



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 09:24 PM EDT

PDB ID : 3P14 / pdb_00003p14
Title : Crystal structure of L-rhamnose isomerase with a novel high thermo-stability from *Bacillus halodurans*
Authors : Doan, T.T.N.; Prabhu, P.; Kang, L.W.; Lee, J.K.
Deposited on : 2010-09-30
Resolution : 2.51 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

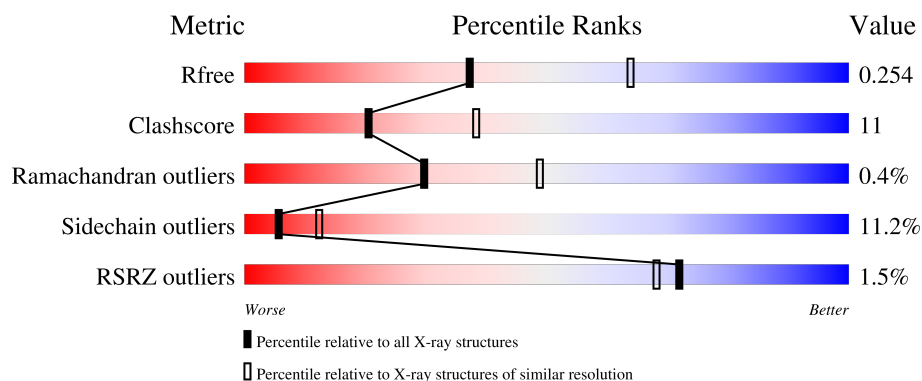
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	424	
1	B	424	
1	C	424	
1	D	424	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13611 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 L-rhamnose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	0	0	0
			3290	2109	566	606	9			
1	B	403	Total	C	N	O	S	0	0	0
			3298	2114	567	607	10			
1	C	402	Total	C	N	O	S	0	0	0
			3294	2112	566	606	10			
1	D	401	Total	C	N	O	S	0	0	0
			3278	2103	561	604	10			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP Q9KCL9
A	-4	HIS	-	expression tag	UNP Q9KCL9
A	-3	HIS	-	expression tag	UNP Q9KCL9
A	-2	HIS	-	expression tag	UNP Q9KCL9
A	-1	HIS	-	expression tag	UNP Q9KCL9
A	0	HIS	-	expression tag	UNP Q9KCL9
B	-5	HIS	-	expression tag	UNP Q9KCL9
B	-4	HIS	-	expression tag	UNP Q9KCL9
B	-3	HIS	-	expression tag	UNP Q9KCL9
B	-2	HIS	-	expression tag	UNP Q9KCL9
B	-1	HIS	-	expression tag	UNP Q9KCL9
B	0	HIS	-	expression tag	UNP Q9KCL9
C	-5	HIS	-	expression tag	UNP Q9KCL9
C	-4	HIS	-	expression tag	UNP Q9KCL9
C	-3	HIS	-	expression tag	UNP Q9KCL9
C	-2	HIS	-	expression tag	UNP Q9KCL9
C	-1	HIS	-	expression tag	UNP Q9KCL9
C	0	HIS	-	expression tag	UNP Q9KCL9
D	-5	HIS	-	expression tag	UNP Q9KCL9
D	-4	HIS	-	expression tag	UNP Q9KCL9
D	-3	HIS	-	expression tag	UNP Q9KCL9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	HIS	-	expression tag	UNP Q9KCL9
D	-1	HIS	-	expression tag	UNP Q9KCL9
D	0	HIS	-	expression tag	UNP Q9KCL9

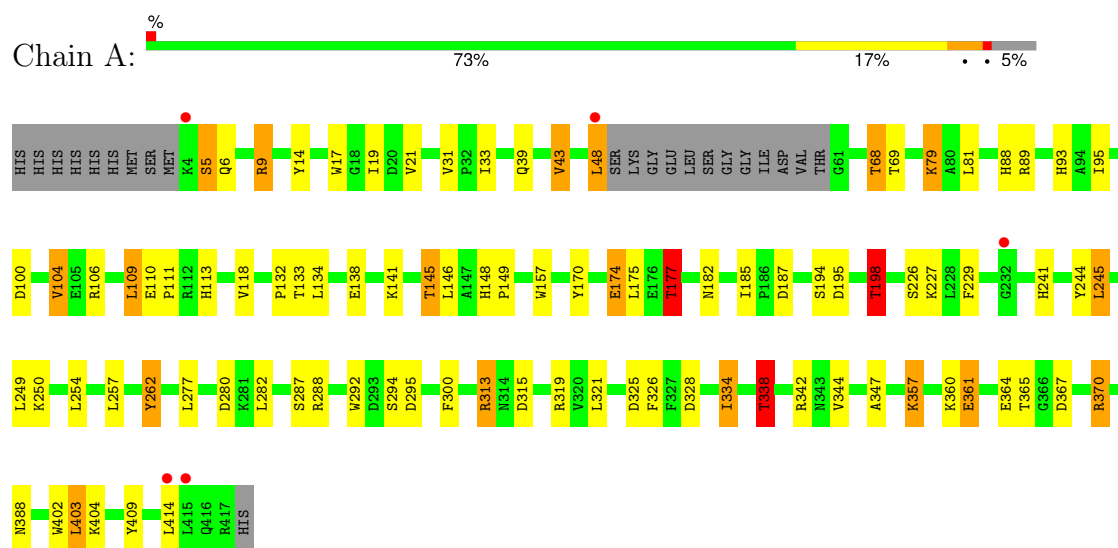
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	114	Total 114	O 114	0	0
2	B	114	Total 114	O 114	0	0
2	C	119	Total 119	O 119	0	0
2	D	104	Total 104	O 104	0	0

