



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:35 AM UTC

PDB ID : 7FF4 / pdb_00007ff4
Title : The crystal structure of Clostridium cellulolyticum LacI family transcriptional regulator Ccel_1438
Authors : Zhang, N.; Ge, H.
Deposited on : 2021-07-22
Resolution : 2.19 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
Xtrriage (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

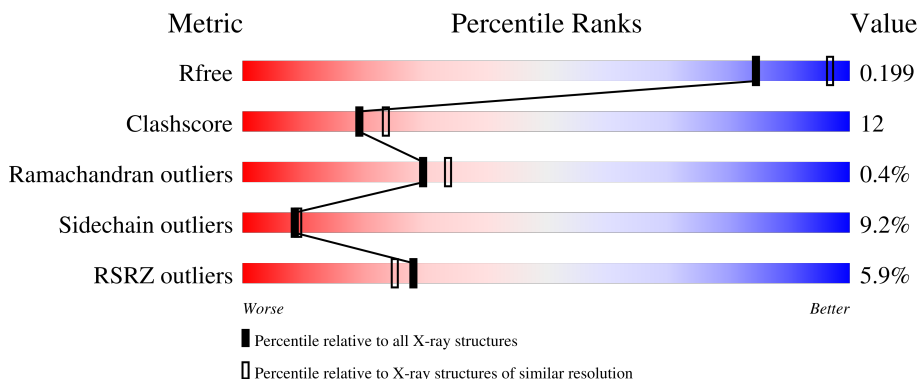
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	343	5% (Poor fit) 62% (0 outliers) 15% (1 outlier) 5% (2 outliers) 18% (Not modelled)
1	B	343	5% (Poor fit) 62% (0 outliers) 16% (1 outlier) • (2 outliers) 19% (Not modelled)
1	C	343	4% (Poor fit) 63% (0 outliers) 13% (1 outlier) 6% (2 outliers) 19% (Not modelled)
1	D	343	6% (Poor fit) 62% (0 outliers) 16% (1 outlier) • (2 outliers) 18% (Not modelled)

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 9043 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 Transcriptional regulator, LacI family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	280	2202	1398	374	422	8	0	0	0
1	B	278	2191	1392	372	419	8	0	0	0
1	C	279	2200	1397	373	421	9	0	1	0
1	D	280	2205	1400	374	422	9	0	1	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	338	HIS	-	expression tag	UNP B8I1W6
A	339	HIS	-	expression tag	UNP B8I1W6
A	340	HIS	-	expression tag	UNP B8I1W6
A	341	HIS	-	expression tag	UNP B8I1W6
A	342	HIS	-	expression tag	UNP B8I1W6
A	343	HIS	-	expression tag	UNP B8I1W6
B	338	HIS	-	expression tag	UNP B8I1W6
B	339	HIS	-	expression tag	UNP B8I1W6
B	340	HIS	-	expression tag	UNP B8I1W6
B	341	HIS	-	expression tag	UNP B8I1W6
B	342	HIS	-	expression tag	UNP B8I1W6
B	343	HIS	-	expression tag	UNP B8I1W6
C	338	HIS	-	expression tag	UNP B8I1W6
C	339	HIS	-	expression tag	UNP B8I1W6
C	340	HIS	-	expression tag	UNP B8I1W6
C	341	HIS	-	expression tag	UNP B8I1W6
C	342	HIS	-	expression tag	UNP B8I1W6
C	343	HIS	-	expression tag	UNP B8I1W6
D	338	HIS	-	expression tag	UNP B8I1W6
D	339	HIS	-	expression tag	UNP B8I1W6
D	340	HIS	-	expression tag	UNP B8I1W6

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Chain	Residue	Modelled	Actual	Comment	Reference
D	341	HIS	-	expression tag	UNP B8I1W6
D	342	HIS	-	expression tag	UNP B8I1W6
D	343	HIS	-	expression tag	UNP B8I1W6

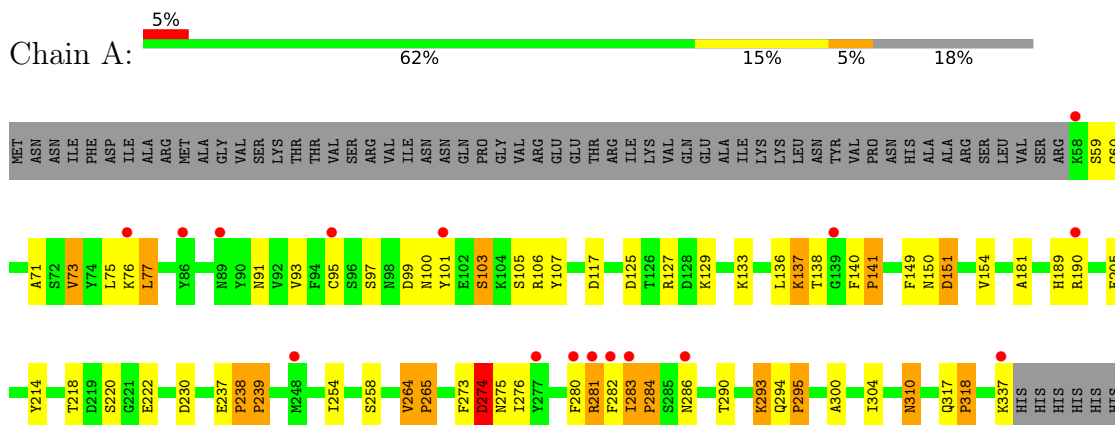
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	89	Total O 89 89	0	0
2	B	53	Total O 53 53	0	0
2	C	62	Total O 62 62	0	0
2	D	41	Total O 41 41	0	0

