



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:27 AM UTC

PDB ID : 2HTB / pdb\_00002htb  
Title : Crystal Structure of a putative mutarotase (YeaD) from *Salmonella typhimurium* in monoclinic form  
Authors : Chittori, S.; Simanshu, D.K.; Savithri, H.S.; Murthy, M.R.N.  
Deposited on : 2006-07-25  
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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

---

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

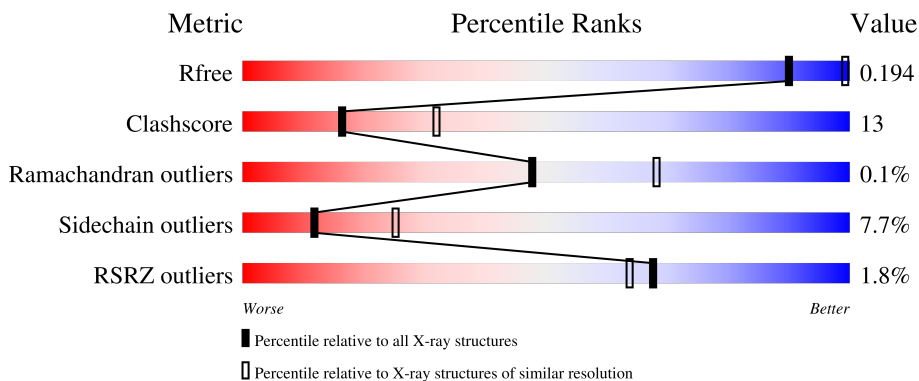
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	309	
1	B	309	
1	C	309	
1	D	309	

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 9434 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 Putative enzyme related to aldose 1-epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	296	2286	1458	395	424	9	0	0	0
1	B	296	2256	1438	387	422	9	0	0	0
1	C	295	2287	1457	397	424	9	0	0	0
1	D	296	2283	1455	396	423	9	0	0	0

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MET	-	cloning artifact	UNP Q8ZPV9
A	-13	ARG	-	cloning artifact	UNP Q8ZPV9
A	-12	GLY	-	cloning artifact	UNP Q8ZPV9
A	-11	SER	-	cloning artifact	UNP Q8ZPV9
A	-10	HIS	-	expression tag	UNP Q8ZPV9
A	-9	HIS	-	expression tag	UNP Q8ZPV9
A	-8	HIS	-	expression tag	UNP Q8ZPV9
A	-7	HIS	-	expression tag	UNP Q8ZPV9
A	-6	HIS	-	expression tag	UNP Q8ZPV9
A	-5	HIS	-	expression tag	UNP Q8ZPV9
A	-4	GLY	-	cloning artifact	UNP Q8ZPV9
A	-3	MET	-	cloning artifact	UNP Q8ZPV9
A	-2	ALA	-	cloning artifact	UNP Q8ZPV9
A	-1	SER	-	cloning artifact	UNP Q8ZPV9
A	0	HIS	-	cloning artifact	UNP Q8ZPV9
B	-14	MET	-	cloning artifact	UNP Q8ZPV9
B	-13	ARG	-	cloning artifact	UNP Q8ZPV9
B	-12	GLY	-	cloning artifact	UNP Q8ZPV9
B	-11	SER	-	cloning artifact	UNP Q8ZPV9
B	-10	HIS	-	expression tag	UNP Q8ZPV9
B	-9	HIS	-	expression tag	UNP Q8ZPV9

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	HIS	-	expression tag	UNP Q8ZPV9
B	-7	HIS	-	expression tag	UNP Q8ZPV9
B	-6	HIS	-	expression tag	UNP Q8ZPV9
B	-5	HIS	-	expression tag	UNP Q8ZPV9
B	-4	GLY	-	cloning artifact	UNP Q8ZPV9
B	-3	MET	-	cloning artifact	UNP Q8ZPV9
B	-2	ALA	-	cloning artifact	UNP Q8ZPV9
B	-1	SER	-	cloning artifact	UNP Q8ZPV9
B	0	HIS	-	cloning artifact	UNP Q8ZPV9
C	-14	MET	-	cloning artifact	UNP Q8ZPV9
C	-13	ARG	-	cloning artifact	UNP Q8ZPV9
C	-12	GLY	-	cloning artifact	UNP Q8ZPV9
C	-11	SER	-	cloning artifact	UNP Q8ZPV9
C	-10	HIS	-	expression tag	UNP Q8ZPV9
C	-9	HIS	-	expression tag	UNP Q8ZPV9
C	-8	HIS	-	expression tag	UNP Q8ZPV9
C	-7	HIS	-	expression tag	UNP Q8ZPV9
C	-6	HIS	-	expression tag	UNP Q8ZPV9
C	-5	HIS	-	expression tag	UNP Q8ZPV9
C	-4	GLY	-	cloning artifact	UNP Q8ZPV9
C	-3	MET	-	cloning artifact	UNP Q8ZPV9
C	-2	ALA	-	cloning artifact	UNP Q8ZPV9
C	-1	SER	-	cloning artifact	UNP Q8ZPV9
C	0	HIS	-	cloning artifact	UNP Q8ZPV9
D	-14	MET	-	cloning artifact	UNP Q8ZPV9
D	-13	ARG	-	cloning artifact	UNP Q8ZPV9
D	-12	GLY	-	cloning artifact	UNP Q8ZPV9
D	-11	SER	-	cloning artifact	UNP Q8ZPV9
D	-10	HIS	-	expression tag	UNP Q8ZPV9
D	-9	HIS	-	expression tag	UNP Q8ZPV9
D	-8	HIS	-	expression tag	UNP Q8ZPV9
D	-7	HIS	-	expression tag	UNP Q8ZPV9
D	-6	HIS	-	expression tag	UNP Q8ZPV9
D	-5	HIS	-	expression tag	UNP Q8ZPV9
D	-4	GLY	-	cloning artifact	UNP Q8ZPV9
D	-3	MET	-	cloning artifact	UNP Q8ZPV9
D	-2	ALA	-	cloning artifact	UNP Q8ZPV9
D	-1	SER	-	cloning artifact	UNP Q8ZPV9
D	0	HIS	-	cloning artifact	UNP Q8ZPV9

