



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:49 AM UTC

PDB ID : 6ACQ / pdb_00006acq
Title : Crystal structure of (S)-3-hydroxybutyryl-CoA dehydrogenase from *Clostridium acetobutylicum*, apo form
Authors : Takenoya, M.; Taguchi, S.; Yajima, S.
Deposited on : 2018-07-27
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:

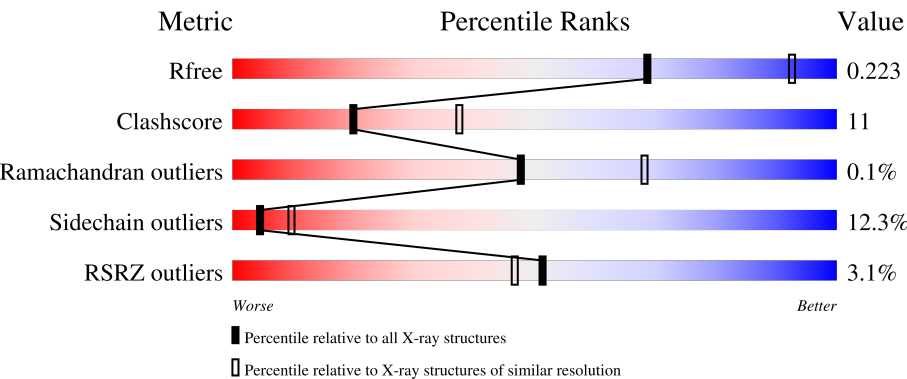
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

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	302	2% 73% 17% • 7%
1	B	302	2% 71% 20% 5% •
1	C	302	% 72% 20% • 7%
1	D	302	8% 60% 26% 6% • 7%
1	E	302	3% 65% 24% 5% • 7%

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Mol	Chain	Length	Quality of chain
1	F	302	% 76% 18%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12991 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-hydroxybutyryl-CoA dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2132	1359	355	404	14			
1	B	289	Total	C	N	O	S	0	0	0
			2189	1394	368	413	14			
1	C	281	Total	C	N	O	S	0	0	0
			2132	1359	355	404	14			
1	D	281	Total	C	N	O	S	0	0	0
			2132	1359	355	404	14			
1	E	281	Total	C	N	O	S	0	0	0
			2132	1359	355	404	14			
1	F	290	Total	C	N	O	S	0	0	0
			2195	1397	369	415	14			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P52041
A	-18	GLY	-	expression tag	UNP P52041
A	-17	SER	-	expression tag	UNP P52041
A	-16	SER	-	expression tag	UNP P52041
A	-15	HIS	-	expression tag	UNP P52041
A	-14	HIS	-	expression tag	UNP P52041
A	-13	HIS	-	expression tag	UNP P52041
A	-12	HIS	-	expression tag	UNP P52041
A	-11	HIS	-	expression tag	UNP P52041
A	-10	HIS	-	expression tag	UNP P52041
A	-9	SER	-	expression tag	UNP P52041
A	-8	SER	-	expression tag	UNP P52041
A	-7	GLY	-	expression tag	UNP P52041
A	-6	LEU	-	expression tag	UNP P52041
A	-5	VAL	-	expression tag	UNP P52041
A	-4	PRO	-	expression tag	UNP P52041
A	-3	ARG	-	expression tag	UNP P52041

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P52041
A	-1	SER	-	expression tag	UNP P52041
A	0	HIS	-	expression tag	UNP P52041
B	-19	MET	-	expression tag	UNP P52041
B	-18	GLY	-	expression tag	UNP P52041
B	-17	SER	-	expression tag	UNP P52041
B	-16	SER	-	expression tag	UNP P52041
B	-15	HIS	-	expression tag	UNP P52041
B	-14	HIS	-	expression tag	UNP P52041
B	-13	HIS	-	expression tag	UNP P52041
B	-12	HIS	-	expression tag	UNP P52041
B	-11	HIS	-	expression tag	UNP P52041
B	-10	HIS	-	expression tag	UNP P52041
B	-9	SER	-	expression tag	UNP P52041
B	-8	SER	-	expression tag	UNP P52041
B	-7	GLY	-	expression tag	UNP P52041
B	-6	LEU	-	expression tag	UNP P52041
B	-5	VAL	-	expression tag	UNP P52041
B	-4	PRO	-	expression tag	UNP P52041
B	-3	ARG	-	expression tag	UNP P52041
B	-2	GLY	-	expression tag	UNP P52041
B	-1	SER	-	expression tag	UNP P52041
B	0	HIS	-	expression tag	UNP P52041
C	-19	MET	-	expression tag	UNP P52041
C	-18	GLY	-	expression tag	UNP P52041
C	-17	SER	-	expression tag	UNP P52041
C	-16	SER	-	expression tag	UNP P52041
C	-15	HIS	-	expression tag	UNP P52041
C	-14	HIS	-	expression tag	UNP P52041
C	-13	HIS	-	expression tag	UNP P52041
C	-12	HIS	-	expression tag	UNP P52041
C	-11	HIS	-	expression tag	UNP P52041
C	-10	HIS	-	expression tag	UNP P52041
C	-9	SER	-	expression tag	UNP P52041
C	-8	SER	-	expression tag	UNP P52041
C	-7	GLY	-	expression tag	UNP P52041
C	-6	LEU	-	expression tag	UNP P52041
C	-5	VAL	-	expression tag	UNP P52041
C	-4	PRO	-	expression tag	UNP P52041
C	-3	ARG	-	expression tag	UNP P52041
C	-2	GLY	-	expression tag	UNP P52041
C	-1	SER	-	expression tag	UNP P52041

