



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:40 AM UTC

PDB ID : 4LK5 / pdb_00004lk5
Title : Crystal structure of a enoyl-CoA hydratase from Mycobacterium avium subsp. paratuberculosis K-10
Authors : Kumaran, D.; Chamala, S.; Evans, B.; Foti, R.; Gizzi, A.; Hillerich, B.; Kar, A.; Lafleur, J.; Seidel, R.; Villigas, G.; Zencheck, W.; Al Obaidi, N.; Almo, S.C.; Swaminathan, S.; New York Structural Genomics Research Consortium (NYSGRC)
Deposited on : 2013-07-06
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
Mogul	:	2022.3.0, CSD as543be (2022)
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)

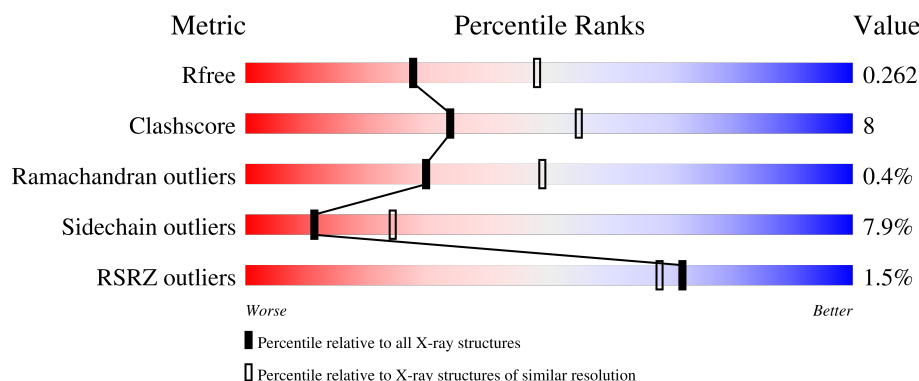
1 Overall quality at a glance

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	292	% 70% 10% • 18%
1	B	292	% 67% 14% •• 17%
1	C	292	% 67% 13% •• 16%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5342 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 enoyl-CoA hydratase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	Se	0	0	0
			1725	1104	288	327	1	5			
1	B	242	Total	C	N	O	S	Se	0	0	0
			1755	1124	293	332	1	5			
1	C	246	Total	C	N	O	S	Se	0	0	0
			1791	1143	303	339	1	5			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MSE	-	expression tag	UNP Q73WM1
A	-20	HIS	-	expression tag	UNP Q73WM1
A	-19	HIS	-	expression tag	UNP Q73WM1
A	-18	HIS	-	expression tag	UNP Q73WM1
A	-17	HIS	-	expression tag	UNP Q73WM1
A	-16	HIS	-	expression tag	UNP Q73WM1
A	-15	HIS	-	expression tag	UNP Q73WM1
A	-14	SER	-	expression tag	UNP Q73WM1
A	-13	SER	-	expression tag	UNP Q73WM1
A	-12	GLY	-	expression tag	UNP Q73WM1
A	-11	VAL	-	expression tag	UNP Q73WM1
A	-10	ASP	-	expression tag	UNP Q73WM1
A	-9	LEU	-	expression tag	UNP Q73WM1
A	-8	GLY	-	expression tag	UNP Q73WM1
A	-7	THR	-	expression tag	UNP Q73WM1
A	-6	GLU	-	expression tag	UNP Q73WM1
A	-5	ASN	-	expression tag	UNP Q73WM1
A	-4	LEU	-	expression tag	UNP Q73WM1
A	-3	TYR	-	expression tag	UNP Q73WM1
A	-2	PHE	-	expression tag	UNP Q73WM1