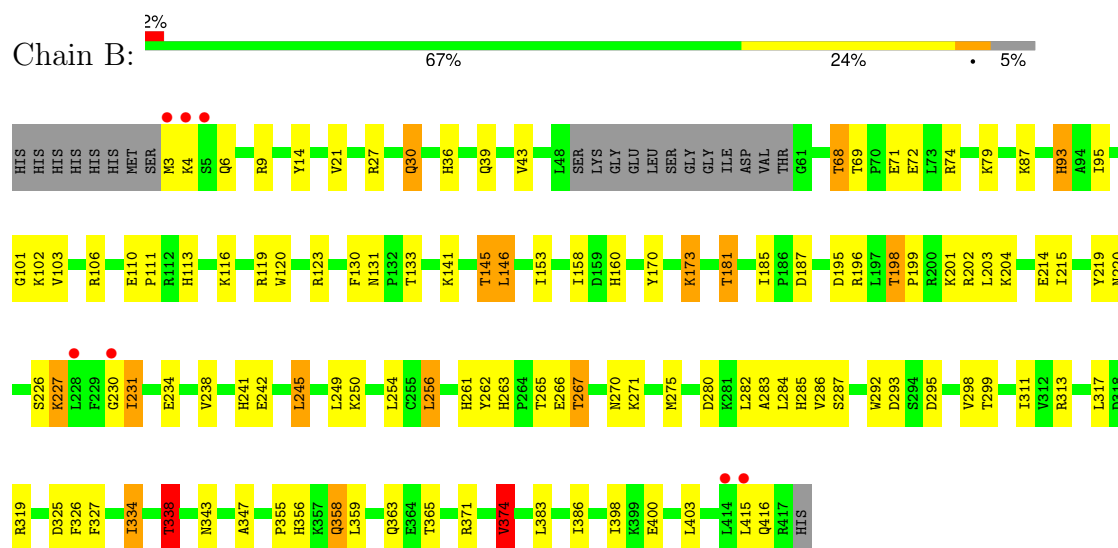
3 Residue-property plots [i](#)

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: L-rhamnose isomerase

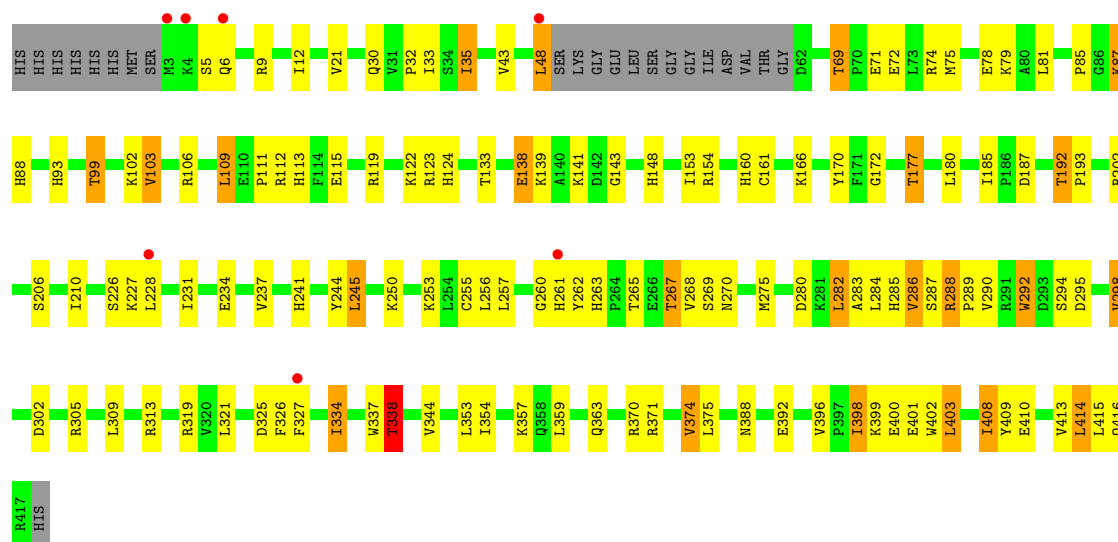


• Molecule 1: L-rhamnose isomerase

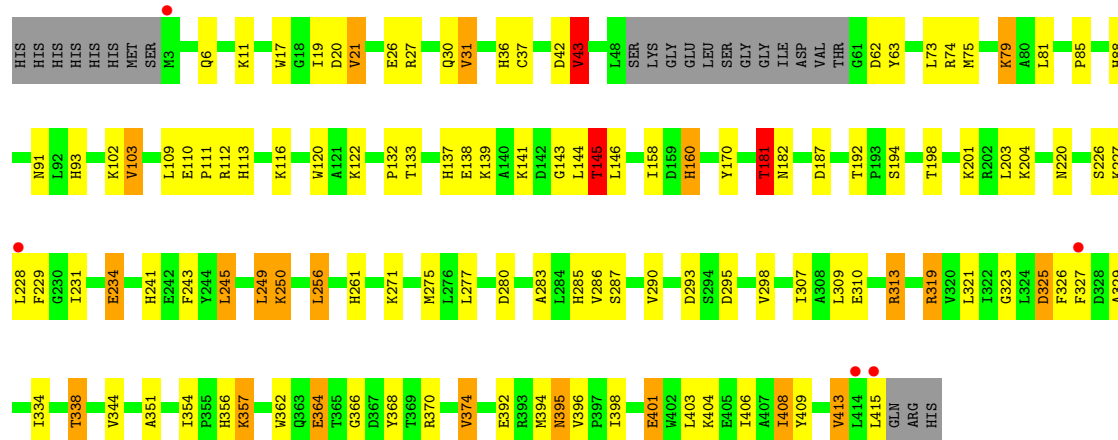


• Molecule 1: L-rhamnose isomerase





• Molecule 1: L-rhamnose isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.24Å 164.88Å 92.01Å 90.00° 115.97° 90.00°	Depositor
Resolution (Å)	45.78 – 2.51 45.78 – 2.51	Depositor EDS
% Data completeness (in resolution range)	97.7 (45.78-2.51) 98.0 (45.78-2.51)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.185 , 0.256 0.186 , 0.254	Depositor DCC
R_{free} test set	3751 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	34.4	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13611	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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 [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	1.08	3/3371 (0.1%)	1.14	7/4564 (0.2%)
1	B	1.06	3/3379 (0.1%)	1.16	12/4574 (0.3%)
1	C	1.08	5/3375 (0.1%)	1.16	6/4569 (0.1%)
1	D	1.08	9/3359 (0.3%)	1.21	13/4548 (0.3%)
All	All	1.08	20/13484 (0.1%)	1.17	38/18255 (0.2%)