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP P52041
D	-19	MET	-	expression tag	UNP P52041
D	-18	GLY	-	expression tag	UNP P52041
D	-17	SER	-	expression tag	UNP P52041
D	-16	SER	-	expression tag	UNP P52041
D	-15	HIS	-	expression tag	UNP P52041
D	-14	HIS	-	expression tag	UNP P52041
D	-13	HIS	-	expression tag	UNP P52041
D	-12	HIS	-	expression tag	UNP P52041
D	-11	HIS	-	expression tag	UNP P52041
D	-10	HIS	-	expression tag	UNP P52041
D	-9	SER	-	expression tag	UNP P52041
D	-8	SER	-	expression tag	UNP P52041
D	-7	GLY	-	expression tag	UNP P52041
D	-6	LEU	-	expression tag	UNP P52041
D	-5	VAL	-	expression tag	UNP P52041
D	-4	PRO	-	expression tag	UNP P52041
D	-3	ARG	-	expression tag	UNP P52041
D	-2	GLY	-	expression tag	UNP P52041
D	-1	SER	-	expression tag	UNP P52041
D	0	HIS	-	expression tag	UNP P52041
E	-19	MET	-	expression tag	UNP P52041
E	-18	GLY	-	expression tag	UNP P52041
E	-17	SER	-	expression tag	UNP P52041
E	-16	SER	-	expression tag	UNP P52041
E	-15	HIS	-	expression tag	UNP P52041
E	-14	HIS	-	expression tag	UNP P52041
E	-13	HIS	-	expression tag	UNP P52041
E	-12	HIS	-	expression tag	UNP P52041
E	-11	HIS	-	expression tag	UNP P52041
E	-10	HIS	-	expression tag	UNP P52041
E	-9	SER	-	expression tag	UNP P52041
E	-8	SER	-	expression tag	UNP P52041
E	-7	GLY	-	expression tag	UNP P52041
E	-6	LEU	-	expression tag	UNP P52041
E	-5	VAL	-	expression tag	UNP P52041
E	-4	PRO	-	expression tag	UNP P52041
E	-3	ARG	-	expression tag	UNP P52041
E	-2	GLY	-	expression tag	UNP P52041
E	-1	SER	-	expression tag	UNP P52041
E	0	HIS	-	expression tag	UNP P52041
F	-19	MET	-	expression tag	UNP P52041

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-18	GLY	-	expression tag	UNP P52041
F	-17	SER	-	expression tag	UNP P52041
F	-16	SER	-	expression tag	UNP P52041
F	-15	HIS	-	expression tag	UNP P52041
F	-14	HIS	-	expression tag	UNP P52041
F	-13	HIS	-	expression tag	UNP P52041
F	-12	HIS	-	expression tag	UNP P52041
F	-11	HIS	-	expression tag	UNP P52041
F	-10	HIS	-	expression tag	UNP P52041
F	-9	SER	-	expression tag	UNP P52041
F	-8	SER	-	expression tag	UNP P52041
F	-7	GLY	-	expression tag	UNP P52041
F	-6	LEU	-	expression tag	UNP P52041
F	-5	VAL	-	expression tag	UNP P52041
F	-4	PRO	-	expression tag	UNP P52041
F	-3	ARG	-	expression tag	UNP P52041
F	-2	GLY	-	expression tag	UNP P52041
F	-1	SER	-	expression tag	UNP P52041
F	0	HIS	-	expression tag	UNP P52041

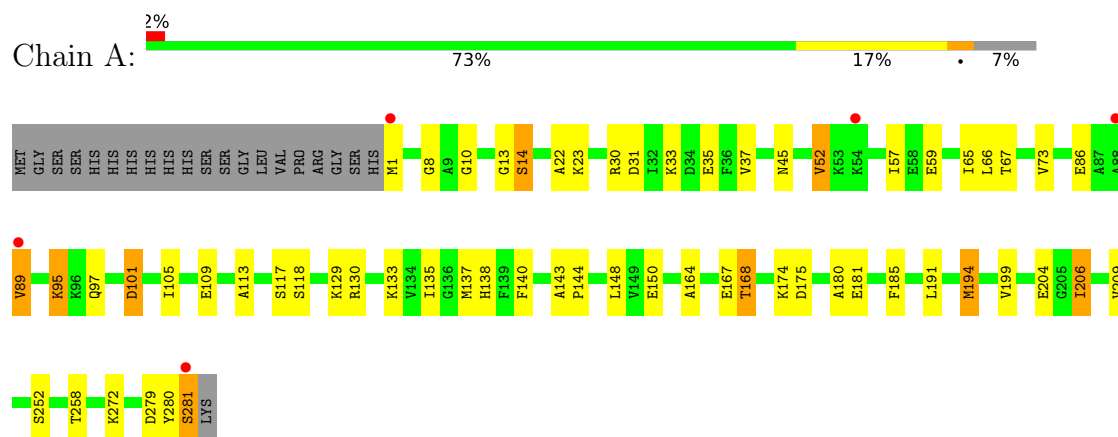
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	23	Total O 23 23	0	0
2	B	8	Total O 8 8	0	0
2	C	16	Total O 16 16	0	0
2	D	4	Total O 4 4	0	0
2	E	8	Total O 8 8	0	0
2	F	20	Total O 20 20	0	0

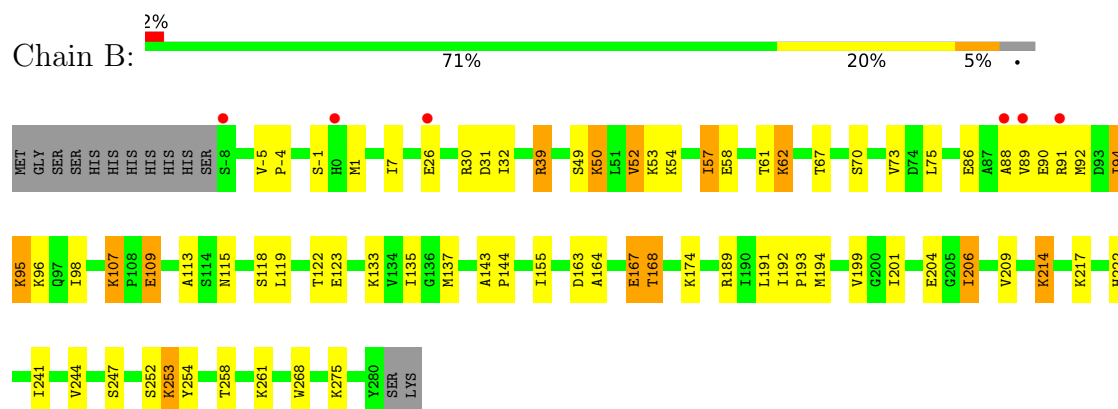
3 Residue-property plots [i](#)

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.