A	-1	GLN	-	expression tag	UNP Q73WM1
A	0	SER	-	expression tag	UNP Q73WM1
B	-21	MSE	-	expression tag	UNP Q73WM1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	HIS	-	expression tag	UNP Q73WM1
B	-19	HIS	-	expression tag	UNP Q73WM1
B	-18	HIS	-	expression tag	UNP Q73WM1
B	-17	HIS	-	expression tag	UNP Q73WM1
B	-16	HIS	-	expression tag	UNP Q73WM1
B	-15	HIS	-	expression tag	UNP Q73WM1
B	-14	SER	-	expression tag	UNP Q73WM1
B	-13	SER	-	expression tag	UNP Q73WM1
B	-12	GLY	-	expression tag	UNP Q73WM1
B	-11	VAL	-	expression tag	UNP Q73WM1
B	-10	ASP	-	expression tag	UNP Q73WM1
B	-9	LEU	-	expression tag	UNP Q73WM1
B	-8	GLY	-	expression tag	UNP Q73WM1
B	-7	THR	-	expression tag	UNP Q73WM1
B	-6	GLU	-	expression tag	UNP Q73WM1
B	-5	ASN	-	expression tag	UNP Q73WM1
B	-4	LEU	-	expression tag	UNP Q73WM1
B	-3	TYR	-	expression tag	UNP Q73WM1
B	-2	PHE	-	expression tag	UNP Q73WM1
B	-1	GLN	-	expression tag	UNP Q73WM1
B	0	SER	-	expression tag	UNP Q73WM1
C	-21	MSE	-	expression tag	UNP Q73WM1
C	-20	HIS	-	expression tag	UNP Q73WM1
C	-19	HIS	-	expression tag	UNP Q73WM1
C	-18	HIS	-	expression tag	UNP Q73WM1
C	-17	HIS	-	expression tag	UNP Q73WM1
C	-16	HIS	-	expression tag	UNP Q73WM1
C	-15	HIS	-	expression tag	UNP Q73WM1
C	-14	SER	-	expression tag	UNP Q73WM1
C	-13	SER	-	expression tag	UNP Q73WM1
C	-12	GLY	-	expression tag	UNP Q73WM1
C	-11	VAL	-	expression tag	UNP Q73WM1
C	-10	ASP	-	expression tag	UNP Q73WM1
C	-9	LEU	-	expression tag	UNP Q73WM1
C	-8	GLY	-	expression tag	UNP Q73WM1
C	-7	THR	-	expression tag	UNP Q73WM1
C	-6	GLU	-	expression tag	UNP Q73WM1
C	-5	ASN	-	expression tag	UNP Q73WM1
C	-4	LEU	-	expression tag	UNP Q73WM1
C	-3	TYR	-	expression tag	UNP Q73WM1
C	-2	PHE	-	expression tag	UNP Q73WM1
C	-1	GLN	-	expression tag	UNP Q73WM1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP Q73WM1

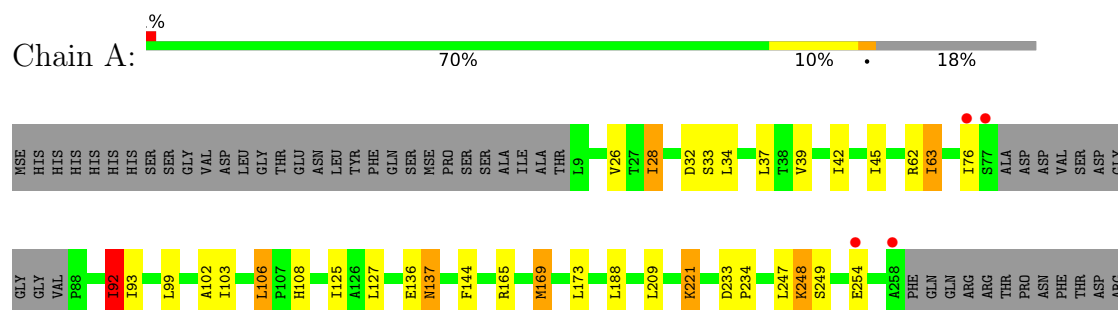
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	29	Total	O	0	0
			29	29		
2	B	26	Total	O	0	0
			26	26		
2	C	16	Total	O	0	0
			16	16		

