
1:A:18:VAL:HG11	1:A:303:LEU:HD12	2.00	0.42
1:B:83:ILE:HG22	1:B:85:GLY:H	1.85	0.42
1:B:188:VAL:CG1	1:B:327:ALA:HB3	2.49	0.42
1:B:147:VAL:HG23	1:B:173:ALA:HB2	2.01	0.42
1:B:293:ALA:C	1:B:296:PRO:HD2	2.45	0.42
1:B:310:PRO:HB3	1:B:334:LYS:HG3	2.02	0.41
1:A:91:ASN:O	1:B:96:GLN:HB3	2.20	0.41
1:A:186:PRO:CG	1:A:244:ALA:HB1	2.50	0.41
1:A:252:ILE:HD11	1:A:274:LEU:HD22	2.01	0.41
1:A:20:ALA:HB1	1:A:297:LEU:HD23	2.03	0.41
1:A:254:VAL:HB	1:A:314:ALA:HB2	2.01	0.41
1:A:331:ARG:HG2	1:B:4:ILE:HD11	2.02	0.41
1:B:252:ILE:HD12	1:B:252:ILE:HG21	1.86	0.41
1:A:98:PRO:HB2	1:B:322:GLY:HA3	2.02	0.41
1:A:321:ALA:O	1:A:323:LEU:N	2.54	0.41
1:B:25:ARG:O	1:B:53:ARG:HA	2.20	0.41
1:B:93:HIS:CE1	1:B:95:LEU:HB2	2.56	0.41
1:B:188:VAL:CG1	1:B:241:ALA:HB2	2.51	0.40
1:B:224:PRO:HD3	2:B:389:HOH:O	2.21	0.40
1:B:261:ASN:HD21	1:B:263:ARG:HB2	1.86	0.40
1:A:14:GLY:HA2	1:A:184:ILE:HG13	2.04	0.40
1:A:110:ALA:O	1:A:112:GLY:N	2.54	0.40
1:A:239:ARG:O	1:A:243:ASP:HB2	2.22	0.40
1:B:111:LYS:HG2	2:B:365:HOH:O	2.21	0.40
1:B:211:ASN:N	1:B:212:PRO:HD3	2.36	0.40

There are no symmetry-related clashes.

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	332/356 (93%)	304 (92%)	24 (7%)	4 (1%)	10	20
1	B	332/356 (93%)	310 (93%)	19 (6%)	3 (1%)	14	27
All	All	664/712 (93%)	614 (92%)	43 (6%)	7 (1%)	11	22

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	321	ALA
1	A	111	LYS
1	A	321	ALA
1	A	333	PRO
1	B	163	ASN
1	B	259	GLN
1	A	322	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	244/262 (93%)	220 (90%)	24 (10%)	7	16
1	B	244/262 (93%)	222 (91%)	22 (9%)	9	19
All	All	488/524 (93%)	442 (91%)	46 (9%)	8	18

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	25	ARG
1	A	36	ILE
1	A	39	SER
1	A	41	GLU
1	A	66	THR
1	A	73	LEU
1	A	78	LEU
1	A	97	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	98	PRO
1	A	104	VAL
1	A	166	ILE
1	A	200	ARG
1	A	211	ASN
1	A	221	LEU
1	A	228	ARG
1	A	252	ILE
1	A	266	GLU
1	A	291	SER
1	A	309	LYS
1	A	315	LEU
1	A	324	SER
1	A	328	GLN
1	A	334	LYS
1	B	15	LEU
1	B	22	ARG
1	B	25	ARG
1	B	36	ILE
1	B	66	THR
1	B	73	LEU
1	B	103	MET
1	B	158	MET
1	B	178	GLU
1	B	181	PHE
1	B	216	ARG
1	B	221	LEU
1	B	228	ARG
1	B	252	ILE
1	B	263	ARG
1	B	266	GLU
1	B	272	LEU
1	B	301	GLU
1	B	315	LEU
1	B	324	SER
1	B	328	GLN
1	B	334	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	75	ASN
1	A	93	HIS
1	A	96	GLN
1	A	182	GLN
1	A	201	GLN
1	A	258	HIS
1	A	259	GLN
1	A	265	ASN
1	B	34	GLN
1	B	93	HIS
1	B	201	GLN
1	B	211	ASN
1	B	251	GLN
1	B	258	HIS
1	B	261	ASN
1	B	265	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 ⓘ

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	334/356 (93%)	0.22	9 (2%) 56 51	18, 38, 58, 81	0
1	B	334/356 (93%)	0.39	10 (2%) 52 48	18, 43, 62, 79	0
All	All	668/712 (93%)	0.31	19 (2%) 55 50	18, 41, 60, 81	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	232	PHE	3.5
1	A	334	LYS	3.1
1	B	252	ILE	3.0
1	A	321	ALA	3.0
1	B	322	GLY	2.9
1	B	1	MET	2.9
1	A	283	ASP	2.7
1	B	122	CYS	2.6
1	A	7	THR	2.6
1	A	306	GLY	2.6
1	B	282	ASN	2.4
1	A	110	ALA	2.4
1	A	213	SER	2.3
1	A	44	TYR	2.3
1	B	284	ILE	2.2
1	B	219	VAL	2.1
1	A	2	THR	2.0
1	B	231	ALA	2.0
1	B	266	GLU	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.