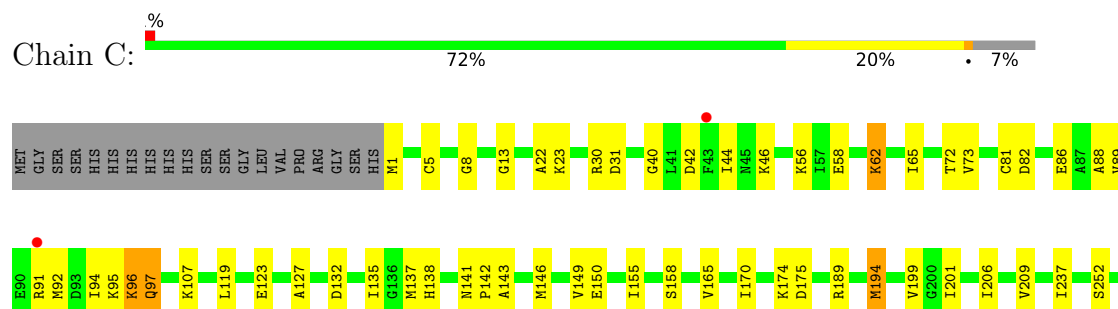
- Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



- Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase

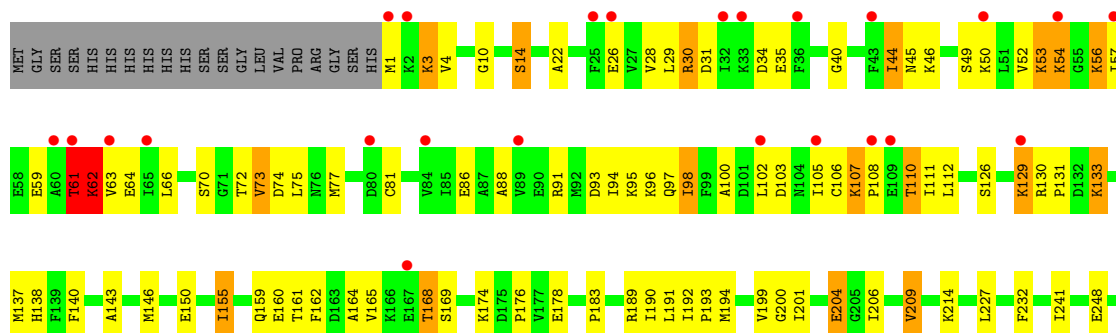


- Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase

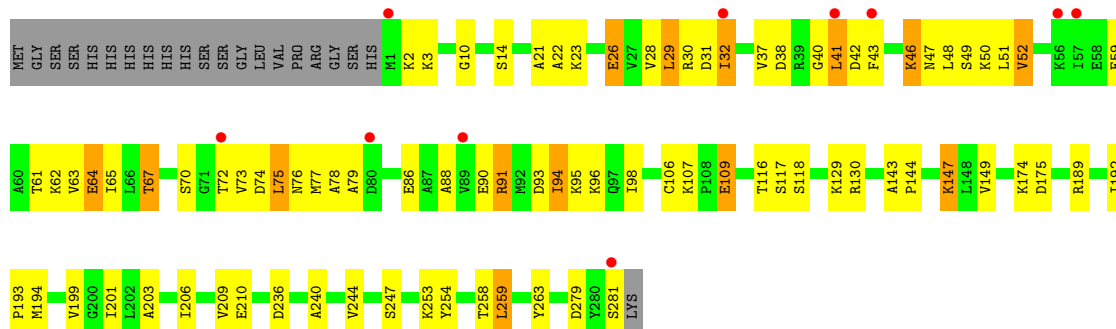




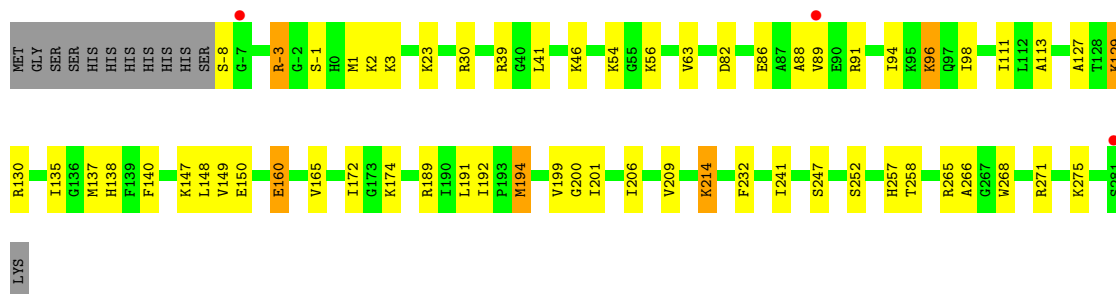
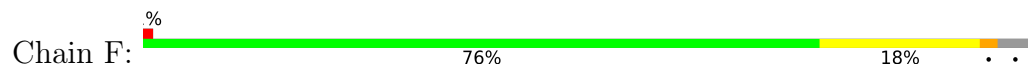
• Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



• Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



• Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	133.43Å 134.48Å 153.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.90 – 2.50 47.90 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.90-2.50) 99.6 (47.90-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0218	Depositor
R, R_{free}	0.199 , 0.222 0.203 , 0.223	Depositor DCC
R_{free} test set	4780 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	48.4	Xtriage
Anisotropy	0.430	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.010 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12991	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 3.18% 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	0.80	0/2162	0.96	1/2914 (0.0%)
1	B	0.72	0/2221	0.94	0/2994
1	C	0.75	0/2162	0.91	0/2914
1	D	0.66	0/2162	0.95	1/2914 (0.0%)
1	E	0.69	0/2162	0.95	2/2914 (0.1%)
1	F	0.79	0/2227	0.95	1/3002 (0.0%)
All	All	0.74	0/13096	0.94	5/17652 (0.0%)

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	D	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	62	LYS	N-CA-C	-5.65	102.91	111.56
1	F	98	ILE	CB-CA-C	-5.31	105.08	111.88
1	A	185	PHE	CB-CA-C	-5.31	110.43	116.54
1	E	88	ALA	CA-C-N	5.15	127.40	120.50
1	E	88	ALA	C-N-CA	5.15	127.40	120.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	61	THR	Peptide

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	2132	0	2209	39	0
1	B	2189	0	2267	45	0
1	C	2132	0	2209	41	0
1	D	2132	0	2209	92	0
1	E	2132	0	2209	73	0
1	F	2195	0	2272	34	0
2	A	23	0	0	0	0
2	B	8	0	0	0	0
2	C	16	0	0	1	0
2	D	4	0	0	0	0
2	E	8	0	0	0	0
2	F	20	0	0	2	0
All	All	12991	0	13375	300	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 11.