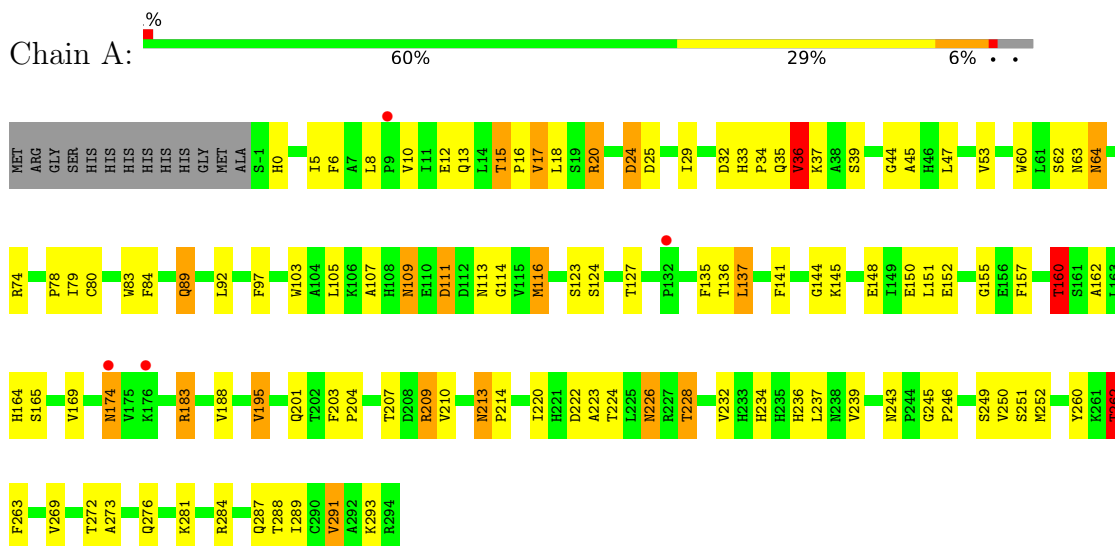
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	98	Total O 98 98	0	0
2	B	45	Total O 45 45	0	0
2	C	75	Total O 75 75	0	0
2	D	104	Total O 104 104	0	0

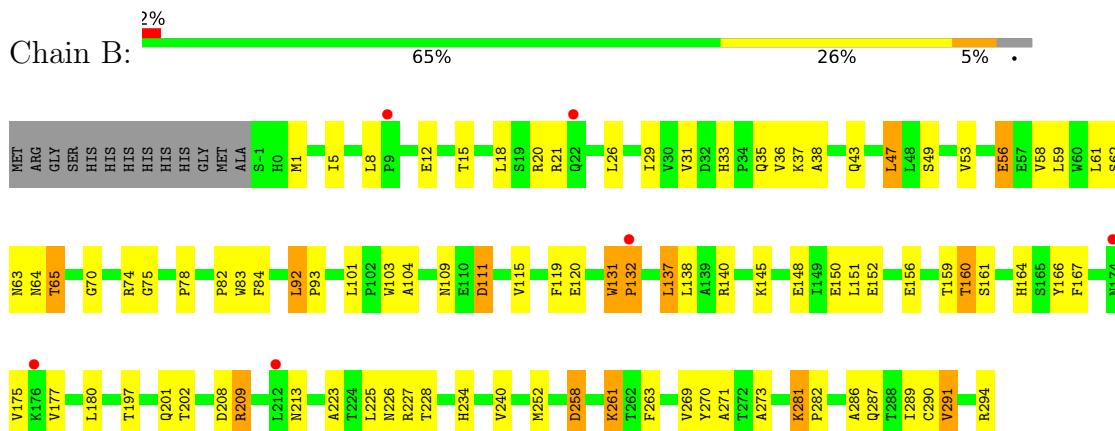
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

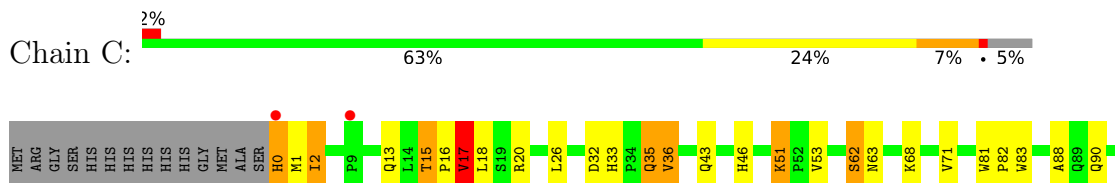
- Molecule 1: Putative enzyme related to aldose 1-epimerase