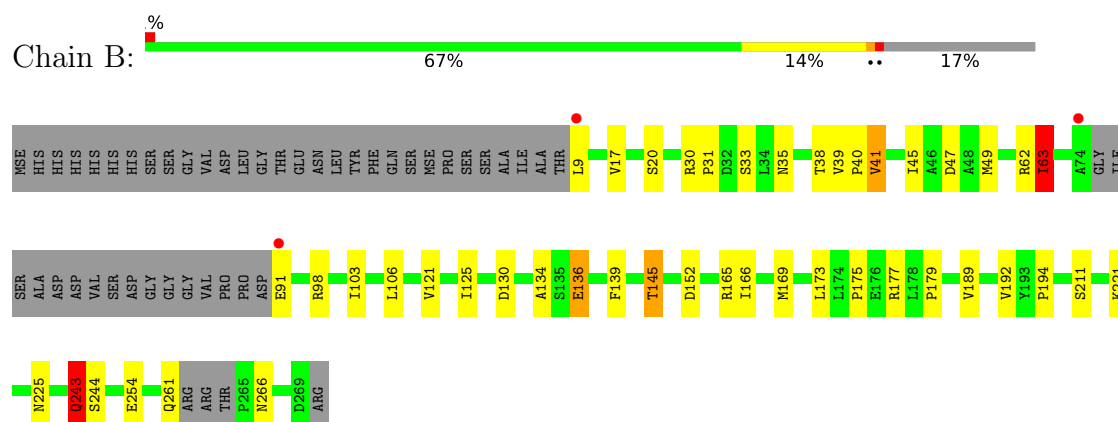
3 Residue-property plots [i](#)

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.