All (300) 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:194:MET:HE2	1:C:194:MET:SD	1.63	1.38
1:B:194:MET:CE	1:C:194:MET:SD	2.24	1.25
1:D:194:MET:SD	1:E:194:MET:CE	2.34	1.16
1:D:102:LEU:CD2	1:D:112:LEU:HD22	1.82	1.10
1:D:102:LEU:HD23	1:D:112:LEU:HD22	1.35	1.05
1:D:30:ARG:HD2	1:D:30:ARG:C	1.81	1.02
1:C:86:GLU:OE1	1:C:95:LYS:HE3	1.61	1.01
1:E:29:LEU:HD12	1:E:30:ARG:N	1.77	0.99
1:E:29:LEU:HD12	1:E:29:LEU:C	1.85	0.98
1:E:46:LYS:HE2	1:E:50:LYS:HE3	1.45	0.95
1:D:30:ARG:HD2	1:D:31:ASP:N	1.82	0.93
1:D:194:MET:SD	1:E:194:MET:HE1	2.08	0.93
1:D:102:LEU:CD2	1:D:112:LEU:CD2	2.47	0.92
1:E:41:LEU:HD12	1:E:41:LEU:H	1.38	0.89
1:D:194:MET:SD	1:E:194:MET:HE2	2.11	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:86:GLU:OE1	1:D:95:LYS:NZ	2.06	0.87
1:D:102:LEU:HD21	1:D:112:LEU:CD2	2.05	0.86
1:A:89:VAL:O	1:A:95:LYS:NZ	2.08	0.86
1:D:10:GLY:O	1:D:14:SER:OG	1.94	0.85
1:C:40:GLY:O	1:C:44:ILE:HG13	1.79	0.83
1:B:1:MET:HE3	1:B:168:THR:HG21	1.61	0.82
1:E:86:GLU:OE2	1:E:95:LYS:HE3	1.80	0.81
1:E:37:VAL:O	1:E:41:LEU:HD12	1.81	0.80
1:A:164:ALA:O	1:A:168:THR:CG2	2.30	0.79
1:D:53:LYS:HG3	1:D:53:LYS:O	1.83	0.78
1:A:86:GLU:CD	1:A:95:LYS:HE3	2.08	0.77
1:B:143:ALA:O	1:B:174:LYS:NZ	2.18	0.77
1:D:86:GLU:OE2	1:D:88:ALA:HB3	1.86	0.75
1:D:232:PHE:HD1	1:D:271:ARG:NH1	1.85	0.75
1:A:86:GLU:OE1	1:A:95:LYS:HE3	1.88	0.73
1:D:81:CYS:O	1:D:110:THR:HG23	1.87	0.73
1:E:29:LEU:C	1:E:29:LEU:CD1	2.59	0.73
1:E:10:GLY:O	1:E:14:SER:OG	2.08	0.72
1:D:232:PHE:CD1	1:D:271:ARG:NH1	2.58	0.72
1:A:164:ALA:O	1:A:168:THR:HG22	1.90	0.71
1:F:-1:SER:O	1:F:2:LYS:NZ	2.24	0.70
1:D:30:ARG:CD	1:D:31:ASP:N	2.54	0.70
1:D:81:CYS:O	1:D:110:THR:CG2	2.39	0.70
1:B:95:LYS:HE2	1:B:119:LEU:HD12	1.72	0.70
1:D:107:LYS:O	1:D:110:THR:OG1	2.09	0.70
1:B:107:LYS:HD2	1:B:109:GLU:HG2	1.75	0.69
1:F:-3:ARG:HG2	1:F:-3:ARG:HH11	1.58	0.69
1:D:164:ALA:O	1:D:168:THR:HG22	1.92	0.69
1:E:41:LEU:HD12	1:E:41:LEU:N	2.07	0.68
1:B:163:ASP:O	1:B:167:GLU:HG3	1.93	0.68
1:D:102:LEU:HD12	1:D:106:CYS:SG	2.34	0.68
1:E:37:VAL:O	1:E:41:LEU:CD1	2.41	0.68
1:C:8:GLY:O	1:C:13:GLY:HA3	1.94	0.68
1:A:22:ALA:HA	1:A:65:ILE:HG12	1.77	0.67
1:D:73:VAL:HG23	1:D:73:VAL:O	1.92	0.67
1:E:143:ALA:O	1:E:174:LYS:NZ	2.24	0.66
1:B:194:MET:HE1	1:C:194:MET:SD	2.31	0.66
1:D:102:LEU:CD1	1:D:106:CYS:SG	2.84	0.66
1:D:232:PHE:HD1	1:D:271:ARG:HH11	1.43	0.65
1:C:86:GLU:OE2	1:C:88:ALA:N	2.27	0.65
1:E:203:ALA:HB2	1:E:259:LEU:HD12	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:30:ARG:C	1:D:30:ARG:CD	2.64	0.64
1:A:52:VAL:HG12	1:A:57:ILE:O	1.97	0.64
1:B:217:LYS:HG2	1:B:222:HIS:O	1.97	0.64
1:E:72:THR:HG22	1:E:73:VAL:N	2.12	0.64
1:A:52:VAL:HG11	1:A:59:GLU:HA	1.80	0.63
1:D:3:LYS:HD2	1:D:81:CYS:SG	2.38	0.63
1:A:280:TYR:O	1:A:281:SER:C	2.41	0.63
1:A:59:GLU:O	1:A:59:GLU:HG3	1.99	0.63
1:E:86:GLU:CD	1:E:95:LYS:HE3	2.24	0.62
1:B:1:MET:HE3	1:B:168:THR:CG2	2.29	0.62
1:B:49:SER:OG	1:B:62:LYS:NZ	2.31	0.62
1:E:201:ILE:HG23	1:E:206:ILE:HB	1.80	0.62
1:D:52:VAL:HG12	1:D:57:ILE:HG13	1.80	0.61
1:A:194:MET:SD	1:F:194:MET:SD	2.98	0.61
1:F:232:PHE:CD1	1:F:271:ARG:NH1	2.68	0.61
1:D:130:ARG:N	1:D:131:PRO:CD	2.64	0.61
1:D:159:GLN:O	1:D:159:GLN:HG3	1.99	0.61
1:F:1:MET:HE1	1:F:111:ILE:HD12	1.82	0.61
1:A:164:ALA:O	1:A:168:THR:HG23	2.00	0.61
1:E:59:GLU:O	1:E:63:VAL:HG23	2.01	0.61
1:E:63:VAL:O	1:E:67:THR:HG23	2.01	0.61
1:E:3:LYS:HG3	1:E:26:GLU:HG2	1.82	0.60
1:D:30:ARG:HD2	1:D:31:ASP:CA	2.32	0.60
1:D:100:ALA:O	1:D:103:ASP:HB3	2.02	0.60
1:E:107:LYS:HD2	1:E:109:GLU:HG2	1.84	0.59
1:F:1:MET:CE	1:F:111:ILE:HD12	2.32	0.59
1:B:122:THR:HG23	1:B:155:ILE:HG22	1.82	0.59
1:D:201:ILE:HG23	1:D:206:ILE:HB	1.84	0.59
1:D:111:ILE:HD13	1:D:161:THR:HG23	1.84	0.59
1:A:143:ALA:O	1:A:174:LYS:NZ	2.30	0.59
1:E:49:SER:HA	1:E:52:VAL:HG23	1.84	0.58
1:E:91:ARG:HD2	1:E:94:ILE:HD13	1.86	0.58
1:D:165:VAL:HA	1:D:168:THR:HG23	1.85	0.58
1:E:72:THR:HG22	1:E:74:ASP:N	2.19	0.58
1:D:169:SER:HB2	1:D:176:PRO:HG3	1.86	0.57
1:A:86:GLU:OE2	1:A:95:LYS:HE2	2.05	0.57
1:B:164:ALA:O	1:B:168:THR:HG23	2.03	0.57
1:D:214:LYS:HE3	1:E:175:ASP:OD2	2.05	0.56
1:C:272:LYS:HG2	2:C:308:HOH:O	2.05	0.56
1:D:61:THR:HA	1:D:64:GLU:HB3	1.89	0.55
1:D:137:MET:HE3	1:D:165:VAL:HG13	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:73:VAL:O	1:D:73:VAL:CG2	2.54	0.55
1:E:46:LYS:HE2	1:E:50:LYS:CE	2.29	0.55
1:A:86:GLU:CD	1:A:95:LYS:CE	2.80	0.54
1:E:259:LEU:HD21	1:E:263:TYR:HE1	1.72	0.54
1:C:95:LYS:HD2	1:C:119:LEU:CD1	2.37	0.54
1:C:95:LYS:HD2	1:C:119:LEU:HD12	1.89	0.54
1:A:8:GLY:O	1:A:13:GLY:HA3	2.08	0.54