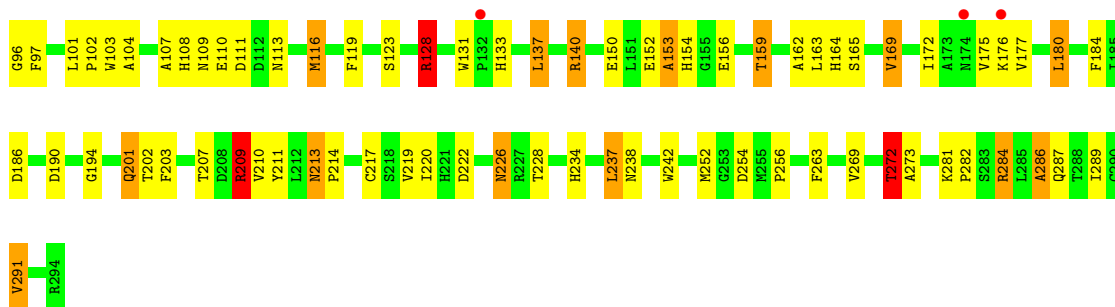


- Molecule 1: Putative enzyme related to aldose 1-epimerase

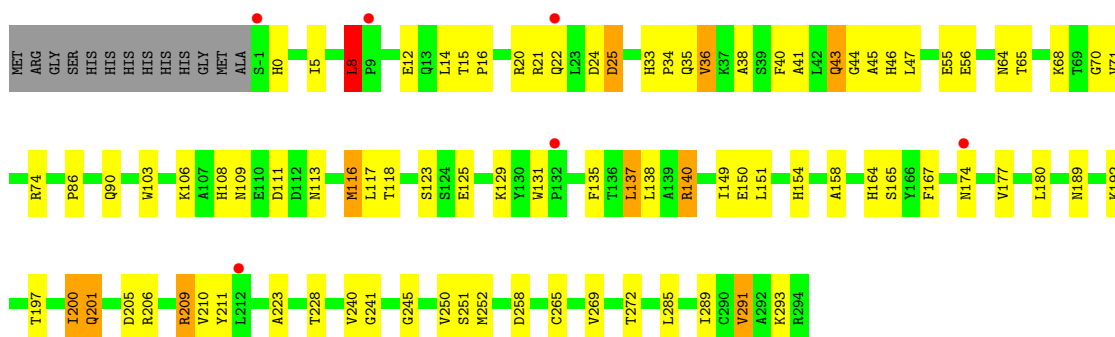


- Molecule 1: Putative enzyme related to aldose 1-epimerase





- Molecule 1: Putative enzyme related to aldose 1-epimerase



## 4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	219.55Å 47.60Å 138.09Å 90.00° 91.82° 90.00°	Depositor
Resolution (Å)	27.80 – 2.50 27.80 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.1 (27.80-2.50) 99.1 (27.80-2.50)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.45 (at 2.50Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.189 , 0.240 0.186 , 0.194	Depositor DCC
$R_{free}$ test set	2526 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.5	Xtrriage
Anisotropy	0.025	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 35.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.001 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9434	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	1.64	19/2347 (0.8%)	1.47	25/3204 (0.8%)
1	B	1.50	10/2317 (0.4%)	1.43	18/3172 (0.6%)
1	C	1.61	17/2348 (0.7%)	1.47	31/3206 (1.0%)
1	D	1.60	17/2345 (0.7%)	1.45	18/3206 (0.6%)
All	All	1.59	63/9357 (0.7%)	1.45	92/12788 (0.7%)

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	C	0	1

All (63) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	195	VAL	CA-CB	7.76	1.64	1.54
1	D	117	LEU	CA-C	7.68	1.62	1.52
1	C	210	VAL	CA-CB	7.53	1.62	1.53
1	C	53	VAL	CA-CB	7.50	1.63	1.54
1	C	153	ALA	CA-CB	-7.35	1.43	1.53
1	D	131	TRP	C-O	-7.10	1.17	1.24
1	A	136	THR	C-O	7.01	1.31	1.24
1	A	10	VAL	CA-CB	6.81	1.62	1.54
1	B	132	PRO	CA-C	-6.58	1.47	1.53
1	C	140	ARG	CA-C	-6.47	1.45	1.52
1	D	241	GLY	C-O	6.44	1.31	1.23
1	C	119	PHE	CA-C	-6.42	1.45	1.52
1	A	53	VAL	CA-CB	6.34	1.60	1.54
1	D	138	LEU	CA-C	-6.33	1.45	1.52
1	B	111	ASP	C-O	6.26	1.30	1.24
1	A	107	ALA	CA-CB	6.24	1.64	1.53

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	273	ALA	CA-C	6.13	1.58	1.52
1	C	2	ILE	CA-CB	6.10	1.61	1.54
1	A	79	ILE	C-O	6.07	1.30	1.24
1	C	20	ARG	CA-CB	-6.05	1.44	1.53
1	D	22	GLN	N-CA	-6.04	1.38	1.46
1	B	223	ALA	CA-CB	5.85	1.63	1.53
1	A	262	THR	CA-CB	5.80	1.62	1.53
1	B	63	ASN	C-O	-5.75	1.16	1.24
1	B	36	VAL	CA-CB	5.68	1.61	1.55
1	D	109	ASN	N-CA	-5.68	1.39	1.46
1	B	290	CYS	N-CA	5.66	1.52	1.45
1	C	286	ALA	C-O	5.61	1.30	1.23
1	A	210	VAL	CA-CB	5.59	1.60	1.53
1	C	46	HIS	N-CA	-5.58	1.39	1.46
1	A	18	LEU	C-O	5.57	1.30	1.24
1	A	20	ARG	N-CA	-5.49	1.39	1.46
1	A	47	LEU	N-CA	-5.49	1.39	1.46
1	A	80	CYS	C-O	5.41	1.30	1.24
1	D	118	THR	N-CA	-5.40	1.39	1.46
1	C	88	ALA	CA-CB	5.37	1.62	1.53
1	C	104	ALA	CA-CB	5.37	1.61	1.53
1	D	86	PRO	C-O	-5.35	1.17	1.23
1	C	71	VAL	CA-CB	5.31	1.60	1.54
1	C	128	ARG	CA-C	5.31	1.59	1.52
1	C	273	ALA	CA-C	5.25	1.57	1.52
1	D	269	VAL	C-O	5.23	1.29	1.23
1	D	43	GLN	C-O	5.21	1.30	1.24
1	C	140	ARG	N-CA	5.21	1.52	1.46
1	B	209	ARG	N-CA	-5.21	1.39	1.46
1	A	116	MET	CB-CG	5.16	1.68	1.52
1	A	281	LYS	CA-C	5.16	1.57	1.52
1	A	273	ALA	CA-CB	5.15	1.60	1.53
1	A	251	SER	N-CA	5.14	1.52	1.46
1	B	37	LYS	C-O	5.13	1.30	1.24
1	C	119	PHE	C-O	-5.12	1.17	1.23
1	D	252	MET	C-O	5.11	1.30	1.24
1	A	39	SER	C-O	5.09	1.30	1.23
1	B	202	THR	CA-CB	5.09	1.61	1.53
1	D	149	ILE	CA-CB	5.09	1.60	1.54
1	D	116	MET	CA-C	5.09	1.58	1.52
1	C	217	CYS	N-CA	-5.08	1.40	1.46
1	D	38	ALA	CA-CB	5.07	1.62	1.53

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	228	THR	C-O	5.07	1.30	1.23
1	D	41	ALA	CA-CB	-5.06	1.45	1.53
1	A	111	ASP	CB-CG	-5.04	1.39	1.52
1	D	158	ALA	CA-CB	-5.02	1.45	1.53
1	D	71	VAL	C-O	-5.01	1.18	1.24