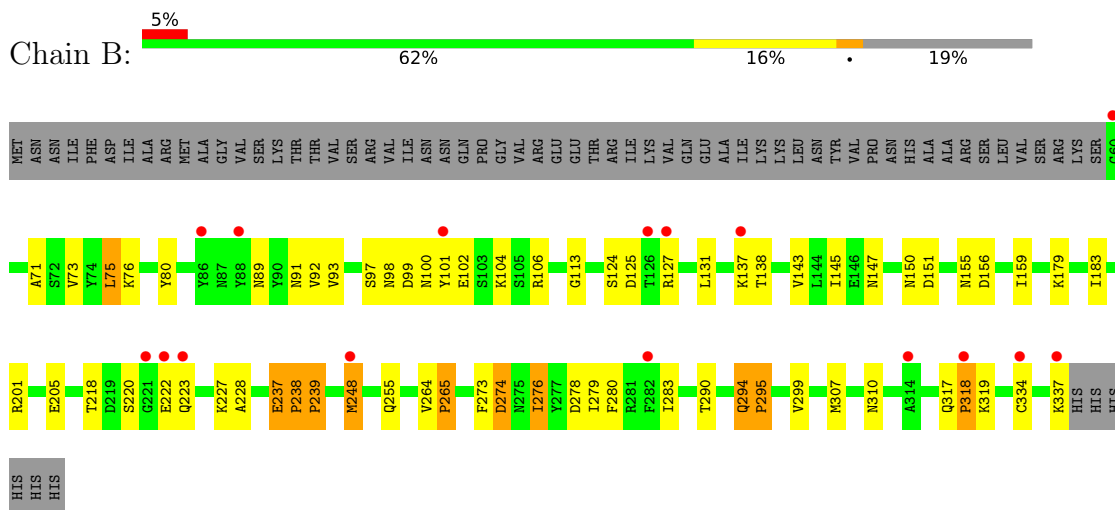
3 Residue-property plots [i](#)

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.

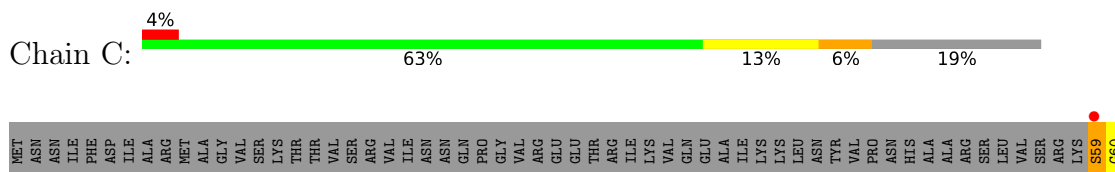
- Molecule 1: Transcriptional regulator, LacI family

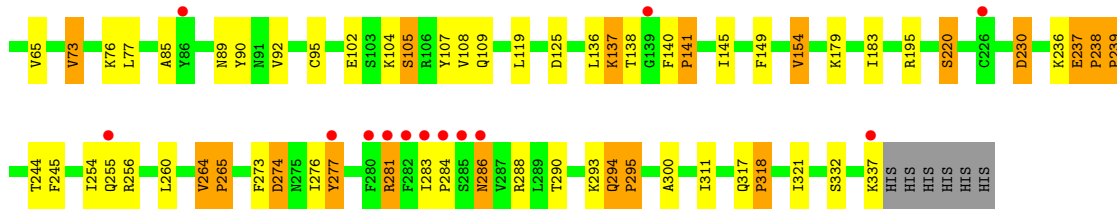


- Molecule 1: Transcriptional regulator, LacI family

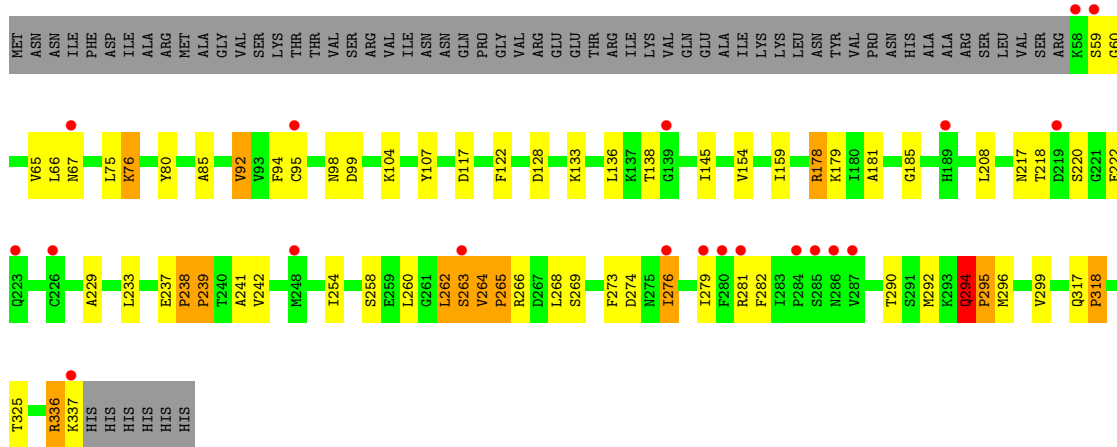


- Molecule 1: Transcriptional regulator, LacI family





● Molecule 1: Transcriptional regulator, LacI family



4 Data and refinement statistics i

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	194.52Å 194.52Å 118.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.63 – 2.19 48.63 – 2.19	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.63-2.19) 99.4 (48.63-2.19)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 2.18Å)	Xtrriage
Refinement program	PHENIX 1.14 3260	Depositor
R, R_{free}	0.178 , 0.208 0.184 , 0.199	Depositor DCC
R_{free} test set	4313 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	48.9	Xtrriage
Anisotropy	0.294	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.006 for $-2/3^*h-1/3^*k-4/3^*l,-1/3^*h-2/3^*k+4/3^*l,-1/3^*h+1/3^*k+1/3^*l$ 0.004 for $-h,1/3^*h-1/3^*k-4/3^*l,-1/3^*h-2/3^*k+1/3^*l$ 0.002 for $-1/3^*h+1/3^*k+4/3^*l,-k,2/3^*h+1/3^*k+1/3^*l$ 0.004 for $-h,2/3^*h+1/3^*k+4/3^*l,1/3^*h+2/3^*k-1/3^*l$ 0.009 for $-1/3^*h-2/3^*k+4/3^*l,-2/3^*h-1/3^*k-4/3^*l,1/3^*h-1/3^*k-1/3^*l$ 0.003 for $1/3^*h+2/3^*k-4/3^*l,-k,-2/3^*h-1/3^*k-1/3^*l$ 0.014 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9043	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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	1.51	13/2242 (0.6%)	1.67	34/3035 (1.1%)
1	B	1.36	9/2231 (0.4%)	1.56	24/3020 (0.8%)
1	C	1.43	10/2243 (0.4%)	1.59	27/3036 (0.9%)
1	D	1.35	9/2248 (0.4%)	1.56	22/3043 (0.7%)
All	All	1.41	41/8964 (0.5%)	1.60	107/12134 (0.9%)