1:A:86:GLU:OE2	1:A:95:LYS:CE	2.56	0.54
1:B:109:GLU:HA	1:B:133:LYS:HE2	1.91	0.53
1:E:72:THR:HG22	1:E:74:ASP:H	1.74	0.53
1:C:135:ILE:HG21	1:C:165:VAL:HG21	1.89	0.53
1:E:40:GLY:O	1:E:43:PHE:HB3	2.07	0.53
1:E:49:SER:CA	1:E:52:VAL:HG23	2.39	0.53
1:F:1:MET:HE2	1:F:82:ASP:O	2.08	0.53
1:D:91:ARG:HG2	1:D:93:ASP:OD1	2.09	0.53
1:D:143:ALA:O	1:D:174:LYS:NZ	2.33	0.53
1:E:48:LEU:HA	1:E:51:LEU:HD12	1.90	0.53
1:B:50:LYS:O	1:B:54:LYS:HG3	2.09	0.52
1:B:113:ALA:HA	1:B:135:ILE:O	2.10	0.52
1:D:30:ARG:CG	1:D:30:ARG:HH11	2.22	0.52
1:C:96:LYS:HG2	1:C:127:ALA:HB2	1.92	0.52
1:F:86:GLU:OE1	1:F:88:ALA:N	2.31	0.52
1:E:259:LEU:CD2	1:E:263:TYR:CE1	2.92	0.52
1:A:31:ASP:O	1:A:73:VAL:HA	2.10	0.52
1:E:41:LEU:H	1:E:41:LEU:CD1	2.17	0.52
1:C:30:ARG:HA	1:C:72:THR:O	2.09	0.52
1:C:22:ALA:HA	1:C:65:ILE:HG12	1.92	0.51
1:D:102:LEU:HD11	1:D:106:CYS:SG	2.50	0.51
1:D:138:HIS:HB3	1:D:150:GLU:HB2	1.92	0.51
1:E:30:ARG:HD3	1:E:31:ASP:O	2.09	0.51
1:E:130:ARG:O	1:E:130:ARG:HG3	2.10	0.51
1:E:49:SER:HA	1:E:52:VAL:CG2	2.41	0.51
1:A:194:MET:HG3	1:F:194:MET:HG3	1.92	0.51
1:B:214:LYS:HE3	1:C:175:ASP:OD1	2.10	0.51
1:C:201:ILE:HG23	1:C:206:ILE:HB	1.93	0.51
1:F:192:ILE:HG13	1:F:241:ILE:HG21	1.93	0.51
1:F:113:ALA:HA	1:F:135:ILE:O	2.11	0.51
1:C:1:MET:HE2	1:C:82:ASP:HB3	1.93	0.50
1:D:192:ILE:HB	1:D:193:PRO:HD3	1.92	0.50
1:E:91:ARG:HD2	1:E:94:ILE:CD1	2.41	0.50
1:C:97:GLN:HA	1:C:97:GLN:NE2	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:LYS:HD2	1:B:109:GLU:CG	2.40	0.50
1:E:192:ILE:HB	1:E:193:PRO:HD3	1.93	0.50
1:B:115:ASN:CG	1:B:115:ASN:O	2.55	0.50
1:A:175:ASP:CG	1:F:214:LYS:HE3	2.36	0.50
1:B:194:MET:CE	1:C:194:MET:CG	2.89	0.49
1:C:1:MET:HE2	1:C:82:ASP:CB	2.42	0.49
1:D:62:LYS:O	1:D:66:LEU:HD12	2.13	0.49
1:A:140:PHE:CE2	1:A:148:LEU:HD23	2.48	0.49
1:B:123:GLU:HG3	1:F:265:ARG:HG2	1.93	0.49
1:D:169:SER:CB	1:D:176:PRO:HG3	2.43	0.49
1:C:30:ARG:HD3	1:C:31:ASP:N	2.27	0.49
1:F:-3:ARG:HH11	1:F:-3:ARG:CG	2.25	0.49
1:D:45:ASN:O	1:D:49:SER:OG	2.26	0.49
1:E:259:LEU:CD2	1:E:263:TYR:HE1	2.25	0.49
1:A:191:LEU:O	1:A:194:MET:HB2	2.13	0.48
1:B:52:VAL:HA	1:B:57:ILE:O	2.12	0.48
1:F:140:PHE:CE1	1:F:148:LEU:HD23	2.48	0.48
1:D:52:VAL:CG1	1:D:57:ILE:HG13	2.44	0.48
1:D:40:GLY:O	1:D:44:ILE:HG13	2.14	0.48
1:D:52:VAL:HG12	1:D:57:ILE:CG1	2.43	0.48
1:D:130:ARG:N	1:D:131:PRO:HD2	2.29	0.48
1:D:192:ILE:HG13	1:D:241:ILE:HG21	1.95	0.48
1:F:172:ILE:HG13	1:F:174:LYS:HG2	1.95	0.48
1:F:138:HIS:HB3	1:F:150:GLU:HB2	1.96	0.48
1:E:49:SER:OG	1:E:62:LYS:NZ	2.47	0.48
1:C:31:ASP:O	1:C:73:VAL:HA	2.14	0.47
1:B:143:ALA:N	1:B:144:PRO:CD	2.77	0.47
1:A:89:VAL:O	1:A:89:VAL:CG2	2.62	0.47
1:B:86:GLU:OE2	1:B:88:ALA:N	2.46	0.47
1:A:143:ALA:N	1:A:144:PRO:CD	2.78	0.47
1:D:74:ASP:HB3	1:D:77:MET:CG	2.44	0.47
1:E:28:VAL:HG12	1:E:29:LEU:N	2.30	0.47
1:E:253:LYS:HD3	1:E:254:TYR:CZ	2.50	0.47
1:F:147:LYS:HB3	1:F:147:LYS:HE2	1.55	0.47
1:B:268:TRP:CZ3	1:D:155:ILE:HD13	2.50	0.47
1:C:92:MET:HE3	1:C:123:GLU:CD	2.40	0.47
1:D:30:ARG:HD3	1:D:31:ASP:O	2.15	0.47
1:D:102:LEU:HG	1:D:112:LEU:HD21	1.96	0.47
1:D:133:LYS:HE3	1:D:133:LYS:HB3	1.53	0.47
1:D:194:MET:SD	1:E:194:MET:HE3	2.46	0.46
1:E:61:THR:O	1:E:64:GLU:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253:LYS:HD3	1:B:254:TYR:CZ	2.49	0.46
1:D:72:THR:OG1	1:D:73:VAL:N	2.46	0.46
1:D:133:LYS:HA	1:D:161:THR:HG21	1.96	0.46
1:D:74:ASP:HB3	1:D:77:MET:HG2	1.96	0.46
1:D:93:ASP:OD1	1:D:93:ASP:N	2.48	0.46
1:D:140:PHE:O	1:D:146:MET:HB2	2.15	0.46
1:E:38:ASP:O	1:E:42:ASP:HB3	2.14	0.46
1:B:201:ILE:HG23	1:B:206:ILE:HB	1.98	0.46
1:E:37:VAL:HG12	1:E:41:LEU:HD11	1.97	0.46
1:B:30:ARG:HD3	1:B:31:ASP:N	2.31	0.46
1:C:143:ALA:O	1:C:174:LYS:NZ	2.37	0.46
1:E:79:ALA:HA	1:E:106:CYS:HA	1.98	0.46
1:B:122:THR:CG2	1:B:155:ILE:HG22	2.45	0.45
1:F:201:ILE:HG23	1:F:206:ILE:HB	1.97	0.45
1:F:268:TRP:CD2	1:F:275:LYS:HD3	2.50	0.45
1:B:155:ILE:HG21	1:F:266:ALA:HB2	1.98	0.45
1:A:86:GLU:OE1	1:A:95:LYS:CE	2.62	0.45
1:A:113:ALA:HA	1:A:135:ILE:O	2.16	0.45
1:A:138:HIS:HB3	1:A:150:GLU:HB2	1.97	0.45
1:D:164:ALA:O	1:D:168:THR:CG2	2.61	0.45
1:F:135:ILE:HG21	1:F:165:VAL:HG21	1.99	0.45
1:D:102:LEU:HD21	1:D:112:LEU:HD23	1.95	0.45
1:F:96:LYS:HG2	1:F:127:ALA:HB2	1.98	0.45
1:B:-5:VAL:HG13	1:B:-4:PRO:HD2	1.99	0.45
1:E:91:ARG:CG	1:E:93:ASP:OD1	2.65	0.45
1:B:192:ILE:HG13	1:B:241:ILE:HG21	1.99	0.45
1:E:90:GLU:OE2	1:E:116:THR:OG1	2.31	0.45
1:E:30:ARG:HH12	1:E:32:ILE:HD12	1.81	0.44
1:E:72:THR:CG2	1:E:73:VAL:N	2.78	0.44
1:E:49:SER:O	1:E:52:VAL:HG23	2.17	0.44
1:B:-1:SER:HB2	1:B:168:THR:HA	1.98	0.44
1:F:160:GLU:HG3	2:F:315:HOH:O	2.16	0.44