All (92) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	36	VAL	CB-CA-C	-9.68	96.90	111.80
1	A	145	LYS	N-CA-C	-8.97	102.11	113.23
1	C	272	THR	N-CA-C	8.79	121.81	111.71
1	C	213	ASN	N-CA-C	8.49	123.89	108.94
1	B	15	THR	CB-CA-C	-8.48	100.41	110.76
1	B	101	LEU	CA-C-N	-8.44	111.03	119.90
1	B	101	LEU	C-N-CA	-8.44	111.03	119.90
1	C	17	VAL	CB-CA-C	8.42	118.91	110.65
1	B	65	THR	CA-C-N	-8.34	112.27	120.52
1	B	65	THR	C-N-CA	-8.34	112.27	120.52
1	A	195	VAL	CB-CA-C	8.03	119.56	110.41
1	C	36	VAL	CB-CA-C	-7.92	96.72	110.65
1	B	281	LYS	CA-C-N	7.79	127.80	119.78
1	B	281	LYS	C-N-CA	7.79	127.80	119.78
1	D	245	GLY	CA-C-N	7.79	127.07	118.97
1	D	245	GLY	C-N-CA	7.79	127.07	118.97
1	C	35	GLN	N-CA-C	7.75	122.50	112.89
1	C	15	THR	CA-C-N	-7.64	112.33	119.82
1	C	15	THR	C-N-CA	-7.64	112.33	119.82
1	C	226	ASN	N-CA-C	7.64	121.60	111.28
1	B	161	SER	N-CA-C	7.47	120.10	108.96
1	C	291	VAL	CB-CA-C	-7.42	99.28	110.55
1	D	113	ASN	N-CA-C	7.42	120.81	111.69
1	B	29	ILE	N-CA-C	-7.41	97.74	108.11
1	A	291	VAL	CB-CA-C	-7.36	98.92	110.69
1	D	36	VAL	CB-CA-C	-7.19	98.00	111.30
1	D	8	LEU	CA-C-N	-7.12	113.47	120.52
1	D	8	LEU	C-N-CA	-7.12	113.47	120.52
1	A	97	PHE	N-CA-C	-7.11	103.83	114.64
1	A	8	LEU	CA-C-N	-7.09	113.01	120.03
1	A	8	LEU	C-N-CA	-7.09	113.01	120.03
1	B	26	LEU	N-CA-C	6.77	120.28	109.24
1	C	51	LYS	CA-C-N	6.61	126.58	119.78

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	51	LYS	C-N-CA	6.61	126.58	119.78
1	C	101	LEU	CA-C-N	-6.54	112.93	119.99
1	C	101	LEU	C-N-CA	-6.54	112.93	119.99
1	B	175	VAL	N-CA-C	6.33	117.04	108.17
1	D	64	ASN	CB-CA-C	-6.31	98.95	109.55
1	A	0	HIS	N-CA-C	-6.30	104.49	111.36
1	C	62	SER	N-CA-C	6.26	119.28	109.96
1	B	197	THR	N-CA-C	-6.22	104.28	113.61
1	B	64	ASN	CB-CA-C	-6.19	99.75	110.09
1	C	203	PHE	CA-C-N	-6.17	113.33	119.56
1	C	203	PHE	C-N-CA	-6.17	113.33	119.56
1	A	15	THR	N-CA-C	-6.08	102.66	108.07
1	C	254	ASP	N-CA-C	6.08	120.31	113.02
1	D	125	GLU	N-CA-C	-6.03	104.70	111.28
1	B	145	LYS	N-CA-C	-6.02	105.43	112.89
1	D	201	GLN	CB-CA-C	-6.00	98.59	109.38
1	D	291	VAL	CB-CA-C	-5.99	101.60	110.82
1	C	209	ARG	NE-CZ-NH1	5.97	127.47	121.50
1	C	165	SER	N-CA-C	5.92	119.19	110.30
1	C	163	LEU	N-CA-C	-5.92	97.66	108.02
1	A	226	ASN	N-CA-C	5.91	119.26	111.28
1	A	92	LEU	CA-C-N	-5.86	114.72	120.52
1	A	92	LEU	C-N-CA	-5.86	114.72	120.52
1	D	36	VAL	N-CA-C	5.82	116.03	108.35
1	B	131	TRP	CA-C-N	5.72	129.02	120.96
1	B	131	TRP	C-N-CA	5.72	129.02	120.96
1	D	25	ASP	CA-C-N	-5.71	115.12	123.00
1	D	25	ASP	C-N-CA	-5.71	115.12	123.00
1	C	209	ARG	NE-CZ-NH2	-5.69	114.08	119.20
1	C	62	SER	CB-CA-C	-5.69	101.96	110.16
1	D	258	ASP	N-CA-C	5.69	120.21	113.16
1	D	241	GLY	CA-C-N	-5.66	112.64	122.37
1	D	241	GLY	C-N-CA	-5.66	112.64	122.37
1	A	174	ASN	CB-CA-C	-5.57	102.43	111.06
1	C	140	ARG	NE-CZ-NH2	-5.56	114.20	119.20
1	A	210	VAL	CB-CA-C	-5.55	104.31	111.53
1	A	114	GLY	N-CA-C	5.54	118.55	110.63
1	D	200	ILE	N-CA-C	5.54	115.83	107.75
1	A	165	SER	N-CA-C	5.43	118.19	110.10
1	A	5	ILE	N-CA-C	5.41	115.54	110.30
1	A	74	ARG	NE-CZ-NH2	-5.37	114.37	119.20
1	B	92	LEU	N-CA-C	5.37	118.07	110.24

*Continued on next page...*

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	213	ASN	CA-C-N	-5.34	115.07	120.31
1	A	213	ASN	C-N-CA	-5.34	115.07	120.31
1	A	281	LYS	N-CA-C	5.28	120.14	108.27
1	C	284	ARG	N-CA-C	5.26	116.98	109.14
1	C	213	ASN	CA-C-N	5.24	125.69	120.14
1	C	213	ASN	C-N-CA	5.24	125.69	120.14
1	C	140	ARG	CG-CD-NE	-5.23	100.49	112.00
1	C	180	LEU	N-CA-C	5.22	119.52	113.20
1	B	150	GLU	N-CA-C	5.20	116.87	108.34
1	C	17	VAL	N-CA-CB	-5.20	105.09	112.12
1	A	105	LEU	N-CA-C	-5.18	98.88	107.99
1	C	26	LEU	N-CA-C	5.18	117.68	109.24
1	A	111	ASP	CB-CA-C	-5.13	99.35	111.65
1	A	263	PHE	N-CA-C	5.11	116.83	109.07
1	A	160	THR	N-CA-C	-5.08	100.43	109.06
1	C	133	HIS	N-CA-C	5.07	117.98	110.48
1	D	209	ARG	N-CA-C	5.06	117.64	108.69

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	0	HIS	Peptide

## 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	2286	0	2209	74	0
1	B	2256	0	2150	61	0
1	C	2287	0	2221	66	0
1	D	2283	0	2205	43	0
2	A	98	0	0	3	0
2	B	45	0	0	1	0
2	C	75	0	0	7	0
2	D	104	0	0	1	0
All	All	9434	0	8785	233	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 13.