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	318	PRO	N-CA	18.29	1.68	1.47
1	C	318	PRO	N-CA	18.26	1.68	1.47
1	A	318	PRO	N-CA	18.17	1.68	1.47
1	A	239	PRO	N-CA	17.72	1.68	1.47
1	C	239	PRO	N-CA	17.69	1.68	1.47
1	D	239	PRO	N-CA	17.56	1.68	1.47
1	A	141	PRO	N-CA	17.39	1.69	1.47
1	B	239	PRO	N-CA	17.19	1.68	1.47
1	D	295	PRO	N-CA	17.06	1.69	1.47
1	B	295	PRO	N-CA	16.98	1.69	1.47
1	C	295	PRO	N-CA	16.87	1.68	1.47
1	C	141	PRO	N-CA	16.83	1.68	1.47
1	A	295	PRO	N-CA	16.83	1.68	1.47
1	B	318	PRO	N-CA	16.59	1.68	1.47
1	A	284	PRO	N-CA	16.24	1.69	1.47
1	D	265	PRO	N-CA	13.23	1.69	1.47
1	B	265	PRO	N-CA	13.13	1.69	1.47
1	C	265	PRO	N-CA	12.81	1.69	1.47
1	A	265	PRO	N-CA	12.56	1.68	1.47
1	D	238	PRO	N-CA	10.98	1.69	1.46
1	C	238	PRO	N-CA	10.62	1.68	1.46
1	A	238	PRO	N-CA	10.61	1.68	1.46
1	B	238	PRO	N-CA	10.49	1.68	1.46
1	D	237	GLU	C-N	9.68	1.44	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	237	GLU	C-N	9.30	1.44	1.33
1	D	317	GLN	C-N	9.25	1.44	1.33
1	B	237	GLU	C-N	9.21	1.44	1.33
1	A	317	GLN	C-N	9.21	1.44	1.33
1	C	317	GLN	C-N	9.15	1.44	1.33
1	A	237	GLU	C-N	9.12	1.44	1.33
1	D	238	PRO	C-N	8.40	1.44	1.33
1	B	238	PRO	C-N	7.95	1.44	1.33
1	A	238	PRO	C-N	7.91	1.43	1.33
1	C	238	PRO	C-N	7.80	1.43	1.33
1	A	283	ILE	C-N	7.71	1.44	1.34
1	D	264	VAL	C-N	5.58	1.44	1.34
1	B	238	PRO	C-O	-5.39	1.18	1.24
1	A	295	PRO	C-O	-5.38	1.17	1.24
1	B	264	VAL	C-N	5.26	1.44	1.34
1	C	264	VAL	C-N	5.23	1.44	1.34
1	A	264	VAL	C-N	5.18	1.44	1.34