1:B:92:MET:HG3	1:B:96:LYS:HD2	1.99	0.44
1:C:132:ASP:HB2	1:C:158:SER:HB3	1.99	0.44
1:D:3:LYS:HG2	1:D:26:GLU:HB3	2.00	0.44
1:E:30:ARG:CZ	1:E:75:LEU:HD23	2.48	0.44
1:E:49:SER:C	1:E:52:VAL:HG23	2.42	0.44
1:A:194:MET:HG3	1:F:194:MET:CG	2.48	0.44
1:D:52:VAL:CG1	1:D:57:ILE:CG1	2.96	0.44
1:D:34:ASP:O	1:D:35:GLU:C	2.60	0.44
1:E:86:GLU:OE2	1:E:95:LYS:CE	2.60	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:61:THR:O	1:E:62:LYS:C	2.61	0.44
1:F:-3:ARG:CG	1:F:-3:ARG:NH1	2.81	0.44
1:B:191:LEU:O	1:B:194:MET:HB3	2.18	0.43
1:D:214:LYS:HE3	1:E:175:ASP:CG	2.42	0.43
1:B:189:ARG:HG2	1:C:201:ILE:HD11	1.99	0.43
1:D:30:ARG:HH11	1:D:30:ARG:HG2	1.83	0.43
1:B:192:ILE:N	1:B:193:PRO:CD	2.81	0.43
1:C:268:TRP:CD2	1:C:275:LYS:HD3	2.53	0.43
1:C:5:CYS:HB2	1:C:81:CYS:SG	2.59	0.43
1:D:105:ILE:O	1:D:105:ILE:HG22	2.19	0.43
1:A:101:ASP:O	1:A:105:ILE:HG13	2.18	0.43
1:D:54:LYS:C	1:D:56:LYS:H	2.27	0.43
1:F:130:ARG:HG3	1:F:130:ARG:O	2.19	0.43
1:E:240:ALA:O	1:E:244:VAL:HG23	2.19	0.43
1:A:33:LYS:O	1:A:37:VAL:HG23	2.19	0.43
1:A:89:VAL:O	1:A:89:VAL:HG23	2.18	0.43
1:D:209:VAL:HB	1:D:227:LEU:HD13	1.99	0.43
1:F:160:GLU:HG3	1:F:160:GLU:H	1.58	0.43
1:A:30:ARG:HD3	1:A:30:ARG:C	2.44	0.43
1:B:91:ARG:HD2	1:B:94:ILE:CD1	2.48	0.43
1:C:86:GLU:OE2	1:C:88:ALA:HB3	2.18	0.43
1:D:200:GLY:O	1:D:204:GLU:HG3	2.19	0.42
1:A:10:GLY:O	1:A:14:SER:OG	2.36	0.42
1:B:39:ARG:HE	1:B:39:ARG:HB3	1.48	0.42
1:C:62:LYS:O	1:C:62:LYS:HG3	2.19	0.42
1:C:141:ASN:OD1	1:C:142:PRO:CA	2.67	0.42
1:E:21:ALA:C	1:E:23:LYS:H	2.26	0.42
1:F:200:GLY:HA2	1:F:257:HIS:CD2	2.54	0.42
1:D:190:ILE:CG2	1:E:194:MET:HG3	2.50	0.42
1:B:194:MET:HE3	1:C:194:MET:CG	2.49	0.42
1:E:72:THR:HG21	1:E:77:MET:CG	2.49	0.42
1:D:86:GLU:OE2	1:D:88:ALA:CB	2.64	0.42
1:D:162:PHE:CE1	1:D:178:GLU:HG3	2.54	0.42
1:A:206:ILE:N	1:A:206:ILE:CD1	2.83	0.42
1:D:56:LYS:HD2	1:D:56:LYS:HA	1.58	0.42
1:A:130:ARG:O	1:A:130:ARG:HG3	2.19	0.41
1:C:42:ASP:O	1:C:46:LYS:HG3	2.20	0.41
1:D:30:ARG:CG	1:D:30:ARG:NH1	2.83	0.41
1:E:30:ARG:NH2	1:E:75:LEU:HD23	2.33	0.41
1:F:129:LYS:HA	1:F:129:LYS:HD3	1.78	0.41
1:E:78:ALA:O	1:E:79:ALA:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:ASN:ND2	1:A:66:LEU:HD11	2.35	0.41
1:D:129:LYS:C	1:D:131:PRO:HD2	2.44	0.41
1:F:30:ARG:HD3	1:F:30:ARG:C	2.45	0.41
1:C:97:GLN:NE2	1:C:97:GLN:CA	2.84	0.41
1:C:97:GLN:CA	1:C:97:GLN:HE21	2.32	0.41
1:C:141:ASN:OD1	1:C:141:ASN:C	2.63	0.41
1:D:102:LEU:HD21	1:D:112:LEU:HD22	1.72	0.41
1:E:22:ALA:HA	1:E:65:ILE:HG12	2.02	0.41
1:D:77:MET:HE2	1:D:77:MET:HA	2.03	0.41
1:A:204:GLU:CB	1:A:206:ILE:HD13	2.50	0.41
1:E:143:ALA:N	1:E:144:PRO:HD2	2.35	0.41
1:B:261:LYS:NZ	1:D:248:GLU:O	2.49	0.41
1:C:138:HIS:HB3	1:C:150:GLU:HB2	2.03	0.41
1:D:30:ARG:CD	1:D:31:ASP:O	2.69	0.41
1:C:141:ASN:OD1	1:C:142:PRO:N	2.53	0.41
1:D:105:ILE:O	1:D:105:ILE:CG2	2.67	0.41
1:F:191:LEU:O	1:F:194:MET:HB2	2.21	0.41
1:D:137:MET:HE3	1:D:165:VAL:CG1	2.51	0.41
1:E:72:THR:HG21	1:E:77:MET:HG3	2.03	0.41
1:E:74:ASP:OD1	1:E:76:ASN:ND2	2.35	0.41
1:A:175:ASP:OD1	1:F:214:LYS:HE3	2.20	0.40
1:D:30:ARG:HD3	1:D:73:VAL:HA	2.03	0.40
1:F:1:MET:HE2	2:F:303:HOH:O	2.21	0.40
1:A:180:ALA:O	1:A:181:GLU:C	2.64	0.40
1:B:31:ASP:O	1:B:73:VAL:HA	2.21	0.40
1:D:28:VAL:HG12	1:D:29:LEU:N	2.36	0.40
1:E:203:ALA:CB	1:E:259:LEU:HD12	2.49	0.40
1:E:236:ASP:OD1	1:E:236:ASP:N	2.54	0.40
1:D:191:LEU:O	1:D:194:MET:HB3	2.21	0.40
1:E:37:VAL:CG1	1:E:41:LEU:HD11	2.50	0.40
1:B:214:LYS:HE3	1:C:175:ASP:CG	2.46	0.40
1:C:30:ARG:HD3	1:C:30:ARG:C	2.47	0.40
1:C:142:PRO:HD2	1:C:146:MET:HG2	2.02	0.40
1:D:97:GLN:O	1:D:98:ILE:C	2.63	0.40
1:B:89:VAL:HG22	1:B:90:GLU:H	1.87	0.40
1:B:118:SER:HB2	1:B:244:VAL:HG21	2.04	0.40
1:D:183:PRO:O	1:D:189:ARG:HD3	2.22	0.40
1:E:118:SER:HB2	1:E:244:VAL:HG21	2.03	0.40
1:E:147:LYS:HD3	1:E:175:ASP:HB2	2.02	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	279/302 (92%)	269 (96%)	10 (4%)	0	100	100
1	B	287/302 (95%)	280 (98%)	7 (2%)	0	100	100
1	C	279/302 (92%)	267 (96%)	12 (4%)	0	100	100
1	D	279/302 (92%)	261 (94%)	16 (6%)	2 (1%)	18	34
1	E	279/302 (92%)	263 (94%)	16 (6%)	0	100	100
1	F	288/302 (95%)	283 (98%)	5 (2%)	0	100	100
All	All	1691/1812 (93%)	1623 (96%)	66 (4%)	2 (0%)	48	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	22	ALA
1	D	108	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	A	231/249 (93%)	204 (88%)	27 (12%)	5	11
1	B	237/249 (95%)	205 (86%)	32 (14%)	4	8
1	C	231/249 (93%)	207 (90%)	24 (10%)	7	14
1	D	231/249 (93%)	198 (86%)	33 (14%)	3	6
1	E	231/249 (93%)	201 (87%)	30 (13%)	4	8