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	408	ILE	CA-CB	7.78	1.64	1.54
1	A	338	THR	CA-CB	6.69	1.64	1.53
1	D	43	VAL	CA-CB	6.50	1.63	1.54
1	A	334	ILE	CA-CB	6.32	1.62	1.54
1	D	406	ILE	CA-CB	6.31	1.62	1.54
1	D	307	ILE	CA-CB	6.07	1.60	1.54
1	D	325	ASP	CA-C	-6.06	1.47	1.53
1	C	338	THR	CA-CB	5.97	1.62	1.53
1	D	31	VAL	CA-CB	5.91	1.59	1.53
1	D	181	THR	CA-CB	5.84	1.60	1.53
1	B	347	ALA	CA-CB	-5.78	1.44	1.53
1	D	374	VAL	CA-CB	5.62	1.62	1.54
1	D	192	THR	CA-CB	5.54	1.62	1.53
1	C	392	GLU	CG-CD	5.50	1.65	1.52
1	B	374	VAL	CA-CB	5.40	1.61	1.54
1	C	288	ARG	CA-C	5.25	1.59	1.52
1	A	198	THR	CA-CB	5.25	1.61	1.53
1	B	338	THR	CA-CB	5.20	1.61	1.53
1	C	298	VAL	CA-CB	5.10	1.59	1.54
1	D	103	VAL	CA-CB	5.08	1.59	1.53

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	30	GLN	CA-C-N	-10.28	115.19	122.59
1	D	30	GLN	C-N-CA	-10.28	115.19	122.59
1	B	334	ILE	N-CA-C	-7.66	102.65	111.00
1	D	88	HIS	N-CA-C	7.23	118.04	108.07
1	A	338	THR	CB-CA-C	7.01	122.77	110.85
1	D	245	LEU	N-CA-C	-6.88	103.86	111.36
1	A	5	SER	N-CA-C	-6.76	103.71	112.23
1	A	177	THR	CA-C-N	-6.76	113.02	119.85
1	A	177	THR	C-N-CA	-6.76	113.02	119.85
1	C	192	THR	CA-C-N	6.33	126.30	120.03
1	C	192	THR	C-N-CA	6.33	126.30	120.03
1	D	31	VAL	CB-CA-C	6.15	117.29	110.71
1	C	334	ILE	N-CA-C	-6.14	104.52	110.72
1	D	362	TRP	CA-CB-CG	6.02	125.04	113.60
1	B	30	GLN	CA-C-N	-6.01	118.26	122.59
1	B	30	GLN	C-N-CA	-6.01	118.26	122.59
1	B	356	HIS	N-CA-C	6.01	118.32	111.11
1	A	388	ASN	N-CA-C	5.78	117.58	111.28
1	D	228	LEU	N-CA-C	-5.73	104.64	111.69
1	D	198	THR	CA-C-N	-5.66	113.19	119.19
1	D	198	THR	C-N-CA	-5.66	113.19	119.19
1	B	198	THR	CA-C-N	-5.64	113.08	119.28
1	B	198	THR	C-N-CA	-5.64	113.08	119.28
1	A	17	TRP	N-CA-C	-5.59	106.50	113.38
1	B	262	TYR	N-CA-C	5.56	117.96	109.95
1	B	313	ARG	CB-CG-CD	-5.54	98.55	111.30
1	B	313	ARG	CB-CA-C	-5.45	99.88	110.46
1	D	160	HIS	N-CA-C	5.36	117.13	111.28
1	D	20	ASP	N-CA-C	5.34	117.69	107.75
1	C	388	ASN	N-CA-C	5.33	117.09	111.28
1	D	63	TYR	N-CA-C	5.18	115.56	109.60
1	C	177	THR	CB-CA-C	5.17	116.37	109.65
1	B	131	ASN	CA-C-N	-5.07	114.50	119.92
1	B	131	ASN	C-N-CA	-5.07	114.50	119.92
1	A	262	TYR	N-CA-C	5.05	117.53	109.65
1	C	292	TRP	N-CA-C	-5.02	101.78	108.86
1	B	158	ILE	N-CA-C	-5.01	105.61	110.42
1	D	158	ILE	N-CA-C	-5.00	105.83	110.53