All (233) 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:C:201:GLN:HG3	1:C:211:TYR:OH	1.54	1.06
1:B:84:PHE:H	1:B:160:THR:HG23	1.26	0.98
1:A:183:ARG:HH11	1:A:183:ARG:HG2	1.29	0.96
1:A:127:THR:HG22	2:A:299:HOH:O	1.67	0.93
1:D:140:ARG:HG2	1:D:140:ARG:HH11	1.36	0.88
1:B:84:PHE:H	1:B:160:THR:CG2	1.87	0.87
1:A:89:GLN:HE21	1:C:107:ALA:HA	1.40	0.85
1:C:111:ASP:HB3	1:C:113:ASN:H	1.42	0.85
1:B:201:GLN:HE22	1:B:209:ARG:HH11	1.22	0.84
1:D:140:ARG:HH11	1:D:140:ARG:CG	1.89	0.84
1:A:89:GLN:NE2	1:C:107:ALA:HA	1.95	0.81
1:B:33:HIS:HD2	1:B:35:GLN:H	1.28	0.80
1:A:111:ASP:HB3	1:A:113:ASN:H	1.47	0.80
1:D:33:HIS:HD2	1:D:35:GLN:H	1.29	0.79
1:A:36:VAL:HG22	1:A:144:GLY:HA2	1.66	0.78
1:A:89:GLN:NE2	1:C:108:HIS:H	1.82	0.77
1:B:33:HIS:CD2	1:B:35:GLN:H	2.02	0.77
1:D:33:HIS:CD2	1:D:35:GLN:H	2.04	0.76
1:B:84:PHE:N	1:B:160:THR:HG23	2.00	0.75
1:B:83:TRP:HA	1:B:160:THR:HG22	1.69	0.75
1:A:89:GLN:HE21	1:C:107:ALA:CA	1.99	0.74
1:A:36:VAL:CG2	1:A:144:GLY:HA2	2.18	0.73
1:B:18:LEU:HD21	1:B:115:VAL:HG11	1.69	0.73
1:A:60:TRP:CD1	1:A:262:THR:HG22	2.24	0.72
1:A:84:PHE:H	1:A:160:THR:CG2	2.03	0.71
1:A:174:ASN:ND2	1:A:224:THR:OG1	2.23	0.71
1:A:84:PHE:H	1:A:160:THR:HG22	1.54	0.71
1:A:183:ARG:HG2	1:A:183:ARG:NH1	2.06	0.71
1:D:164:HIS:O	1:D:164:HIS:CD2	2.44	0.70
1:A:103:TRP:CH2	1:A:137:LEU:HD13	2.27	0.69
1:A:209:ARG:HD2	2:A:321:HOH:O	1.93	0.69
1:C:103:TRP:CH2	1:C:137:LEU:HD13	2.28	0.69
1:C:164:HIS:O	1:C:164:HIS:CD2	2.46	0.69
1:A:60:TRP:NE1	1:A:262:THR:HG22	2.09	0.68
1:D:103:TRP:CH2	1:D:137:LEU:HD13	2.27	0.68
1:D:164:HIS:O	1:D:164:HIS:HD2	1.76	0.68
1:A:83:TRP:HA	1:A:160:THR:HG22	1.75	0.68

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:GLN:HE22	1:C:108:HIS:H	1.42	0.68
1:B:201:GLN:HE22	1:B:209:ARG:NH1	1.92	0.67
1:B:137:LEU:HD22	1:B:151:LEU:HD22	1.77	0.66
1:A:33:HIS:HD2	1:A:35:GLN:H	1.44	0.66
1:D:108:HIS:HA	1:D:116:MET:O	1.97	0.65
1:C:207:THR:O	1:C:242:TRP:HA	1.97	0.65
1:C:164:HIS:O	1:C:164:HIS:HD2	1.78	0.64
1:D:140:ARG:HG2	1:D:140:ARG:NH1	2.12	0.64
1:A:33:HIS:CD2	1:A:35:GLN:H	2.16	0.64
1:B:33:HIS:HE1	1:B:111:ASP:O	1.81	0.64
1:C:109:ASN:HD22	1:C:110:GLU:H	1.44	0.64
1:B:140:ARG:HH21	1:B:152:GLU:CD	2.07	0.63
1:C:1:MET:H	1:C:63:ASN:ND2	1.97	0.62
1:C:238:ASN:HD21	1:C:272:THR:HB	1.65	0.62
1:D:201:GLN:CG	1:D:211:TYR:OH	2.48	0.61
1:D:201:GLN:OE1	1:D:209:ARG:NH1	2.32	0.61
1:C:159:THR:HG21	2:C:309:HOH:O	1.99	0.61
1:D:5:ILE:HA	1:D:8:LEU:HD22	1.81	0.61
1:C:33:HIS:HD2	1:C:35:GLN:H	1.49	0.60
1:D:14:LEU:HD12	1:D:14:LEU:N	2.17	0.60
1:C:33:HIS:CD2	1:C:35:GLN:H	2.20	0.59
1:D:140:ARG:HD2	1:D:150:GLU:HB2	1.84	0.59
1:D:123:SER:HB3	1:D:135:PHE:CD2	2.36	0.59
1:A:124:SER:H	1:A:127:THR:HB	1.67	0.59
1:C:109:ASN:ND2	1:C:110:GLU:H	2.00	0.59
1:C:17:VAL:HG22	1:C:32:ASP:O	2.03	0.59
1:B:83:TRP:CA	1:B:160:THR:HG22	2.32	0.58
1:A:84:PHE:N	1:A:160:THR:HG22	2.18	0.58
1:C:33:HIS:HE1	1:C:111:ASP:O	1.86	0.58
1:D:33:HIS:HE1	1:D:111:ASP:O	1.85	0.58
1:C:201:GLN:NE2	1:C:202:THR:O	2.36	0.58
1:C:234:HIS:HB2	1:C:286:ALA:HB3	1.86	0.57
1:D:201:GLN:HG3	1:D:211:TYR:OH	2.04	0.57
1:D:103:TRP:HH2	1:D:137:LEU:HD13	1.68	0.57
1:B:1:MET:HB3	1:B:61:LEU:HD23	1.86	0.57
1:B:5:ILE:O	1:B:8:LEU:HB2	2.03	0.57
1:B:92:LEU:HB3	1:B:93:PRO:HD2	1.86	0.57
1:B:47:LEU:HD13	1:B:59:LEU:HD11	1.87	0.57
1:B:47:LEU:HD13	1:B:59:LEU:CD1	2.35	0.57
1:A:12:GLU:OE2	1:B:20:ARG:HD2	2.05	0.56
1:A:89:GLN:NE2	1:C:108:HIS:N	2.54	0.55