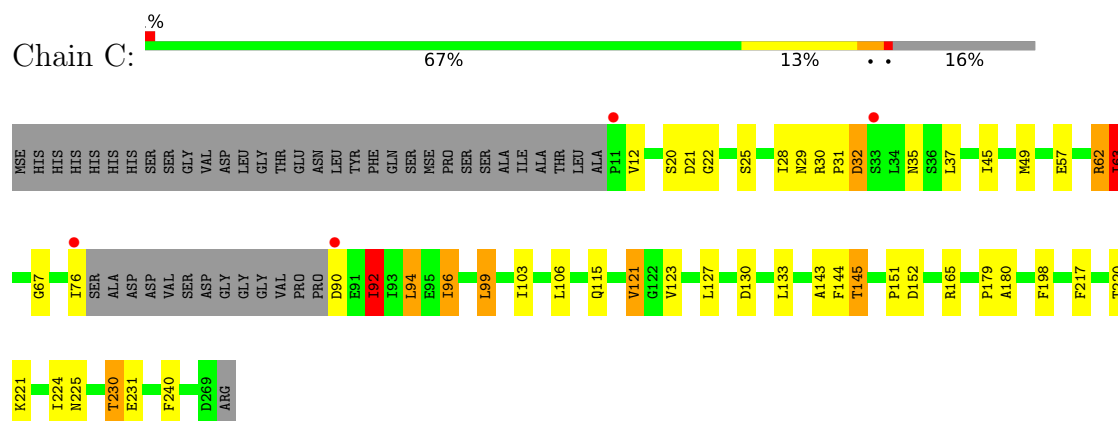
• Molecule 1: enoyl-CoA hydratase



• Molecule 1: enoyl-CoA hydratase



• Molecule 1: enoyl-CoA hydratase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	78.33Å 78.33Å 212.93Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.04 – 2.50 49.04 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.04-2.50) 100.0 (49.04-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.62 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.199 , 0.258 0.213 , 0.262	Depositor DCC
R_{free} test set	1360 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 22.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5342	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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	0.94	0/1748	1.15	13/2379 (0.5%)
1	B	0.91	0/1778	1.08	4/2417 (0.2%)
1	C	0.95	0/1815	1.11	6/2467 (0.2%)
All	All	0.94	0/5341	1.11	23/7263 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	63	ILE	CB-CA-C	-6.67	100.67	110.81
1	A	103	ILE	CB-CA-C	-6.46	103.70	111.97
1	A	249	SER	CA-C-N	6.27	125.49	118.97
1	A	249	SER	C-N-CA	6.27	125.49	118.97
1	A	28	ILE	CB-CA-C	6.06	118.42	110.91
1	A	137	ASN	CB-CA-C	-5.74	100.92	110.68
1	C	96	ILE	CB-CA-C	-5.73	104.55	111.88
1	B	166	ILE	N-CA-C	-5.67	105.32	110.82
1	A	108	HIS	CA-C-N	5.62	126.50	119.98
1	A	108	HIS	C-N-CA	5.62	126.50	119.98
1	B	125	ILE	CB-CA-C	-5.59	104.81	111.97
1	C	165	ARG	N-CA-C	5.59	117.05	111.07
1	A	92	ILE	N-CA-CB	5.57	117.07	110.55
1	A	144	PHE	N-CA-C	5.53	118.02	111.33
1	C	198	PHE	N-CA-C	5.51	117.72	111.11
1	B	243	GLN	CB-CA-C	-5.51	102.23	110.88
1	A	125	ILE	CB-CA-C	-5.48	104.96	111.97
1	C	92	ILE	CB-CA-C	-5.14	105.30	111.88
1	A	92	ILE	CB-CA-C	-5.13	105.40	111.97
1	B	63	ILE	CB-CA-C	-5.11	103.00	110.62
1	C	12	VAL	N-CA-C	5.10	115.62	108.17
1	A	136	GLU	CA-C-N	5.02	127.26	120.38
1	A	136	GLU	C-N-CA	5.02	127.26	120.38

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1725	0	1777	25	0
1	B	1755	0	1797	33	0
1	C	1791	0	1833	28	0
2	A	29	0	0	3	0
2	B	26	0	0	2	0
2	C	16	0	0	5	0
All	All	5342	0	5407	82	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 8.