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	3290	0	3244	68	0
1	B	3298	0	3253	74	0
1	C	3294	0	3250	87	0
1	D	3278	0	3232	72	0
2	A	114	0	0	2	0
2	B	114	0	0	1	0
2	C	119	0	0	2	0
2	D	104	0	0	2	0
All	All	13611	0	12979	293	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 (293) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:TYR:HB3	1:A:19:ILE:CD1	1.62	1.27
1:A:14:TYR:CB	1:A:19:ILE:HD11	1.64	1.26
1:B:145:THR:HG22	1:B:146:LEU:H	1.11	1.14
1:A:145:THR:HG22	1:A:146:LEU:H	1.13	1.08
1:C:99:THR:HG21	1:C:102:LYS:O	1.51	1.08
1:C:99:THR:CG2	1:C:102:LYS:O	2.03	1.07
1:C:286:VAL:HG23	1:C:298:VAL:HG21	1.29	1.07
1:A:357:LYS:HD3	1:A:357:LYS:N	1.68	1.05
1:D:357:LYS:H	1:D:357:LYS:CD	1.69	1.05
1:D:357:LYS:H	1:D:357:LYS:HD3	0.91	1.05
1:A:357:LYS:HD3	1:A:357:LYS:H	0.87	1.03
1:B:227:LYS:HG3	1:B:261:HIS:CE1	1.92	1.02
1:D:357:LYS:HD3	1:D:357:LYS:N	1.77	0.99
1:A:89:ARG:HH12	1:A:177:THR:HG21	1.25	0.98
1:B:110:GLU:H	1:B:113:HIS:HD2	1.07	0.98
1:A:357:LYS:H	1:A:357:LYS:CD	1.76	0.98
1:B:145:THR:CG2	1:B:146:LEU:H	1.79	0.95
1:B:133:THR:H	1:B:160:HIS:HE1	1.13	0.95
1:C:69:THR:HG22	1:C:72:GLU:H	1.31	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:286:VAL:HG23	1:D:298:VAL:HG21	1.50	0.94
1:D:110:GLU:H	1:D:113:HIS:HD2	1.18	0.90
1:B:286:VAL:HG23	1:B:298:VAL:HG21	1.54	0.89
1:B:287:SER:HB3	1:B:325:ASP:O	1.73	0.89
1:D:133:THR:H	1:D:160:HIS:HE1	1.20	0.87
1:A:287:SER:HB3	1:A:325:ASP:O	1.75	0.86
1:A:14:TYR:HB3	1:A:19:ILE:HD11	0.87	0.85
1:A:145:THR:HG22	1:A:146:LEU:N	1.87	0.85
1:B:145:THR:HG22	1:B:146:LEU:N	1.93	0.83
1:A:14:TYR:HB3	1:A:19:ILE:CG1	2.09	0.82
1:D:356:HIS:HD2	2:D:492:HOH:O	1.61	0.82
1:A:110:GLU:H	1:A:113:HIS:HD2	1.24	0.82
1:D:145:THR:CG2	1:D:146:LEU:H	1.93	0.81
1:B:271:LYS:O	1:B:275:MET:HG2	1.81	0.81
1:A:364:GLU:OE2	1:C:202:ARG:NH2	2.14	0.79
1:A:195:ASP:OD2	1:A:198:THR:HG23	1.82	0.79
1:C:286:VAL:HG23	1:C:298:VAL:CG2	2.09	0.78
1:A:145:THR:CG2	1:A:146:LEU:H	1.94	0.78
1:A:145:THR:HB	1:A:187:ASP:OD1	1.84	0.77
1:B:363:GLN:HE21	1:B:371:ARG:HH11	1.32	0.77
1:A:280:ASP:O	1:A:319:ARG:HD2	1.85	0.77
1:C:288:ARG:HD2	2:C:455:HOH:O	1.83	0.76
1:C:228:LEU:HD13	1:C:263:HIS:CE1	2.21	0.76
1:B:363:GLN:NE2	1:D:194:SER:H	1.83	0.76
1:A:110:GLU:H	1:A:113:HIS:CD2	2.04	0.75
1:B:110:GLU:H	1:B:113:HIS:CD2	1.99	0.75
1:B:68:THR:HG22	1:B:69:THR:HG23	1.69	0.75
1:B:363:GLN:HE22	1:D:194:SER:H	1.31	0.75
1:D:413:VAL:HG12	1:D:413:VAL:O	1.87	0.75
1:A:145:THR:CG2	1:A:146:LEU:N	2.47	0.74
1:B:280:ASP:O	1:B:319:ARG:HD2	1.88	0.74
1:C:267:THR:HG22	1:C:270:ASN:H	1.53	0.73
1:C:309:LEU:O	1:C:313:ARG:HB2	1.89	0.73
1:C:257:LEU:HD22	1:C:262:TYR:OH	1.90	0.72
1:A:68:THR:HG22	1:A:69:THR:HG23	1.72	0.71
1:C:286:VAL:CG2	1:C:298:VAL:HG21	2.16	0.71
1:D:145:THR:HG23	1:D:146:LEU:H	1.55	0.71
1:D:271:LYS:O	1:D:275:MET:HG2	1.91	0.70
1:B:145:THR:HB	1:B:187:ASP:OD1	1.90	0.70
1:B:119:ARG:HH12	1:B:123:ARG:HH21	1.40	0.69
1:D:334:ILE:O	1:D:338:THR:HG23	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:295:ASP:HB3	1:C:326:PHE:HA	1.74	0.69
1:B:68:THR:HB	1:B:72:GLU:OE1	1.92	0.69
1:B:145:THR:CG2	1:B:146:LEU:N	2.44	0.69
1:D:287:SER:HB3	1:D:325:ASP:O	1.93	0.68
1:B:110:GLU:N	1:B:113:HIS:HD2	1.87	0.68
1:C:35:ILE:HD13	1:C:337:TRP:CD1	2.28	0.68
1:C:115:GLU:HG2	1:C:119:ARG:NH2	2.09	0.67
1:B:119:ARG:NH1	1:B:123:ARG:HH21	1.92	0.67
1:B:133:THR:N	1:B:160:HIS:HE1	1.89	0.66