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:169:VAL:HG11	1:C:175:VAL:HG21	1.89	0.55
1:C:116:MET:HB2	2:C:364:HOH:O	2.06	0.55
1:D:55:GLU:HB3	2:D:373:HOH:O	2.05	0.55
1:C:209:ARG:HD2	2:C:302:HOH:O	2.06	0.55
1:A:60:TRP:HE1	1:A:262:THR:HG22	1.71	0.55
1:A:276:GLN:HE22	1:C:154:HIS:HE1	1.54	0.54
1:B:103:TRP:CH2	1:B:137:LEU:HD13	2.42	0.54
1:C:109:ASN:HD22	1:C:110:GLU:N	2.04	0.54
1:A:174:ASN:HB3	1:A:223:ALA:HB3	1.89	0.54
1:A:6:PHE:CE1	1:A:37:LYS:HB3	2.43	0.53
1:D:209:ARG:O	1:D:240:VAL:HA	2.08	0.53
1:B:226:ASN:HD21	1:B:294:ARG:CB	2.22	0.53
1:C:222:ASP:C	1:C:222:ASP:OD1	2.51	0.52
1:B:201:GLN:NE2	1:B:209:ARG:HH11	2.01	0.52
1:B:226:ASN:O	1:B:294:ARG:N	2.36	0.52
1:C:1:MET:H	1:C:63:ASN:HD22	1.55	0.52
1:A:207:THR:CG2	1:A:209:ARG:HD3	2.38	0.52
1:A:246:PRO:O	1:A:250:VAL:HG23	2.09	0.52
1:D:44:GLY:O	1:D:45:ALA:C	2.51	0.52
1:B:82:PRO:HB2	1:B:159:THR:HG22	1.91	0.52
1:B:137:LEU:CD2	1:B:151:LEU:HD22	2.40	0.52
1:B:234:HIS:HB2	1:B:286:ALA:HB3	1.92	0.52
1:A:152:GLU:OE2	1:A:284:ARG:NH1	2.44	0.51
1:D:33:HIS:HD2	1:D:35:GLN:N	2.03	0.51
1:D:33:HIS:CD2	1:D:34:PRO:HD2	2.45	0.51
1:A:116:MET:HA	1:A:141:PHE:O	2.10	0.51
1:D:135:PHE:HA	1:D:154:HIS:O	2.11	0.51
1:B:74:ARG:NH2	1:B:252:MET:HE3	2.26	0.51
1:C:43:GLN:HA	1:C:102:PRO:HA	1.93	0.51
1:A:78:PRO:HD2	1:A:164:HIS:O	2.12	0.50
1:C:15:THR:HB	1:C:16:PRO:CD	2.41	0.50
1:A:287:GLN:HE21	1:A:289:ILE:HD11	1.77	0.50
1:C:201:GLN:CG	1:C:211:TYR:OH	2.44	0.50
1:B:74:ARG:CZ	1:B:252:MET:CE	2.90	0.50
1:C:172:ILE:HD13	1:C:263:PHE:O	2.12	0.50
1:D:20:ARG:HG2	1:D:21:ARG:N	2.27	0.50
1:A:15:THR:HB	1:A:16:PRO:HD2	1.94	0.49
1:A:33:HIS:HE1	1:A:111:ASP:O	1.95	0.49
1:B:226:ASN:ND2	1:B:294:ARG:CB	2.75	0.49
1:A:89:GLN:HE21	1:C:107:ALA:CB	2.25	0.49
1:C:184:PHE:CZ	1:C:194:GLY:HA3	2.48	0.49