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	238/249 (96%)	212 (89%)	26 (11%)	6	13
All	All	1399/1494 (94%)	1227 (88%)	172 (12%)	4	10

All (172) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	14	SER
1	A	23	LYS
1	A	35	GLU
1	A	52	VAL
1	A	67	THR
1	A	89	VAL
1	A	95	LYS
1	A	97	GLN
1	A	101	ASP
1	A	109	GLU
1	A	117	SER
1	A	118	SER
1	A	129	LYS
1	A	133	LYS
1	A	137	MET
1	A	167	GLU
1	A	168	THR
1	A	194	MET
1	A	199	VAL
1	A	206	ILE
1	A	209	VAL
1	A	252	SER
1	A	258	THR
1	A	272	LYS
1	A	279	ASP
1	A	281	SER
1	B	7	ILE
1	B	26	GLU
1	B	32	ILE
1	B	39	ARG
1	B	50	LYS
1	B	52	VAL
1	B	53	LYS
1	B	57	ILE
1	B	58	GLU

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Mol	Chain	Res	Type
1	B	61	THR
1	B	62	LYS
1	B	67	THR
1	B	70	SER
1	B	75	LEU
1	B	94	ILE
1	B	95	LYS
1	B	98	ILE
1	B	107	LYS
1	B	109	GLU
1	B	137	MET
1	B	167	GLU
1	B	168	THR
1	B	199	VAL
1	B	204	GLU
1	B	206	ILE
1	B	209	VAL
1	B	214	LYS
1	B	247	SER
1	B	252	SER
1	B	253	LYS
1	B	258	THR
1	B	275	LYS
1	C	23	LYS
1	C	56	LYS
1	C	58	GLU
1	C	62	LYS
1	C	89	VAL
1	C	91	ARG
1	C	94	ILE
1	C	96	LYS
1	C	97	GLN
1	C	107	LYS
1	C	137	MET
1	C	149	VAL
1	C	155	ILE
1	C	170	ILE
1	C	189	ARG
1	C	194	MET
1	C	199	VAL
1	C	209	VAL
1	C	237	ILE