1:B:256:LEU:HD22	1:B:283:ALA:CB	2.25	0.66
1:A:5:SER:O	1:A:9:ARG:HG2	1.95	0.66
1:B:181:THR:CG2	1:B:220:ASN:OD1	2.42	0.66
1:A:89:ARG:HH12	1:A:177:THR:CG2	2.07	0.65
1:A:48:LEU:C	1:A:48:LEU:HD22	2.21	0.65
1:D:181:THR:CG2	1:D:220:ASN:OD1	2.45	0.65
1:D:256:LEU:HD21	1:D:285:HIS:CE1	2.32	0.65
1:A:19:ILE:HD12	1:A:19:ILE:O	1.97	0.65
1:D:181:THR:HG22	1:D:220:ASN:OD1	1.97	0.65
1:C:287:SER:HB3	1:C:325:ASP:O	1.97	0.65
1:C:106:ARG:HA	1:C:109:LEU:HD22	1.80	0.64
1:D:145:THR:CG2	1:D:146:LEU:N	2.60	0.64
1:A:104:VAL:CG2	1:A:109:LEU:HD13	2.27	0.63
1:B:93:HIS:HD2	1:B:133:THR:OG1	1.81	0.63
1:C:99:THR:HG23	1:C:102:LYS:O	1.96	0.62
1:C:282:LEU:HD22	1:C:283:ALA:N	2.15	0.61
1:A:295:ASP:HB3	1:A:326:PHE:HA	1.81	0.61
1:B:287:SER:CB	1:B:325:ASP:O	2.46	0.61
1:C:35:ILE:CD1	1:C:337:TRP:CD1	2.83	0.61
1:B:295:ASP:HB3	1:B:326:PHE:HA	1.83	0.61
1:A:134:LEU:HG	1:A:185:ILE:HG22	1.82	0.60
1:B:146:LEU:HD22	1:B:185:ILE:HG21	1.82	0.60
1:B:231:ILE:HD13	1:B:292:TRP:CE3	2.37	0.60
1:A:106:ARG:HA	1:A:109:LEU:HD22	1.84	0.59
1:C:75:MET:HE3	1:C:409:TYR:HE1	1.67	0.59
1:B:119:ARG:HH12	1:B:123:ARG:NH2	2.01	0.59
1:D:398:ILE:O	1:D:401:GLU:HB2	2.03	0.59
1:A:357:LYS:N	1:A:357:LYS:CD	2.50	0.59
1:B:241:HIS:O	1:B:245:LEU:HB2	2.02	0.59
1:D:133:THR:N	1:D:160:HIS:HE1	1.98	0.59
1:B:202:ARG:NH2	1:D:364:GLU:OE1	2.35	0.59
1:B:256:LEU:HD22	1:B:283:ALA:HB1	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:THR:HG23	1:D:146:LEU:N	2.18	0.58
1:C:148:HIS:O	1:C:154:ARG:HD3	2.03	0.58
1:D:110:GLU:H	1:D:113:HIS:CD2	2.09	0.58
1:D:413:VAL:O	1:D:413:VAL:CG1	2.50	0.58
1:C:133:THR:H	1:C:160:HIS:HE1	1.50	0.58
1:C:398:ILE:HG13	1:C:399:LYS:HG3	1.85	0.58
1:C:267:THR:CG2	1:C:270:ASN:H	2.17	0.58
1:B:133:THR:H	1:B:160:HIS:CE1	2.06	0.57
1:B:181:THR:HG23	1:B:220:ASN:OD1	2.03	0.57
1:A:19:ILE:HD12	1:A:19:ILE:C	2.29	0.57
1:C:48:LEU:HD22	1:C:103:VAL:HG12	1.87	0.57
1:B:227:LYS:CG	1:B:261:HIS:CE1	2.81	0.57
1:C:206:SER:O	1:C:210:ILE:HG13	2.05	0.57
1:D:137:HIS:HE1	1:D:139:LYS:HD3	1.70	0.57
1:B:263:HIS:HB2	1:B:266:GLU:OE1	2.04	0.56
1:C:180:LEU:HD22	1:C:321:LEU:HD23	1.87	0.56
1:D:357:LYS:CD	1:D:357:LYS:N	2.49	0.56
1:A:334:ILE:O	1:A:338:THR:HG23	2.06	0.56
1:D:133:THR:H	1:D:160:HIS:CE1	2.12	0.56
1:D:256:LEU:HB2	1:D:283:ALA:HB3	1.86	0.56
1:C:227:LYS:HD2	1:C:261:HIS:CE1	2.41	0.56
1:D:227:LYS:HE3	1:D:229:PHE:O	2.05	0.55
1:C:35:ILE:CD1	1:C:337:TRP:NE1	2.69	0.55
1:D:111:PRO:HB3	1:D:170:TYR:CD2	2.41	0.55
1:B:334:ILE:O	1:B:338:THR:HG23	2.06	0.55
1:C:226:SER:O	1:C:261:HIS:ND1	2.29	0.55
1:D:17:TRP:HB2	1:D:19:ILE:HD12	1.89	0.54
1:A:194:SER:H	1:C:363:GLN:NE2	2.05	0.54
1:B:256:LEU:HD22	1:B:283:ALA:HB3	1.89	0.54
1:C:231:ILE:CD1	1:C:292:TRP:CE3	2.90	0.54
1:D:404:LYS:O	1:D:408:ILE:HG23	2.08	0.54
1:A:288:ARG:HD2	2:A:462:HOH:O	2.08	0.54
1:C:398:ILE:O	1:C:401:GLU:HG2	2.08	0.54
1:C:115:GLU:HG2	1:C:119:ARG:HH22	1.73	0.53
1:A:241:HIS:CE1	1:A:245:LEU:HD12	2.43	0.53
1:C:282:LEU:HD22	1:C:283:ALA:H	1.74	0.53
1:C:228:LEU:CD1	1:C:263:HIS:CE1	2.90	0.53
1:C:280:ASP:O	1:C:319:ARG:HD2	2.09	0.53
1:B:363:GLN:HE22	1:D:194:SER:N	2.05	0.52
1:C:363:GLN:HE21	1:C:371:ARG:HH11	1.56	0.52
1:D:143:GLY:O	1:D:187:ASP:HA	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:ARG:NH2	1:C:113:HIS:HE1	2.08	0.52
1:C:402:TRP:CZ3	1:C:403:LEU:HD13	2.45	0.52
1:C:309:LEU:HD21	1:C:354:ILE:HG13	1.92	0.52