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:140:ARG:NH2	1:C:152:GLU:OE2	2.46	0.49
1:B:78:PRO:HD2	1:B:164:HIS:HB3	1.94	0.49
1:A:243:ASN:ND2	1:A:245:GLY:H	2.11	0.49
1:D:15:THR:HB	1:D:16:PRO:HD2	1.95	0.48
1:A:232:VAL:HB	1:A:288:THR:HB	1.95	0.48
1:C:238:ASN:ND2	1:C:272:THR:HB	2.29	0.48
1:C:162:ALA:HB2	1:C:269:VAL:HG22	1.96	0.48
1:B:83:TRP:HA	1:B:160:THR:O	2.14	0.48
1:C:51:LYS:HE2	2:C:339:HOH:O	2.13	0.48
1:B:281:LYS:HA	1:B:282:PRO:HD2	1.54	0.47
1:B:56:GLU:HB3	1:B:227:ARG:NH1	2.29	0.47
1:D:205:ASP:O	1:D:206:ARG:C	2.56	0.47
1:B:58:VAL:HG21	1:B:291:VAL:HG21	1.96	0.47
1:B:177:VAL:HG12	1:B:180:LEU:HD11	1.95	0.47
1:D:167:PHE:HE2	1:D:289:ILE:HD13	1.80	0.47
1:B:269:VAL:CG1	1:B:270:TYR:N	2.78	0.47
1:A:203:PHE:N	1:A:204:PRO:CD	2.78	0.47
1:B:38:ALA:HA	1:B:49:SER:O	2.15	0.47
1:B:109:ASN:HA	2:B:297:HOH:O	2.14	0.47
1:C:169:VAL:CG1	1:C:175:VAL:HG21	2.45	0.47
1:C:281:LYS:HA	1:C:282:PRO:HD2	1.79	0.47
1:B:43:GLN:HE22	1:B:70:GLY:H	1.63	0.46
1:C:201:GLN:OE1	1:C:209:ARG:NH2	2.48	0.46
1:A:60:TRP:CH2	1:A:62:SER:HB2	2.50	0.46
1:A:214:PRO:HG3	1:A:239:VAL:HG23	1.97	0.46
1:A:24:ASP:HB3	1:A:25:ASP:H	1.15	0.46
1:C:180:LEU:CD1	1:C:201:GLN:HB2	2.45	0.46
1:B:258:ASP:O	1:B:261:LYS:HG3	2.16	0.46
1:B:269:VAL:HG12	1:B:270:TYR:N	2.30	0.46
1:D:65:THR:HB	1:D:74:ARG:HB3	1.97	0.46
1:B:62:SER:HB3	1:B:65:THR:HB	1.97	0.46
1:B:287:GLN:HE21	1:B:289:ILE:HD11	1.81	0.45
1:C:15:THR:HB	1:C:16:PRO:HD2	1.98	0.45
1:B:131:TRP:HA	1:B:132:PRO:HD3	1.64	0.45
1:A:20:ARG:HD2	1:B:12:GLU:OE2	2.17	0.45
1:D:177:VAL:O	1:D:200:ILE:HA	2.17	0.45
1:A:44:GLY:O	1:A:45:ALA:C	2.58	0.44
1:A:141:PHE:HA	1:A:148:GLU:O	2.17	0.44
1:D:174:ASN:HB3	1:D:223:ALA:HB3	1.99	0.44
1:D:189:ASN:O	1:D:192:LYS:HG3	2.17	0.44
1:A:33:HIS:CD2	1:A:34:PRO:HD2	2.53	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:PRO:HD3	1:C:97:PHE:CE1	2.52	0.44
1:B:18:LEU:CD2	1:B:31:VAL:HG22	2.48	0.44
1:B:33:HIS:HD2	1:B:35:GLN:N	2.04	0.44
1:C:162:ALA:CB	1:C:269:VAL:HG22	2.47	0.44
1:C:287:GLN:HE21	1:C:289:ILE:HD11	1.82	0.44
1:A:109:ASN:C	1:A:109:ASN:ND2	2.75	0.44
1:B:160:THR:OG1	1:B:271:ALA:HA	2.17	0.44
1:A:84:PHE:HB3	1:A:160:THR:HG23	1.99	0.44
1:C:0:HIS:H2	1:C:2:ILE:H	1.64	0.44
1:A:13:GLN:NE2	1:A:15:THR:O	2.50	0.44
1:C:180:LEU:HD12	1:C:201:GLN:HB2	2.00	0.44
1:C:153:ALA:O	1:C:282:PRO:CB	2.66	0.44
1:A:17:VAL:O	1:A:17:VAL:CG1	2.66	0.43
1:B:208:ASP:C	1:B:209:ARG:HG3	2.42	0.43
1:A:17:VAL:HG22	1:A:32:ASP:O	2.18	0.43
1:A:252:MET:O	2:A:306:HOH:O	2.21	0.43
1:A:36:VAL:HG21	1:A:144:GLY:HA2	1.96	0.43
1:C:177:VAL:HA	1:C:219:VAL:O	2.19	0.43
1:B:140:ARG:NH2	1:B:152:GLU:OE1	2.51	0.43
1:D:12:GLU:OE2	1:D:20:ARG:NH1	2.52	0.43
1:A:188:VAL:O	1:A:188:VAL:HG12	2.19	0.43
1:B:8:LEU:O	1:B:21:ARG:HD3	2.19	0.43
1:A:222:ASP:C	1:A:222:ASP:OD1	2.61	0.43
1:B:35:GLN:HE21	1:B:35:GLN:HA	1.82	0.43
1:A:207:THR:HG23	1:A:209:ARG:HD3	2.00	0.42
1:D:33:HIS:CG	1:D:34:PRO:HD2	2.54	0.42
1:C:83:TRP:CE2	1:C:96:GLY:HA2	2.54	0.42
1:C:111:ASP:HB3	1:C:113:ASN:N	2.22	0.42
1:B:209:ARG:O	1:B:240:VAL:HA	2.19	0.42
1:C:62:SER:OG	1:C:256:PRO:HD3	2.19	0.42
1:D:180:LEU:HD13	1:D:211:TYR:CZ	2.55	0.42
1:A:64:ASN:HD22	1:A:64:ASN:HA	1.49	0.42
1:B:83:TRP:CB	1:B:160:THR:HG22	2.49	0.42
1:C:213:ASN:HA	1:C:214:PRO:HD2	1.65	0.42
1:A:103:TRP:CD1	1:A:103:TRP:N	2.84	0.42
1:D:165:SER:O	1:D:265:CYS:HA	2.19	0.42
1:B:84:PHE:N	1:B:160:THR:CG2	2.68	0.42
1:C:33:HIS:HD2	1:C:35:GLN:N	2.15	0.42
1:D:285:LEU:HD12	1:D:285:LEU:HA	1.88	0.42
1:A:155:GLY:HA3	1:A:157:PHE:CE2	2.55	0.42
1:B:75:GLY:HA2	1:B:166:TYR:CD1	2.54	0.42