All (107) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	317	GLN	CA-C-N	17.27	138.41	119.93
1	A	317	GLN	C-N-CA	17.27	138.41	119.93
1	D	238	PRO	CA-C-N	17.04	137.31	119.76
1	D	238	PRO	C-N-CA	17.04	137.31	119.76
1	D	317	GLN	CA-C-N	16.94	138.06	119.93
1	D	317	GLN	C-N-CA	16.94	138.06	119.93
1	C	317	GLN	CA-C-N	16.81	137.92	119.93
1	C	317	GLN	C-N-CA	16.81	137.92	119.93
1	C	237	GLU	CA-C-N	16.34	137.22	120.38
1	C	237	GLU	C-N-CA	16.34	137.22	120.38
1	C	238	PRO	CA-C-N	16.28	136.53	119.76
1	C	238	PRO	C-N-CA	16.28	136.53	119.76
1	D	237	GLU	CA-C-N	16.16	137.02	120.38
1	D	237	GLU	C-N-CA	16.16	137.02	120.38
1	A	283	ILE	CA-C-N	16.14	136.97	119.05
1	A	283	ILE	C-N-CA	16.14	136.97	119.05
1	B	237	GLU	CA-C-N	16.04	136.90	120.38
1	B	237	GLU	C-N-CA	16.04	136.90	120.38
1	A	238	PRO	CA-C-N	15.96	136.19	119.76
1	A	238	PRO	C-N-CA	15.96	136.19	119.76
1	B	238	PRO	CA-C-N	15.78	136.01	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	238	PRO	C-N-CA	15.78	136.01	119.76
1	A	237	GLU	CA-C-N	15.73	136.58	120.38
1	A	237	GLU	C-N-CA	15.73	136.58	120.38
1	B	317	GLN	CA-C-N	14.98	138.56	119.84
1	B	317	GLN	C-N-CA	14.98	138.56	119.84
1	B	294	GLN	CA-C-N	14.67	138.18	119.84
1	B	294	GLN	C-N-CA	14.67	138.18	119.84
1	D	294	GLN	CA-C-N	14.43	137.88	119.84
1	D	294	GLN	C-N-CA	14.43	137.88	119.84
1	A	294	GLN	CA-C-N	14.31	137.73	119.84
1	A	294	GLN	C-N-CA	14.31	137.73	119.84
1	C	140	PHE	CA-C-N	13.33	136.50	119.84
1	C	140	PHE	C-N-CA	13.33	136.50	119.84
1	A	140	PHE	CA-C-N	13.32	136.49	119.84
1	A	140	PHE	C-N-CA	13.32	136.49	119.84
1	C	294	GLN	CA-C-N	9.68	137.54	120.88
1	C	294	GLN	C-N-CA	9.68	137.54	120.88
1	B	150	ASN	CB-CA-C	-9.66	95.63	110.90
1	C	238	PRO	N-CA-C	-9.16	99.52	110.70
1	A	318	PRO	CA-N-CD	-8.63	99.91	112.00
1	B	318	PRO	CA-N-CD	-8.47	100.14	112.00
1	C	295	PRO	N-CA-C	-8.07	101.59	112.48
1	A	238	PRO	N-CA-C	-8.06	100.87	110.70
1	D	318	PRO	CA-N-CD	-8.03	100.76	112.00
1	B	238	PRO	N-CA-C	-7.85	101.12	110.70
1	B	239	PRO	N-CA-C	-7.82	98.94	111.14
1	A	239	PRO	N-CA-C	-7.67	99.17	111.14
1	C	141	PRO	N-CA-C	-7.65	96.70	112.47
1	D	239	PRO	N-CA-C	-7.64	99.22	111.14
1	C	239	PRO	N-CA-C	-7.59	99.29	111.14
1	D	239	PRO	CA-N-CD	-7.52	101.47	112.00
1	C	238	PRO	CA-N-CD	-7.45	101.57	112.00
1	C	318	PRO	CA-N-CD	-7.45	101.57	112.00
1	D	265	PRO	CA-N-CD	-7.18	101.44	111.50
1	A	238	PRO	CA-N-CD	-7.11	102.05	112.00
1	B	238	PRO	CA-N-CD	-7.07	102.11	112.00
1	D	238	PRO	CA-N-CD	-6.96	102.26	112.00
1	A	284	PRO	CA-N-CD	-6.88	102.37	112.00
1	A	141	PRO	CA-N-CD	-6.69	102.63	112.00
1	A	295	PRO	CA-N-CD	-6.68	102.65	112.00
1	B	295	PRO	CA-N-CD	-6.66	102.68	112.00
1	C	239	PRO	CA-N-CD	-6.64	102.71	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	273	PHE	CA-C-N	6.64	134.21	121.54
1	C	273	PHE	C-N-CA	6.64	134.21	121.54
1	A	265	PRO	CA-N-CD	-6.61	102.25	111.50
1	C	141	PRO	CA-N-CD	-6.56	102.82	112.00
1	B	239	PRO	CA-N-CD	-6.54	102.84	112.00
1	C	295	PRO	CA-N-CD	-6.54	102.84	112.00
1	B	283	ILE	CB-CA-C	6.52	115.72	109.33
1	D	295	PRO	CA-N-CD	-6.49	102.91	112.00
1	A	318	PRO	N-CA-C	-6.46	100.73	111.26
1	C	295	PRO	CB-CA-C	6.42	121.30	112.89
1	D	238	PRO	N-CA-C	-6.38	102.92	110.70
1	B	113	GLY	CA-C-N	-6.35	115.97	123.08
1	B	113	GLY	C-N-CA	-6.35	115.97	123.08
1	B	265	PRO	CA-N-CD	-6.31	102.66	111.50
1	A	239	PRO	CA-N-CD	-6.25	103.25	112.00
1	A	273	PHE	CA-C-N	6.17	133.32	121.54
1	A	273	PHE	C-N-CA	6.17	133.32	121.54
1	B	273	PHE	CA-C-N	6.01	133.01	121.54
1	B	273	PHE	C-N-CA	6.01	133.01	121.54
1	C	265	PRO	CA-N-CD	-5.92	103.21	111.50
1	C	141	PRO	CB-CA-C	5.84	121.20	111.56
1	B	283	ILE	N-CA-C	-5.77	104.46	109.19
1	D	318	PRO	N-CA-C	-5.76	101.88	111.26
1	A	290	THR	N-CA-C	-5.73	101.37	109.96
1	C	290	THR	N-CA-C	-5.72	101.38	109.96
1	D	273	PHE	CA-C-N	5.65	132.33	121.54
1	D	273	PHE	C-N-CA	5.65	132.33	121.54
1	A	150	ASN	CA-C-N	-5.65	114.76	123.17
1	A	150	ASN	C-N-CA	-5.65	114.76	123.17
1	A	141	PRO	N-CA-C	-5.64	100.85	112.47
1	A	274	ASP	N-CA-CB	5.64	120.02	110.49
1	A	274	ASP	CA-CB-CG	5.47	118.07	112.60
1	A	150	ASN	CB-CA-C	-5.46	99.24	110.38
1	D	290	THR	N-CA-C	-5.39	101.88	109.96
1	A	141	PRO	CB-CA-C	5.36	120.41	111.56
1	C	318	PRO	N-CA-C	-5.29	102.65	111.26
1	B	318	PRO	N-CA-C	-5.25	101.66	112.47
1	B	290	THR	N-CA-C	-5.20	102.16	109.96
1	A	281	ARG	N-CA-C	-5.18	107.48	112.97
1	A	140	PHE	O-C-N	-5.12	116.75	121.30
1	D	264	VAL	CA-C-N	5.11	139.26	127.00
1	D	264	VAL	C-N-CA	5.11	139.26	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	273	PHE	CA-C-O	-5.03	114.36	120.54
1	D	273	PHE	O-C-N	5.00	129.26	123.41

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	2202	0	2165	63	0
1	B	2191	0	2158	53	0
1	C	2200	0	2168	50	0
1	D	2205	0	2170	58	0
2	A	89	0	0	4	0
2	B	53	0	0	0	0
2	C	62	0	0	0	0
2	D	41	0	0	0	0
All	All	9043	0	8661	205	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 12.