1:C:231:ILE:HD13	1:C:292:TRP:CZ3	2.45	0.51
1:D:295:ASP:HB3	1:D:326:PHE:HA	1.93	0.51
1:B:106:ARG:HB3	1:B:160:HIS:CD2	2.45	0.51
1:D:203:LEU:HD23	1:D:243:PHE:CE2	2.46	0.51
1:B:111:PRO:HB3	1:B:170:TYR:CD2	2.46	0.50
1:A:227:LYS:HE2	1:A:229:PHE:O	2.11	0.50
1:D:290:VAL:HG23	1:D:290:VAL:O	2.09	0.50
1:D:145:THR:HG22	1:D:146:LEU:H	1.70	0.50
1:A:39:GLN:HE22	1:A:328:ASP:H	1.59	0.50
1:B:130:PHE:O	1:B:181:THR:HA	2.11	0.50
1:D:42:ASP:C	1:D:43:VAL:HG23	2.37	0.50
1:D:250:LYS:HD2	1:D:250:LYS:O	2.12	0.50
1:A:334:ILE:O	1:A:338:THR:CG2	2.59	0.50
1:D:231:ILE:HA	1:D:234:GLU:OE2	2.12	0.50
1:B:359:LEU:HD21	1:B:374:VAL:HG22	1.93	0.49
1:A:257:LEU:CD2	1:A:262:TYR:OH	2.60	0.49
1:A:95:ILE:HG23	1:A:133:THR:HG21	1.93	0.49
1:A:313:ARG:HD3	2:C:481:HOH:O	2.12	0.49
1:C:255:CYS:HB2	1:C:275:MET:HE1	1.94	0.49
1:D:11:LYS:HG3	1:D:21:VAL:HG22	1.94	0.49
1:A:367:ASP:OD2	1:A:370:ARG:HG3	2.12	0.49
1:C:85:PRO:HD3	1:C:396:VAL:HG21	1.94	0.49
1:C:257:LEU:CD2	1:C:262:TYR:OH	2.59	0.49
1:D:79:LYS:HE3	1:D:79:LYS:HA	1.95	0.49
1:C:231:ILE:HD13	1:C:292:TRP:CE3	2.47	0.49
1:A:404:LYS:HA	1:A:404:LYS:HE3	1.95	0.49
1:C:226:SER:HB3	1:C:244:TYR:HD2	1.77	0.49
1:D:137:HIS:CE1	1:D:139:LYS:HD3	2.48	0.49
1:D:309:LEU:O	1:D:313:ARG:HB3	2.12	0.48
1:B:267:THR:CG2	1:B:270:ASN:H	2.26	0.48
1:D:37:CYS:HB2	1:D:73:LEU:HD21	1.94	0.48
1:D:226:SER:O	1:D:261:HIS:ND1	2.37	0.48
1:A:361:GLU:O	1:A:365:THR:OG1	2.27	0.48
1:C:334:ILE:O	1:C:338:THR:HG23	2.14	0.48
1:C:410:GLU:HA	1:C:414:LEU:HB2	1.95	0.48
1:C:359:LEU:HD21	1:C:374:VAL:HG22	1.95	0.48
1:A:338:THR:O	1:A:342:ARG:HB2	2.14	0.47
1:C:112:ARG:HH21	1:C:113:HIS:HE1	1.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:TYR:O	1:A:19:ILE:HG13	2.14	0.47
1:A:134:LEU:HD13	1:A:157:TRP:CE3	2.49	0.47
1:B:242:GLU:OE2	1:D:241:HIS:HE1	1.97	0.47
1:C:231:ILE:HD11	1:C:292:TRP:CE3	2.49	0.47
1:B:355:PRO:HB2	1:B:358:GLN:HG3	1.96	0.47
1:C:172:GLY:HA2	1:C:177:THR:O	2.15	0.47
1:D:145:THR:HG22	1:D:187:ASP:OD1	2.14	0.47
1:C:256:LEU:HD21	1:C:285:HIS:CG	2.49	0.47
1:D:249:LEU:HD23	1:D:249:LEU:HA	1.71	0.47
1:A:79:LYS:HG3	1:A:409:TYR:CD1	2.49	0.47
1:B:95:ILE:HG23	1:B:133:THR:HG21	1.95	0.47
1:D:394:MET:O	1:D:395:ASN:HB2	2.15	0.47
1:A:367:ASP:CG	1:A:370:ARG:HG2	2.40	0.47
1:B:196:ARG:HD2	1:D:310:GLU:OE2	2.14	0.47
1:A:257:LEU:HD23	1:A:262:TYR:OH	2.14	0.47
1:B:74:ARG:HD2	1:B:120:TRP:CG	2.50	0.47
1:A:6:GLN:HA	1:A:6:GLN:OE1	2.15	0.46
1:C:111:PRO:HB3	1:C:170:TYR:CG	2.50	0.46
1:D:62:ASP:HB2	1:D:329:ALA:O	2.15	0.46
1:D:227:LYS:HD2	1:D:261:HIS:CE1	2.50	0.46
1:C:112:ARG:NH2	1:C:113:HIS:CE1	2.83	0.46
1:D:351:ALA:O	1:D:356:HIS:HE1	1.99	0.46
1:A:33:ILE:O	1:A:88:HIS:HB3	2.15	0.46
1:A:402:TRP:CZ3	1:A:403:LEU:HD13	2.51	0.46
1:B:241:HIS:CE1	1:B:245:LEU:HD12	2.51	0.46
1:D:79:LYS:HD2	1:D:409:TYR:CD1	2.51	0.46
1:A:367:ASP:OD1	1:A:370:ARG:HG2	2.16	0.46
1:C:282:LEU:HD22	1:C:282:LEU:C	2.40	0.46
1:D:74:ARG:HD2	1:D:120:TRP:CD1	2.51	0.46
1:C:357:LYS:HA	1:C:357:LYS:HD3	1.79	0.45
1:A:148:HIS:CG	1:A:149:PRO:HD2	2.52	0.45
1:B:36:HIS:NE2	1:B:39:GLN:HB2	2.31	0.45
1:C:79:LYS:HD2	1:C:79:LYS:HA	1.75	0.45
1:B:173:LYS:HG2	1:B:219:TYR:CE1	2.51	0.45
1:B:198:THR:HB	1:B:199:PRO:HD3	1.98	0.45
1:B:226:SER:O	1:B:261:HIS:ND1	2.37	0.45
1:C:87:LYS:HE3	1:C:124:HIS:O	2.16	0.45
1:C:185:ILE:HD12	1:C:187:ASP:CG	2.42	0.45
1:A:104:VAL:HG23	1:A:109:LEU:HD13	1.97	0.44
1:A:404:LYS:HA	1:A:404:LYS:CE	2.47	0.44