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Mol	Chain	Res	Type
1	C	252	SER
1	C	258	THR
1	C	259	LEU
1	C	271	ARG
1	C	273	SER
1	D	1	MET
1	D	3	LYS
1	D	4	VAL
1	D	14	SER
1	D	30	ARG
1	D	44	ILE
1	D	46	LYS
1	D	50	LYS
1	D	53	LYS
1	D	54	LYS
1	D	56	LYS
1	D	59	GLU
1	D	61	THR
1	D	62	LYS
1	D	63	VAL
1	D	70	SER
1	D	73	VAL
1	D	75	LEU
1	D	94	ILE
1	D	96	LYS
1	D	98	ILE
1	D	107	LYS
1	D	110	THR
1	D	126	SER
1	D	129	LYS
1	D	133	LYS
1	D	155	ILE
1	D	160	GLU
1	D	168	THR
1	D	199	VAL
1	D	204	GLU
1	D	209	VAL
1	D	258	THR
1	E	2	LYS
1	E	26	GLU
1	E	29	LEU
1	E	32	ILE

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Mol	Chain	Res	Type
1	E	41	LEU
1	E	46	LYS
1	E	47	ASN
1	E	52	VAL
1	E	64	GLU
1	E	67	THR
1	E	70	SER
1	E	75	LEU
1	E	91	ARG
1	E	94	ILE
1	E	96	LYS
1	E	98	ILE
1	E	109	GLU
1	E	117	SER
1	E	129	LYS
1	E	147	LYS
1	E	149	VAL
1	E	189	ARG
1	E	199	VAL
1	E	209	VAL
1	E	210	GLU
1	E	247	SER
1	E	258	THR
1	E	259	LEU
1	E	279	ASP
1	E	281	SER
1	F	-8	SER
1	F	-3	ARG
1	F	3	LYS
1	F	23	LYS
1	F	39	ARG
1	F	41	LEU
1	F	46	LYS
1	F	54	LYS
1	F	56	LYS
1	F	63	VAL
1	F	89	VAL
1	F	91	ARG
1	F	94	ILE
1	F	96	LYS
1	F	129	LYS
1	F	137	MET

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Mol	Chain	Res	Type
1	F	149	VAL
1	F	160	GLU
1	F	189	ARG
1	F	194	MET
1	F	199	VAL
1	F	209	VAL
1	F	214	LYS
1	F	247	SER
1	F	252	SER
1	F	258	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	45	ASN
1	B	141	ASN
1	B	222	HIS
1	C	97	GLN
1	D	97	GLN
1	E	45	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

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	A	281/302 (93%)	-0.35	5 (1%) 67 64	28, 41, 71, 92	0
1	B	289/302 (95%)	-0.01	6 (2%) 63 59	33, 51, 77, 101	0
1	C	281/302 (93%)	-0.13	3 (1%) 78 75	30, 49, 74, 94	0
1	D	281/302 (93%)	0.52	25 (8%) 15 13	33, 67, 108, 125	0
1	E	281/302 (93%)	0.19	10 (3%) 46 41	33, 58, 94, 111	0
1	F	290/302 (96%)	-0.33	3 (1%) 79 76	30, 41, 65, 82	0
All	All	1703/1812 (93%)	-0.02	52 (3%) 51 47	28, 48, 92, 125	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	89	VAL	7.0
1	D	1	MET	6.3
1	D	61	THR	3.8
1	B	89	VAL	3.6
1	A	89	VAL	3.4
1	D	54	LYS	3.4
1	E	89	VAL	3.3
1	B	0	HIS	3.0
1	D	102	LEU	3.0
1	D	60	ALA	3.0
1	D	26	GLU	2.8
1	D	33	LYS	2.8
1	D	108	PRO	2.7
1	C	43	PHE	2.7
1	A	1	MET	2.7
1	D	80	ASP	2.6
1	D	32	ILE	2.6
1	A	281	SER	2.6
1	F	281	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	72	THR	2.6
1	D	63	VAL	2.6
1	D	36	PHE	2.6
1	B	-8	SER	2.6
1	A	88	ALA	2.5
1	A	54	LYS	2.5
1	F	89	VAL	2.4
1	E	43	PHE	2.4
1	B	26	GLU	2.4
1	E	41	LEU	2.4
1	F	-7	GLY	2.3
1	D	57	ILE	2.3
1	D	129	LYS	2.3
1	D	84	VAL	2.2
1	D	50	LYS	2.2
1	E	281	SER	2.2
1	D	109	GLU	2.2
1	E	32	ILE	2.2
1	D	43	PHE	2.2
1	C	281	SER	2.1
1	D	2	LYS	2.1
1	D	25	PHE	2.1
1	E	57	ILE	2.1
1	D	167	GLU	2.1
1	E	1	MET	2.1
1	E	56	LYS	2.1
1	B	88	ALA	2.1
1	E	80	ASP	2.1
1	B	91	ARG	2.0
1	D	281	SER	2.0
1	C	91	ARG	2.0
1	D	65	ILE	2.0
1	D	105	ILE	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.