All (205) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (\AA)	Clash overlap (\AA)
1:C:265:PRO:CA	1:C:265:PRO:N	1.69	1.48
1:B:238:PRO:CA	1:B:238:PRO:N	1.68	1.47
1:A:284:PRO:N	1:A:284:PRO:CA	1.69	1.46
1:C:238:PRO:N	1:C:238:PRO:CA	1.68	1.46
1:A:318:PRO:N	1:A:318:PRO:CA	1.68	1.45
1:B:295:PRO:N	1:B:295:PRO:CA	1.69	1.44
1:B:265:PRO:N	1:B:265:PRO:CA	1.69	1.42
1:A:141:PRO:N	1:A:141:PRO:CA	1.69	1.42
1:D:295:PRO:N	1:D:295:PRO:CA	1.69	1.42
1:D:318:PRO:N	1:D:318:PRO:CA	1.68	1.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:238:PRO:N	1:D:238:PRO:CA	1.69	1.40
1:A:239:PRO:CA	1:A:239:PRO:N	1.68	1.39
1:C:141:PRO:N	1:C:141:PRO:CA	1.68	1.39
1:C:295:PRO:CA	1:C:295:PRO:N	1.68	1.39
1:C:239:PRO:N	1:C:239:PRO:CA	1.68	1.35
1:B:318:PRO:CA	1:B:318:PRO:N	1.68	1.34
1:D:265:PRO:CA	1:D:265:PRO:N	1.69	1.33
1:C:318:PRO:CA	1:C:318:PRO:N	1.68	1.33
1:B:239:PRO:N	1:B:239:PRO:CA	1.67	1.32
1:A:295:PRO:N	1:A:295:PRO:CA	1.68	1.32
1:D:239:PRO:N	1:D:239:PRO:CA	1.68	1.31
1:A:238:PRO:N	1:A:238:PRO:CA	1.68	1.30
1:A:265:PRO:N	1:A:265:PRO:CA	1.68	1.28
1:A:93:VAL:HG13	1:B:93:VAL:HG13	1.20	1.14
1:C:65:VAL:HA	1:C:95[B]:CYS:SG	1.89	1.12
1:D:229:ALA:O	1:D:233:LEU:HD12	1.54	1.07
1:A:293:LYS:HE2	2:A:465:HOH:O	1.53	1.06
1:D:136:LEU:HD21	1:D:154:VAL:HG12	1.33	1.03
1:D:276:ILE:HG12	1:D:279:ILE:HD12	1.40	1.02
1:D:65:VAL:HA	1:D:95[B]:CYS:SG	2.01	0.99
1:C:65:VAL:HG22	1:C:95[B]:CYS:SG	2.03	0.98
1:A:93:VAL:HG13	1:B:93:VAL:CG1	1.95	0.94
1:D:276:ILE:CG1	1:D:279:ILE:HD12	1.96	0.94
1:A:93:VAL:CG1	1:B:93:VAL:HG13	1.98	0.93
1:D:292:MET:HE2	1:D:325:THR:OG1	1.69	0.90
1:D:136:LEU:HD11	1:D:154:VAL:HG13	1.55	0.88
1:C:65:VAL:HA	1:C:95[B]:CYS:HG	1.38	0.87
1:D:136:LEU:HD11	1:D:154:VAL:CG1	2.07	0.85
1:D:65:VAL:HG22	1:D:95[B]:CYS:SG	2.16	0.84
1:D:292:MET:CE	1:D:325:THR:OG1	2.25	0.83
1:D:336:ARG:HG2	1:D:336:ARG:HH11	1.45	0.81
1:A:136:LEU:HD21	1:A:154:VAL:HG12	1.62	0.81
1:C:137:LYS:HB3	1:C:137:LYS:HZ3	1.46	0.80
1:B:201:ARG:HD2	1:B:205:GLU:OE1	1.82	0.79
1:C:141:PRO:N	1:C:141:PRO:C	2.41	0.79
1:C:286:ASN:OD1	1:C:286:ASN:N	2.15	0.78
1:C:136:LEU:HD11	1:C:154:VAL:HG13	1.66	0.77
1:A:282:PHE:CD1	1:B:248:MET:HG2	2.21	0.76
1:A:136:LEU:HD11	1:A:154:VAL:CG1	2.15	0.76
1:D:76:LYS:HE2	1:D:80:TYR:HE2	1.52	0.75
1:D:239:PRO:N	1:D:239:PRO:C	2.45	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:PRO:N	1:A:239:PRO:C	2.45	0.75
1:B:239:PRO:N	1:B:239:PRO:C	2.44	0.74
1:C:239:PRO:N	1:C:239:PRO:C	2.45	0.74
1:D:254:ILE:HG12	1:D:264:VAL:HG21	1.69	0.74
1:A:282:PHE:CG	1:B:248:MET:HG2	2.22	0.74
1:D:136:LEU:HD21	1:D:154:VAL:CG1	2.15	0.74
1:C:238:PRO:N	1:C:238:PRO:C	2.45	0.73
1:B:265:PRO:HB2	1:B:334:CYS:HB2	1.69	0.73
1:C:137:LYS:HB3	1:C:137:LYS:NZ	2.02	0.73
1:A:275:ASN:HB3	1:A:293:LYS:HE3	1.68	0.73
1:A:318:PRO:N	1:A:318:PRO:C	2.47	0.72
1:C:65:VAL:CA	1:C:95[B]:CYS:SG	2.73	0.72
1:A:282:PHE:HB3	1:B:248:MET:HG2	1.72	0.71
1:B:238:PRO:N	1:B:238:PRO:C	2.48	0.70
1:A:238:PRO:N	1:A:238:PRO:C	2.48	0.70
1:A:141:PRO:N	1:A:141:PRO:C	2.49	0.70
1:B:318:PRO:N	1:B:318:PRO:C	2.49	0.70
1:D:185:GLY:HA2	1:D:220:SER:HB2	1.72	0.70
1:A:91:ASN:OD1	1:B:106:ARG:NH1	2.25	0.70