1:A:132:PRO:HD2	1:A:182:ASN:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35:ILE:HD12	1:C:337:TRP:NE1	2.32	0.44
1:C:267:THR:HG23	1:C:269:SER:OG	2.17	0.44
1:D:287:SER:CB	1:D:325:ASP:O	2.64	0.44
1:B:256:LEU:HD11	1:B:285:HIS:CG	2.53	0.44
1:B:130:PHE:HB3	1:B:181:THR:HB	2.00	0.44
1:B:231:ILE:CD1	1:B:292:TRP:CE3	3.01	0.44
1:C:143:GLY:O	1:C:187:ASP:HA	2.18	0.44
1:C:375:LEU:HD23	1:C:375:LEU:HA	1.90	0.44
1:B:3:MET:O	1:B:6:GLN:HB2	2.17	0.44
1:A:292:TRP:NE1	1:A:294:SER:HB3	2.33	0.44
1:A:118:VAL:HG11	1:A:174:GLU:HG2	2.00	0.43
1:B:203:LEU:HD22	1:B:238:VAL:HG12	2.00	0.43
1:D:280:ASP:O	1:D:319:ARG:HD2	2.18	0.43
1:D:366:GLY:HA2	1:D:368:TYR:CE1	2.54	0.43
1:A:226:SER:HB3	1:A:244:TYR:HD2	1.82	0.43
1:B:227:LYS:HE3	1:B:293:ASP:OD2	2.18	0.43
1:D:201:LYS:HE2	2:D:487:HOH:O	2.17	0.43
1:C:241:HIS:O	1:C:245:LEU:HB2	2.18	0.43
1:A:300:PHE:HE1	1:A:347:ALA:HA	1.84	0.43
1:C:33:ILE:O	1:C:88:HIS:HB3	2.18	0.43
1:C:153:ILE:O	1:C:154:ARG:C	2.61	0.43
1:C:302:ASP:OD1	1:C:305:ARG:NH1	2.52	0.42
1:C:192:THR:HA	1:C:193:PRO:HD2	1.87	0.42
1:C:415:LEU:HD23	1:C:415:LEU:HA	1.71	0.42
1:C:75:MET:HE2	1:C:413:VAL:HG11	2.01	0.42
1:D:138:GLU:OE2	1:D:141:LYS:HE3	2.19	0.42
1:D:145:THR:CG2	1:D:187:ASP:OD1	2.67	0.42
1:A:367:ASP:CG	1:A:370:ARG:CG	2.92	0.42
1:C:9:ARG:O	1:C:12:ILE:HG12	2.19	0.42
1:C:30:GLN:O	1:C:32:PRO:HD3	2.19	0.42
1:D:285:HIS:HA	1:D:323:GLY:O	2.20	0.42
1:D:309:LEU:HD21	1:D:354:ILE:HG13	2.02	0.42
1:A:48:LEU:C	1:A:48:LEU:CD2	2.89	0.42
1:C:133:THR:H	1:C:160:HIS:CE1	2.35	0.42
1:A:367:ASP:OD2	1:A:370:ARG:CG	2.68	0.42
1:C:282:LEU:HD23	1:C:282:LEU:HA	1.67	0.42
1:D:286:VAL:HG23	1:D:298:VAL:CG2	2.36	0.42
1:A:79:LYS:HE2	1:A:409:TYR:HB2	2.01	0.42
1:A:111:PRO:HB3	1:A:170:TYR:CG	2.55	0.42
1:B:79:LYS:HD2	1:B:79:LYS:HA	1.96	0.41
1:B:267:THR:HG22	1:B:270:ASN:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:122:LYS:O	1:C:123:ARG:C	2.61	0.41
1:C:260:GLY:HA2	1:C:289:PRO:HG3	2.02	0.41
1:C:302:ASP:HA	1:C:305:ARG:NH1	2.35	0.41
1:B:400:GLU:CD	1:B:400:GLU:H	2.27	0.41
1:C:161:CYS:CB	1:C:210:ILE:HD13	2.50	0.41
1:B:27:ARG:HH11	1:B:30:GLN:HB2	1.85	0.41
1:C:370:ARG:O	1:C:374:VAL:HG13	2.20	0.41
1:B:231:ILE:HD13	1:B:292:TRP:CZ3	2.55	0.41
1:D:85:PRO:HD3	1:D:396:VAL:HG21	2.03	0.41
1:D:234:GLU:H	1:D:234:GLU:HG2	1.58	0.41
1:B:214:GLU:H	1:B:214:GLU:HG2	1.62	0.41
1:D:132:PRO:HD2	1:D:182:ASN:O	2.20	0.41
1:C:74:ARG:O	1:C:78:GLU:HG3	2.20	0.41
1:C:109:LEU:HD12	1:C:113:HIS:CD2	2.56	0.41
1:B:230:GLY:HA3	2:B:528:HOH:O	2.20	0.41
1:B:185:ILE:HD12	1:B:187:ASP:CG	2.45	0.41
1:B:195:ASP:OD2	1:B:198:THR:OG1	2.34	0.41
1:B:299:THR:HA	1:B:343:ASN:ND2	2.35	0.41
1:C:99:THR:HG22	1:C:102:LYS:H	1.86	0.41
1:C:166:LYS:HD3	1:C:166:LYS:HA	1.93	0.41
1:D:36:HIS:HA	1:D:91:ASN:HB3	2.02	0.41
1:B:311:ILE:HG22	1:B:317:LEU:HD23	2.03	0.41
1:B:14:TYR:CE2	1:B:386:ILE:HD13	2.56	0.40
1:C:141:LYS:HA	1:C:141:LYS:HD2	1.84	0.40
1:A:138:GLU:O	1:A:141:LYS:HG2	2.22	0.40
1:B:226:SER:C	1:B:261:HIS:HD1	2.27	0.40
1:D:408:ILE:HG12	1:D:409:TYR:N	2.37	0.40
1:A:9:ARG:NH1	2:A:522:HOH:O	2.53	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	398/424 (94%)	382 (96%)	15 (4%)	1 (0%)	36	55
1	B	399/424 (94%)	379 (95%)	19 (5%)	1 (0%)	36	55
1	C	398/424 (94%)	380 (96%)	17 (4%)	1 (0%)	36	55
1	D	397/424 (94%)	373 (94%)	20 (5%)	4 (1%)	12	24
All	All	1592/1696 (94%)	1514 (95%)	71 (4%)	7 (0%)	30	49