All (82) 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:169:MSE:SE	2:C:310:HOH:O	2.26	1.04
1:B:49:MSE:HE1	1:B:63:ILE:HD12	1.42	1.01
1:A:169:MSE:SE	2:B:325:HOH:O	2.36	0.92
1:A:106:LEU:O	1:A:221:LYS:HE3	1.74	0.88
1:B:169:MSE:CE	1:B:173:LEU:HD11	2.05	0.87
1:A:76:ILE:HG22	2:A:318:HOH:O	1.80	0.82
1:B:169:MSE:HE2	1:C:130:ASP:HB3	1.62	0.82
1:B:169:MSE:HE1	1:B:173:LEU:HD11	1.65	0.78
1:A:169:MSE:HE2	1:A:173:LEU:CD1	2.15	0.77
1:B:169:MSE:CE	1:B:173:LEU:CD1	2.64	0.75
1:A:169:MSE:CE	1:A:173:LEU:HG	2.19	0.73
1:B:49:MSE:HE1	1:B:63:ILE:CD1	2.18	0.72
1:C:67:GLY:HA3	2:C:314:HOH:O	1.90	0.70
1:A:127:LEU:HD11	1:A:188:LEU:HD11	1.73	0.70
1:C:230:THR:HB	2:C:308:HOH:O	1.93	0.68
1:C:31:PRO:HB3	2:C:314:HOH:O	1.94	0.67
1:A:169:MSE:HE3	1:A:169:MSE:HA	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:GLU:HG3	1:B:194:PRO:HA	1.81	0.62
1:A:137:ASN:HB3	2:A:317:HOH:O	2.01	0.61
1:B:134:ALA:O	1:B:192:VAL:HA	2.02	0.60
1:B:62:ARG:C	1:B:63:ILE:HD13	2.26	0.60
1:C:220:THR:O	1:C:224:ILE:HG13	2.03	0.58
1:B:63:ILE:HD13	1:B:63:ILE:N	2.19	0.57
1:C:92:ILE:O	1:C:96:ILE:HG12	2.05	0.55
1:A:93:ILE:HD11	1:A:247:LEU:HD12	1.88	0.55
1:B:139:PHE:HB3	1:B:177:ARG:HG2	1.88	0.53
1:A:37:LEU:HB3	1:A:42:ILE:HD11	1.91	0.53
1:B:169:MSE:HE3	1:B:173:LEU:HG	1.90	0.52
1:A:169:MSE:CE	1:A:173:LEU:CG	2.88	0.50
1:C:49:MSE:CE	1:C:63:ILE:HD12	2.41	0.50
1:B:169:MSE:HE3	1:B:173:LEU:CD1	2.39	0.50
1:B:145:THR:HB	2:B:313:HOH:O	2.11	0.50
1:A:169:MSE:HE2	1:A:173:LEU:CG	2.42	0.50
1:B:130:ASP:OD2	1:B:221:LYS:HE2	2.11	0.50
1:B:130:ASP:CG	1:B:221:LYS:HE2	2.37	0.49
1:A:165:ARG:HH21	1:B:225:ASN:HD21	1.60	0.49
1:A:169:MSE:HE2	1:A:173:LEU:HG	1.92	0.49
1:C:20:SER:OG	1:C:21:ASP:N	2.43	0.48
1:A:169:MSE:HE3	1:A:169:MSE:CA	2.43	0.48
1:C:45:ILE:HG22	1:C:99:LEU:HD13	1.96	0.47
1:B:165:ARG:HH21	1:C:225:ASN:HD21	1.61	0.47
1:A:45:ILE:HG22	1:A:99:LEU:CD1	2.45	0.46
1:B:30:ARG:N	1:B:31:PRO:CD	2.78	0.46
1:C:94:LEU:HD13	1:C:240:PHE:CE1	2.50	0.46
1:C:144:PHE:CD2	1:C:151:PRO:HA	2.51	0.46
1:C:30:ARG:H	1:C:35:ASN:HD22	1.63	0.46
1:C:22:GLY:HA2	1:C:57:GLU:O	2.16	0.46
1:B:41:VAL:O	1:B:45:ILE:HG13	2.16	0.46
1:C:121:VAL:HG13	1:C:152:ASP:OD1	2.16	0.45
1:C:130:ASP:OD1	1:C:221:LYS:HE2	2.16	0.45
1:A:169:MSE:HE2	1:A:173:LEU:HD11	1.95	0.45
1:B:30:ARG:H	1:B:35:ASN:ND2	2.14	0.45
1:A:102:ALA:HB2	2:A:328:HOH:O	2.16	0.45
1:B:173:LEU:C	1:B:175:PRO:HD3	2.41	0.45
1:B:139:PHE:CD1	1:B:179:PRO:HA	2.52	0.45
1:C:143:ALA:O	1:C:144:PHE:C	2.58	0.45
1:B:39:VAL:HB	1:B:40:PRO:CD	2.47	0.45
1:C:49:MSE:HE1	1:C:63:ILE:HD12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:145:THR:HG21	1:C:217:PHE:HE1	1.82	0.44
1:C:123:VAL:HG12	1:C:127:LEU:HD12	1.99	0.44
1:B:47:ASP:OD1	1:B:98:ARG:NH2	2.44	0.44
1:C:145:THR:HB	2:C:306:HOH:O	2.18	0.44
1:B:169:MSE:HE3	1:B:173:LEU:CG	2.47	0.44
1:C:99:LEU:HD12	1:C:99:LEU:C	2.43	0.43
1:C:103:ILE:O	1:C:221:LYS:NZ	2.52	0.43
1:B:254:GLU:HG2	1:B:266:ASN:O	2.18	0.43
1:B:136:GLU:CG	1:B:194:PRO:HA	2.47	0.43
1:B:243:GLN:HA	1:B:243:GLN:OE1	2.19	0.42
1:A:169:MSE:CE	1:A:173:LEU:CD1	2.92	0.42
1:A:233:ASP:HB2	1:A:234:PRO:HD3	2.01	0.42
1:C:179:PRO:O	1:C:180:ALA:C	2.63	0.42
1:A:248:LYS:HE2	1:A:248:LYS:HB2	1.73	0.42
1:C:28:ILE:HD13	1:C:37:LEU:HD21	2.02	0.42
1:A:169:MSE:HE2	1:A:173:LEU:HD12	1.98	0.41
1:C:25:SER:OG	1:C:62:ARG:HD2	2.19	0.41
1:C:230:THR:HG22	1:C:231:GLU:HG2	2.02	0.41
1:A:26:VAL:O	1:A:63:ILE:HA	2.19	0.41
1:B:130:ASP:OD1	1:B:221:LYS:HE2	2.21	0.41
1:B:103:ILE:O	1:B:221:LYS:NZ	2.55	0.40
1:A:37:LEU:HB3	1:A:42:ILE:CD1	2.51	0.40
1:C:29:ASN:C	1:C:31:PRO:HD3	2.46	0.40
1:A:39:VAL:CG2	1:A:92:ILE:HD12	2.52	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	236/292 (81%)	227 (96%)	9 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	236/292 (81%)	224 (95%)	10 (4%)	2 (1%)	16	31
1	C	242/292 (83%)	230 (95%)	11 (4%)	1 (0%)	30	49
All	All	714/876 (82%)	681 (95%)	30 (4%)	3 (0%)	30	49