1:A:205:GLU:OE1	2:A:401:HOH:O	2.09	0.70
1:A:136:LEU:HD11	1:A:154:VAL:HG13	1.75	0.69
1:C:295:PRO:N	1:C:295:PRO:C	2.50	0.68
1:B:179:LYS:HE3	1:B:237:GLU:HG2	1.75	0.68
1:C:281:ARG:HA	1:C:281:ARG:HH11	1.59	0.67
1:D:336:ARG:HH11	1:D:336:ARG:CG	2.08	0.67
1:D:318:PRO:N	1:D:318:PRO:C	2.49	0.67
1:C:136:LEU:HD21	1:C:154:VAL:HG12	1.76	0.66
1:B:295:PRO:N	1:B:295:PRO:C	2.52	0.66
1:A:106:ARG:NH1	1:B:91:ASN:OD1	2.30	0.65
1:C:73:VAL:HG13	1:C:274:ASP:HB3	1.78	0.65
1:D:295:PRO:N	1:D:295:PRO:C	2.54	0.65
1:C:65:VAL:CG2	1:C:95[B]:CYS:SG	2.82	0.65
1:A:282:PHE:CD1	1:B:248:MET:CG	2.79	0.65
1:A:286:ASN:HB3	1:B:255:GLN:NE2	2.11	0.65
1:D:238:PRO:N	1:D:238:PRO:C	2.52	0.65
1:C:318:PRO:N	1:C:318:PRO:C	2.51	0.64
1:A:295:PRO:N	1:A:295:PRO:C	2.53	0.64
1:C:95[A]:CYS:SG	1:C:107:TYR:CD1	2.88	0.63
1:B:76:LYS:HE3	1:B:278:ASP:OD2	1.98	0.63
1:A:93:VAL:CG1	1:B:93:VAL:CG1	2.67	0.63
1:D:178:ARG:HH11	1:D:178:ARG:HG2	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:277:TYR:OH	1:C:284:PRO:HD2	1.99	0.62
1:D:276:ILE:HG13	1:D:279:ILE:HD12	1.80	0.62
1:A:282:PHE:CB	1:B:248:MET:HG2	2.30	0.62
1:C:283:ILE:HG22	1:C:283:ILE:O	2.01	0.60
1:A:73:VAL:HG22	2:A:411:HOH:O	2.00	0.60
1:A:100:ASN:HB3	1:A:103:SER:HB2	1.85	0.59
1:A:151:ASP:OD1	1:A:151:ASP:N	2.36	0.57
1:C:105:SER:O	1:C:109:GLN:HG3	2.04	0.57
1:C:230:ASP:OD1	1:C:256:ARG:NE	2.37	0.57
1:D:65:VAL:CA	1:D:95[B]:CYS:SG	2.87	0.56
1:A:101:TYR:HB2	1:A:127:ARG:HH11	1.71	0.56
1:D:262:LEU:HD12	1:D:263:SER:H	1.71	0.55
1:A:73:VAL:HG13	1:A:274:ASP:HB3	1.88	0.55
1:B:276:ILE:HD12	1:B:279:ILE:HD12	1.88	0.55
1:C:136:LEU:HD11	1:C:154:VAL:CG1	2.36	0.55
1:D:178:ARG:HD3	1:D:208:LEU:HD21	1.89	0.54
1:C:136:LEU:HD21	1:C:154:VAL:CG1	2.37	0.54
1:A:137:LYS:O	1:A:137:LYS:HD3	2.08	0.54
1:C:73:VAL:CG1	1:C:274:ASP:HB3	2.38	0.53
1:D:242:VAL:HG23	1:D:268:LEU:HD11	1.90	0.53
1:A:282:PHE:HA	1:B:222:GLU:HG3	1.90	0.52
1:A:280:PHE:HE1	1:B:280:PHE:CE2	2.27	0.52
1:A:284:PRO:N	1:A:284:PRO:C	2.59	0.52
1:C:277:TYR:CZ	1:C:284:PRO:HD2	2.45	0.52
1:B:220:SER:O	1:B:220:SER:OG	2.26	0.51
1:C:256:ARG:CZ	1:C:260:LEU:HD11	2.40	0.51
1:B:155:ASN:HD22	1:B:319:LYS:H	1.58	0.51
1:A:141:PRO:HA	1:A:310:ASN:HD22	1.77	0.50
1:B:143:VAL:HG23	1:B:307:MET:HE2	1.93	0.50
1:A:280:PHE:CE1	1:B:280:PHE:CE2	2.99	0.50
1:A:254:ILE:HG12	1:A:264:VAL:HG11	1.94	0.49
1:D:85:ALA:CB	1:D:92:VAL:CG2	2.90	0.49
1:B:276:ILE:HD12	1:B:279:ILE:CD1	2.43	0.49
1:B:337:LYS:O	1:B:337:LYS:HG2	2.12	0.48
1:C:195:ARG:NH1	1:C:245:PHE:O	2.46	0.48
1:A:73:VAL:CG1	1:A:274:ASP:HB3	2.43	0.48
1:B:155:ASN:ND2	1:B:319:LYS:H	2.11	0.48
1:C:65:VAL:CB	1:C:95[B]:CYS:SG	3.02	0.48
1:D:181:ALA:HB2	1:D:239:PRO:HB3	1.96	0.48
1:B:143:VAL:CG2	1:B:307:MET:HE2	2.44	0.48
1:D:65:VAL:CG2	1:D:95[B]:CYS:SG	2.97	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:159:ILE:HD13	1:D:294:GLN:HG2	1.96	0.48
1:A:275:ASN:HD22	1:A:293:LYS:CE	2.28	0.47
1:D:76:LYS:HE2	1:D:80:TYR:CE2	2.41	0.47
1:C:137:LYS:NZ	1:C:137:LYS:CB	2.72	0.47
1:D:178:ARG:HG2	1:D:178:ARG:NH1	2.28	0.47
1:D:282:PHE:C	1:D:282:PHE:CD2	2.90	0.47
1:A:101:TYR:HB2	1:A:127:ARG:NH1	2.29	0.47
1:A:283:ILE:CD1	1:B:223:GLN:HG3	2.46	0.46
1:B:127:ARG:HE	1:B:127:ARG:HB2	1.53	0.46