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	293	ASP
1	D	112	ARG
1	D	43	VAL
1	D	145	THR
1	C	138	GLU
1	B	101	GLY
1	A	43	VAL

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	351/370 (95%)	317 (90%)	34 (10%)	8	17
1	B	352/370 (95%)	310 (88%)	42 (12%)	5	11
1	C	352/370 (95%)	312 (89%)	40 (11%)	5	12
1	D	350/370 (95%)	308 (88%)	42 (12%)	5	10
All	All	1405/1480 (95%)	1247 (89%)	158 (11%)	6	12

All (158) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	21	VAL
1	A	31	VAL
1	A	43	VAL

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Mol	Chain	Res	Type
1	A	48	LEU
1	A	68	THR
1	A	79	LYS
1	A	81	LEU
1	A	93	HIS
1	A	100	ASP
1	A	104	VAL
1	A	109	LEU
1	A	145	THR
1	A	174	GLU
1	A	175	LEU
1	A	177	THR
1	A	198	THR
1	A	245	LEU
1	A	249	LEU
1	A	250	LYS
1	A	254	LEU
1	A	277	LEU
1	A	282	LEU
1	A	313	ARG
1	A	315	ASP
1	A	321	LEU
1	A	338	THR
1	A	344	VAL
1	A	357	LYS
1	A	360	LYS
1	A	361	GLU
1	A	370	ARG
1	A	403	LEU
1	A	414	LEU
1	B	4	LYS
1	B	9	ARG
1	B	21	VAL
1	B	43	VAL
1	B	68	THR
1	B	71	GLU
1	B	87	LYS
1	B	93	HIS
1	B	102	LYS
1	B	103	VAL
1	B	116	LYS
1	B	141	LYS

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Mol	Chain	Res	Type
1	B	145	THR
1	B	146	LEU
1	B	153	ILE
1	B	173	LYS
1	B	181	THR
1	B	201	LYS
1	B	204	LYS
1	B	215	ILE
1	B	227	LYS
1	B	231	ILE
1	B	234	GLU
1	B	245	LEU
1	B	249	LEU
1	B	250	LYS
1	B	254	LEU
1	B	256	LEU
1	B	265	THR
1	B	267	THR
1	B	282	LEU
1	B	284	LEU
1	B	327	PHE
1	B	338	THR
1	B	358	GLN
1	B	365	THR
1	B	374	VAL
1	B	383	LEU
1	B	398	ILE
1	B	403	LEU
1	B	415	LEU
1	B	416	GLN
1	C	5	SER
1	C	6	GLN
1	C	21	VAL
1	C	35	ILE
1	C	43	VAL
1	C	48	LEU
1	C	69	THR
1	C	71	GLU
1	C	81	LEU
1	C	87	LYS
1	C	93	HIS
1	C	99	THR

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Mol	Chain	Res	Type
1	C	103	VAL
1	C	109	LEU
1	C	138	GLU
1	C	139	LYS
1	C	234	GLU
1	C	237	VAL
1	C	245	LEU
1	C	250	LYS
1	C	253	LYS
1	C	265	THR
1	C	267	THR
1	C	268	VAL
1	C	282	LEU
1	C	284	LEU
1	C	286	VAL
1	C	290	VAL
1	C	294	SER
1	C	327	PHE
1	C	338	THR
1	C	344	VAL
1	C	353	LEU
1	C	374	VAL
1	C	398	ILE
1	C	400	GLU
1	C	403	LEU
1	C	408	ILE
1	C	414	LEU
1	C	416	GLN
1	D	6	GLN
1	D	21	VAL
1	D	26	GLU
1	D	27	ARG
1	D	31	VAL
1	D	43	VAL
1	D	75	MET
1	D	79	LYS
1	D	81	LEU
1	D	93	HIS
1	D	102	LYS
1	D	103	VAL
1	D	109	LEU
1	D	116	LYS

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Mol	Chain	Res	Type
1	D	122	LYS
1	D	144	LEU
1	D	145	THR
1	D	181	THR
1	D	204	LYS
1	D	234	GLU
1	D	245	LEU
1	D	249	LEU
1	D	250	LYS
1	D	256	LEU
1	D	277	LEU
1	D	313	ARG
1	D	319	ARG
1	D	321	LEU
1	D	327	PHE
1	D	338	THR
1	D	344	VAL
1	D	357	LYS
1	D	364	GLU
1	D	370	ARG
1	D	374	VAL
1	D	392	GLU
1	D	395	ASN
1	D	401	GLU
1	D	403	LEU
1	D	408	ILE
1	D	413	VAL
1	D	415	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	39	GLN
1	A	113	HIS
1	A	131	ASN
1	A	182	ASN
1	A	209	GLN
1	A	220	ASN
1	A	241	HIS
1	A	388	ASN
1	B	39	GLN

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Mol	Chain	Res	Type
1	B	91	ASN
1	B	93	HIS
1	B	113	HIS
1	B	131	ASN
1	B	155	GLN
1	B	160	HIS
1	B	251	ASN
1	B	358	GLN
1	B	363	GLN
1	B	388	ASN
1	C	16	GLN
1	C	39	GLN
1	C	113	HIS
1	C	131	ASN
1	C	160	HIS
1	C	251	ASN
1	C	314	ASN
1	C	363	GLN
1	C	416	GLN
1	D	16	GLN
1	D	91	ASN
1	D	113	HIS
1	D	160	HIS
1	D	182	ASN
1	D	209	GLN
1	D	314	ASN
1	D	356	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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

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	402/424 (94%)	-0.37	5 (1%) 76 73	17, 30, 45, 66	0
1	B	403/424 (95%)	-0.36	7 (1%) 69 65	17, 30, 50, 85	0
1	C	402/424 (94%)	-0.38	7 (1%) 69 65	19, 30, 45, 81	0
1	D	401/424 (94%)	-0.26	5 (1%) 76 73	18, 31, 50, 71	0
All	All	1608/1696 (94%)	-0.34	24 (1%) 72 68	17, 30, 48, 85	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	415	LEU	4.8
1	C	228	LEU	4.1
1	D	414	LEU	3.6
1	B	228	LEU	3.5
1	B	3	MET	3.4
1	C	3	MET	3.1
1	D	3	MET	3.0
1	C	48	LEU	2.8
1	A	414	LEU	2.8
1	B	414	LEU	2.6
1	B	4	LYS	2.4
1	C	261	HIS	2.4
1	B	415	LEU	2.4
1	B	5	SER	2.4
1	D	327	PHE	2.3
1	A	48	LEU	2.3
1	A	232	GLY	2.2
1	A	4	LYS	2.2
1	A	415	LEU	2.1
1	B	230	GLY	2.1
1	D	228	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	327	PHE	2.1
1	C	6	GLN	2.1
1	C	4	LYS	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.