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	136	GLU
1	C	32	ASP
1	B	20	SER

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	177/214 (83%)	164 (93%)	13 (7%)	13	27
1	B	180/214 (84%)	164 (91%)	16 (9%)	9	20
1	C	184/214 (86%)	170 (92%)	14 (8%)	12	26
All	All	541/642 (84%)	498 (92%)	43 (8%)	11	24

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ILE
1	A	32	ASP
1	A	33	SER
1	A	34	LEU
1	A	62	ARG
1	A	63	ILE
1	A	92	ILE
1	A	106	LEU
1	A	169	MSE
1	A	209	LEU
1	A	221	LYS

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Mol	Chain	Res	Type
1	A	248	LYS
1	A	254	GLU
1	B	9	LEU
1	B	17	VAL
1	B	33	SER
1	B	38	THR
1	B	41	VAL
1	B	63	ILE
1	B	91	GLU
1	B	106	LEU
1	B	121	VAL
1	B	145	THR
1	B	152	ASP
1	B	189	VAL
1	B	211	SER
1	B	243	GLN
1	B	244	SER
1	B	261	GLN
1	C	32	ASP
1	C	62	ARG
1	C	63	ILE
1	C	76	ILE
1	C	90	ASP
1	C	92	ILE
1	C	94	LEU
1	C	99	LEU
1	C	106	LEU
1	C	115	GLN
1	C	121	VAL
1	C	133	LEU
1	C	145	THR
1	C	230	THR

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

Mol	Chain	Res	Type
1	A	108	HIS
1	A	137	ASN
1	B	35	ASN
1	B	225	ASN
1	C	35	ASN
1	C	225	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](#)

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	235/292 (80%)	0.03	4 (1%) 69 65	28, 46, 71, 101	0
1	B	237/292 (81%)	0.01	3 (1%) 75 71	27, 46, 75, 94	0
1	C	241/292 (82%)	-0.09	4 (1%) 69 65	27, 45, 66, 83	0
All	All	713/876 (81%)	-0.02	11 (1%) 72 68	27, 46, 71, 101	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	11	PRO	4.2
1	A	258	ALA	3.5
1	C	76	ILE	3.4
1	A	76	ILE	3.1
1	C	33	SER	2.8
1	B	74	ALA	2.6
1	A	77	SER	2.5
1	A	254	GLU	2.4
1	B	9	LEU	2.2
1	C	90	ASP	2.2
1	B	91	GLU	2.1

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.