1:A:77:LEU:HD23	1:A:300:ALA:HB2	1.96	0.46
1:D:276:ILE:HG13	1:D:279:ILE:CD1	2.45	0.46
1:D:95[A]:CYS:SG	1:D:107:TYR:CD1	3.08	0.45
1:A:125:ASP:O	1:A:149:PHE:HZ	1.99	0.45
1:A:283:ILE:O	1:A:283:ILE:HG22	2.16	0.45
1:C:277:TYR:CD1	1:C:288:ARG:HD2	2.52	0.45
1:D:60:GLY:N	1:D:117:ASP:OD2	2.42	0.45
1:B:101:TYR:OH	1:B:131:LEU:HB2	2.16	0.45
1:B:71:ALA:HA	1:B:75:LEU:HD22	1.98	0.44
1:B:76:LYS:HD3	1:B:80:TYR:HE2	1.83	0.44
1:D:264:VAL:HG13	1:D:268:LEU:O	2.17	0.44
1:D:220:SER:O	1:D:220:SER:OG	2.33	0.44
1:D:85:ALA:CB	1:D:92:VAL:HG22	2.47	0.44
1:B:159:ILE:HD11	1:B:299:VAL:HG21	1.99	0.44
1:D:94:PHE:CD1	1:D:94:PHE:N	2.85	0.44
1:A:71:ALA:HA	1:A:75:LEU:HD22	2.00	0.43
1:A:283:ILE:HD12	1:B:223:GLN:HG3	2.00	0.43
1:A:280:PHE:HE1	1:B:280:PHE:CD2	2.36	0.43
1:C:179:LYS:NZ	1:C:237:GLU:OE1	2.37	0.43
1:A:275:ASN:HD22	1:A:293:LYS:HE3	1.83	0.43
1:B:147:ASN:O	1:B:156:ASP:OD2	2.36	0.43
1:D:292:MET:HE3	1:D:325:THR:OG1	2.15	0.43
1:D:136:LEU:CD2	1:D:154:VAL:HG12	2.25	0.43
1:C:85:ALA:CB	1:C:92:VAL:CG2	2.97	0.43
1:D:76:LYS:NZ	1:D:296:MET:HB2	2.32	0.43
1:A:75:LEU:HD12	1:A:75:LEU:HA	1.81	0.43
1:B:125:ASP:OD1	1:B:127:ARG:HG3	2.18	0.43
1:C:256:ARG:NH2	1:C:260:LEU:HD11	2.34	0.43
1:A:282:PHE:CD1	1:B:248:MET:HG3	2.52	0.42
1:B:75:LEU:HD12	1:B:75:LEU:HA	1.85	0.42
1:D:258:SER:HA	1:D:262:LEU:O	2.19	0.42
1:D:336:ARG:CG	1:D:336:ARG:NH1	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:LYS:CE	1:B:237:GLU:HG2	2.46	0.42
1:C:104:LYS:O	1:C:108:VAL:HG23	2.19	0.42
1:C:281:ARG:HA	1:C:281:ARG:HD2	1.49	0.42
1:A:276:ILE:HD12	1:A:276:ILE:HA	1.89	0.42
1:B:73:VAL:CG1	1:B:274:ASP:HB3	2.50	0.42
1:A:95:CYS:SG	1:A:107:TYR:CD1	3.07	0.42
1:A:181:ALA:HA	1:A:214:TYR:HB3	2.01	0.42
1:A:189:HIS:CD2	1:A:189:HIS:N	2.86	0.42
1:C:90:TYR:OH	1:C:311:ILE:HG21	2.19	0.41
1:D:136:LEU:CD1	1:D:154:VAL:HG13	2.37	0.41
1:C:220:SER:O	1:C:220:SER:OG	2.36	0.41
1:D:276:ILE:CG1	1:D:279:ILE:CD1	2.83	0.41
1:D:67:ASN:O	1:D:99:ASP:N	2.53	0.41
1:C:183:ILE:HD12	1:C:244:THR:HG22	2.02	0.41
1:D:75:LEU:HD12	1:D:75:LEU:HA	1.88	0.41
1:D:241:ALA:HA	1:D:269:SER:O	2.20	0.41
1:A:295:PRO:HA	2:A:478:HOH:O	2.20	0.41
1:B:276:ILE:HA	1:B:279:ILE:HD12	2.03	0.41
1:A:265:PRO:N	1:A:265:PRO:C	2.66	0.41
1:B:183:ILE:HG12	1:B:228:ALA:HB1	2.02	0.41
1:B:159:ILE:HD12	1:B:295:PRO:HD2	2.02	0.41
1:C:59:SER:HB3	1:C:60:GLY:H	1.64	0.41
1:C:125:ASP:O	1:C:149:PHE:HZ	2.04	0.41
1:C:254:ILE:HG12	1:C:264:VAL:HG11	2.03	0.41
1:C:265:PRO:HB3	1:C:332:SER:C	2.46	0.41
1:C:77:LEU:HD12	1:C:300:ALA:HB2	2.03	0.41
1:A:60:GLY:N	1:A:117:ASP:OD2	2.46	0.40
1:A:100:ASN:O	1:A:103:SER:HB2	2.21	0.40
1:D:66:LEU:CD2	1:D:122:PHE:CE2	3.04	0.40
1:D:295:PRO:O	1:D:299:VAL:HG23	2.21	0.40
1:D:104:LYS:NZ	1:D:128:ASP:OD2	2.46	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	278/343 (81%)	273 (98%)	4 (1%)	1 (0%)	30	34
1	B	276/343 (80%)	271 (98%)	4 (1%)	1 (0%)	30	34
1	C	278/343 (81%)	273 (98%)	4 (1%)	1 (0%)	30	34
1	D	279/343 (81%)	273 (98%)	5 (2%)	1 (0%)	30	34
All	All	1111/1372 (81%)	1090 (98%)	17 (2%)	4 (0%)	30	34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	274	ASP
1	B	274	ASP
1	C	274	ASP
1	D	274	ASP