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:ARG:HA	1:C:131:TRP:O	2.20	0.42
1:A:236:HIS:HD1	1:A:236:HIS:H	1.66	0.42
1:A:6:PHE:CZ	1:A:37:LYS:HB3	2.55	0.41
1:C:186:ASP:OD1	1:C:186:ASP:C	2.63	0.41
1:C:284:ARG:HD3	2:C:356:HOH:O	2.20	0.41
1:A:63:ASN:OD1	1:A:63:ASN:N	2.53	0.41
1:C:237:LEU:CD1	1:C:272:THR:HG22	2.50	0.41
1:A:249:SER:HB2	1:A:260:TYR:HB3	2.03	0.41
1:A:62:SER:C	1:A:64:ASN:H	2.28	0.41
1:A:162:ALA:CB	1:A:269:VAL:HG22	2.50	0.41
1:D:14:LEU:HD12	1:D:14:LEU:H	1.85	0.41
1:D:15:THR:HB	1:D:16:PRO:CD	2.49	0.41
1:D:44:GLY:C	1:D:46:HIS:N	2.75	0.41
1:B:225:LEU:O	1:B:226:ASN:C	2.63	0.41
1:B:104:ALA:O	1:B:119:PHE:HA	2.20	0.41
1:B:120:GLU:HB2	1:B:138:LEU:CD2	2.50	0.41
1:D:43:GLN:NE2	1:D:70:GLY:H	2.18	0.41
1:A:17:VAL:HG11	1:A:33:HIS:HB2	2.03	0.41
1:A:201:GLN:HE22	1:A:209:ARG:HH21	1.67	0.41
1:C:150:GLU:HG2	2:C:333:HOH:O	2.20	0.41
1:C:81:TRP:CG	1:C:82:PRO:HA	2.56	0.41
1:C:220:ILE:O	1:C:228:THR:HA	2.21	0.41
1:A:123:SER:HB3	1:A:135:PHE:CD1	2.56	0.40
1:A:272:THR:O	1:C:284:ARG:NH2	2.47	0.40
1:D:40:PHE:CD1	1:D:40:PHE:N	2.89	0.40
1:A:83:TRP:CA	1:A:160:THR:HG22	2.46	0.40
1:A:84:PHE:N	1:A:160:THR:CG2	2.78	0.40
1:B:83:TRP:HA	1:B:160:THR:CG2	2.45	0.40
1:D:24:ASP:HB3	1:D:25:ASP:H	1.55	0.40
1:A:220:ILE:O	1:A:228:THR:HA	2.22	0.40
1:B:167:PHE:O	1:B:263:PHE:HA	2.21	0.40
1:C:252:MET:O	2:C:358:HOH:O	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries

of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	294/309 (95%)	279 (95%)	15 (5%)	0	100	100
1	B	294/309 (95%)	278 (95%)	15 (5%)	1 (0%)	36	55
1	C	293/309 (95%)	278 (95%)	15 (5%)	0	100	100
1	D	294/309 (95%)	278 (95%)	16 (5%)	0	100	100
All	All	1175/1236 (95%)	1113 (95%)	61 (5%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	53	VAL

### 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	241/261 (92%)	219 (91%)	22 (9%)	9	19
1	B	236/261 (90%)	225 (95%)	11 (5%)	23	47
1	C	244/261 (94%)	223 (91%)	21 (9%)	10	21
1	D	243/261 (93%)	223 (92%)	20 (8%)	10	23
All	All	964/1044 (92%)	890 (92%)	74 (8%)	12	25

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	VAL
1	A	24	ASP
1	A	29	ILE
1	A	36	VAL
1	A	64	ASN

*Continued on next page...*

*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	89	GLN
1	A	109	ASN
1	A	137	LEU
1	A	150	GLU
1	A	151	LEU
1	A	160	THR
1	A	169	VAL
1	A	183	ARG
1	A	195	VAL
1	A	209	ARG
1	A	213	ASN
1	A	226	ASN
1	A	234	HIS
1	A	237	LEU
1	A	262	THR
1	A	291	VAL
1	A	293	LYS
1	B	47	LEU
1	B	56	GLU
1	B	137	LEU
1	B	148	GLU
1	B	156	GLU
1	B	160	THR
1	B	213	ASN
1	B	228	THR
1	B	258	ASP
1	B	261	LYS
1	B	291	VAL
1	C	13	GLN
1	C	17	VAL
1	C	18	LEU
1	C	36	VAL
1	C	68	LYS
1	C	90	GLN
1	C	116	MET
1	C	123	SER
1	C	128	ARG
1	C	137	LEU
1	C	156	GLU
1	C	159	THR
1	C	169	VAL
1	C	176	LYS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	190	ASP
1	C	201	GLN
1	C	209	ARG
1	C	226	ASN
1	C	237	LEU
1	C	272	THR
1	C	291	VAL
1	D	0	HIS
1	D	8	LEU
1	D	36	VAL
1	D	47	LEU
1	D	56	GLU
1	D	68	LYS
1	D	90	GLN
1	D	106	LYS
1	D	129	LYS
1	D	137	LEU
1	D	140	ARG
1	D	151	LEU
1	D	197	THR
1	D	210	VAL
1	D	228	THR
1	D	250	VAL
1	D	251	SER
1	D	272	THR
1	D	291	VAL
1	D	293	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	33	HIS
1	A	35	GLN
1	A	64	ASN
1	A	89	GLN
1	A	109	ASN
1	A	113	ASN
1	A	174	ASN
1	A	201	GLN
1	A	243	ASN
1	A	276	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	287	GLN
1	B	33	HIS
1	B	35	GLN
1	B	43	GLN
1	B	64	ASN
1	B	201	GLN
1	B	226	ASN
1	B	235	HIS
1	B	287	GLN
1	C	3	ASN
1	C	22	GLN
1	C	33	HIS
1	C	63	ASN
1	C	109	ASN
1	C	122	GLN
1	C	164	HIS
1	C	174	ASN
1	C	213	ASN
1	C	276	GLN
1	C	287	GLN
1	D	33	HIS
1	D	43	GLN
1	D	63	ASN
1	D	90	GLN
1	D	164	HIS

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	296/309 (95%)	-0.59	4 (1%) 73 70	17, 30, 45, 70	11 (3%)
1	B	296/309 (95%)	-0.42	6 (2%) 65 61	20, 36, 53, 60	12 (4%)
1	C	295/309 (95%)	-0.51	5 (1%) 69 65	17, 32, 46, 65	11 (3%)
1	D	296/309 (95%)	-0.67	6 (2%) 65 61	16, 28, 40, 59	11 (3%)
All	All	1183/1236 (95%)	-0.54	21 (1%) 67 64	16, 32, 47, 70	45 (3%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	174	ASN	6.4
1	C	174	ASN	6.3
1	B	174	ASN	6.3
1	A	174	ASN	5.4
1	C	0	HIS	4.4
1	B	212	LEU	4.3
1	B	132	PRO	3.8
1	D	132	PRO	3.6
1	A	132	PRO	3.5
1	C	9	PRO	3.5
1	C	132	PRO	3.5
1	B	9	PRO	3.5
1	D	212	LEU	3.3
1	D	-1	SER	3.1
1	A	9	PRO	3.1
1	B	176	LYS	2.9
1	C	176	LYS	2.7
1	D	22	GLN	2.6
1	D	9	PRO	2.5
1	B	22	GLN	2.2
1	A	176	LYS	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.