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	239/296 (81%)	215 (90%)	24 (10%)	7	7
1	B	238/296 (80%)	218 (92%)	20 (8%)	10	11
1	C	240/296 (81%)	217 (90%)	23 (10%)	8	8
1	D	240/296 (81%)	219 (91%)	21 (9%)	9	10
All	All	957/1184 (81%)	869 (91%)	88 (9%)	8	9

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	SER
1	A	73	VAL
1	A	76	LYS
1	A	77	LEU

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Mol	Chain	Res	Type
1	A	97	SER
1	A	99	ASP
1	A	103	SER
1	A	105	SER
1	A	129	LYS
1	A	133	LYS
1	A	137	LYS
1	A	138	THR
1	A	151	ASP
1	A	190	ARG
1	A	218	THR
1	A	220	SER
1	A	222	GLU
1	A	230	ASP
1	A	258	SER
1	A	281	ARG
1	A	293	LYS
1	A	304	ILE
1	A	310	ASN
1	A	337	LYS
1	B	75	LEU
1	B	89	ASN
1	B	92	VAL
1	B	97	SER
1	B	98	ASN
1	B	99	ASP
1	B	100	ASN
1	B	102	GLU
1	B	104	LYS
1	B	124	SER
1	B	137	LYS
1	B	138	THR
1	B	145	ILE
1	B	151	ASP
1	B	218	THR
1	B	227	LYS
1	B	248	MET
1	B	276	ILE
1	B	294	GLN
1	B	310	ASN
1	C	59	SER
1	C	73	VAL

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Mol	Chain	Res	Type
1	C	76	LYS
1	C	89	ASN
1	C	102	GLU
1	C	105	SER
1	C	119	LEU
1	C	137	LYS
1	C	138	THR
1	C	145	ILE
1	C	154	VAL
1	C	220	SER
1	C	230	ASP
1	C	236	LYS
1	C	255	GLN
1	C	276	ILE
1	C	277	TYR
1	C	281	ARG
1	C	286	ASN
1	C	293	LYS
1	C	294	GLN
1	C	321	ILE
1	C	337	LYS
1	D	59	SER
1	D	76	LYS
1	D	92	VAL
1	D	98	ASN
1	D	133	LYS
1	D	138	THR
1	D	145	ILE
1	D	178	ARG
1	D	179	LYS
1	D	217	ASN
1	D	218	THR
1	D	222	GLU
1	D	260	LEU
1	D	262	LEU
1	D	263	SER
1	D	266	ARG
1	D	276	ILE
1	D	281	ARG
1	D	294	GLN
1	D	336	ARG
1	D	337	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	189	HIS
1	A	231	GLN
1	A	275	ASN
1	A	329	HIS
1	B	67	ASN
1	B	155	ASN
1	B	206	ASN
1	B	223	GLN
1	B	310	ASN
1	C	255	GLN
1	C	329	HIS
1	D	150	ASN
1	D	294	GLN

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	280/343 (81%)	0.21	16 (5%) 29 26	35, 53, 85, 113	0
1	B	278/343 (81%)	0.35	16 (5%) 29 25	37, 57, 94, 115	0
1	C	279/343 (81%)	0.28	14 (5%) 34 31	36, 56, 87, 109	1 (0%)
1	D	280/343 (81%)	0.49	20 (7%) 22 19	42, 61, 96, 121	1 (0%)
All	All	1117/1372 (81%)	0.33	66 (5%) 28 25	35, 56, 92, 121	2 (0%)

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	59	SER	5.6
1	B	337	LYS	4.9
1	C	283	ILE	4.7
1	C	337	LYS	4.6
1	A	248	MET	4.6
1	C	282	PHE	4.6
1	C	281	ARG	4.3
1	B	101	TYR	4.3
1	D	337	LYS	4.3
1	D	284	PRO	3.7
1	C	277	TYR	3.6
1	C	280	PHE	3.5
1	A	101	TYR	3.5
1	B	221	GLY	3.5
1	D	263	SER	3.5
1	A	281	ARG	3.4
1	A	283	ILE	3.4
1	D	248	MET	3.3
1	A	337	LYS	3.3
1	D	276	ILE	3.3
1	D	285	SER	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	314	ALA	3.2
1	B	248	MET	3.2
1	A	280	PHE	3.2
1	B	223	GLN	2.9
1	D	59	SER	2.9
1	A	58	LYS	2.7
1	B	222	GLU	2.7
1	C	286	ASN	2.6
1	B	137	LYS	2.6
1	D	286	ASN	2.6
1	B	318	PRO	2.6
1	D	58	LYS	2.6
1	D	219	ASP	2.6
1	A	282	PHE	2.5
1	C	285	SER	2.5
1	D	189	HIS	2.5
1	A	139	GLY	2.5
1	D	280	PHE	2.4
1	D	67	ASN	2.4
1	C	226	CYS	2.4
1	C	284	PRO	2.4
1	D	139	GLY	2.4
1	B	126	THR	2.3
1	C	139	GLY	2.3
1	D	279	ILE	2.3
1	B	86	TYR	2.3
1	A	95	CYS	2.3
1	B	127	ARG	2.3
1	B	334	CYS	2.3
1	D	281	ARG	2.3
1	A	286	ASN	2.2
1	D	223	GLN	2.2
1	A	190	ARG	2.2
1	D	95[A]	CYS	2.2
1	C	255	GLN	2.2
1	B	60	GLY	2.2
1	D	226	CYS	2.2
1	B	282	PHE	2.1
1	A	89	ASN	2.1
1	A	76	LYS	2.1
1	A	86	TYR	2.1
1	A	277	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	88	TYR	2.1
1	D	287	VAL	2.1
1	C	86	TYR	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.