



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:02 PM UTC

PDB ID : 2PVP / pdb_00002pvp
Title : Crystal structure of D-Alanine-D-Alanine Ligase from Helicobacter pylori
Authors : Wu, D.; Zhang, L.; Jiang, H.; Shen, X.
Deposited on : 2007-05-10
Resolution : 2.40 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

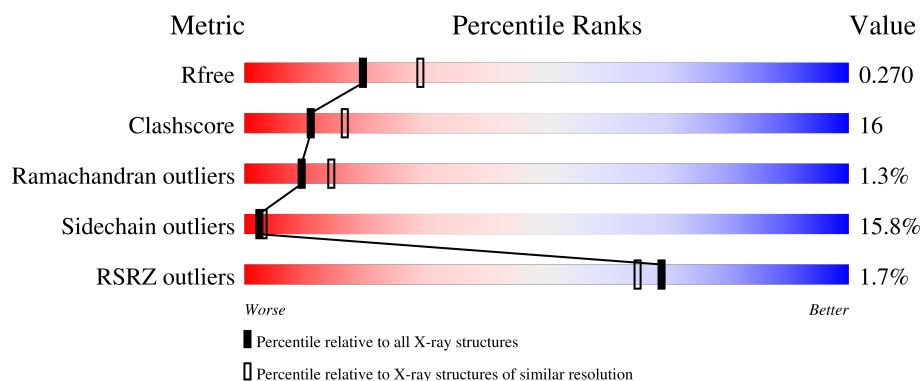
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	367	2% 55% 26% 8% 10%
1	B	367	% 57% 21% 8% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5286 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 D-alanine-D-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2660	1741	422	490	7			
1	B	315	Total	C	N	O	S	0	0	0
			2531	1656	401	467	7			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q2N4T5
A	-18	GLY	-	expression tag	UNP Q2N4T5
A	-17	SER	-	expression tag	UNP Q2N4T5
A	-16	SER	-	expression tag	UNP Q2N4T5
A	-15	HIS	-	expression tag	UNP Q2N4T5
A	-14	HIS	-	expression tag	UNP Q2N4T5
A	-13	HIS	-	expression tag	UNP Q2N4T5
A	-12	HIS	-	expression tag	UNP Q2N4T5
A	-11	HIS	-	expression tag	UNP Q2N4T5
A	-10	HIS	-	expression tag	UNP Q2N4T5
A	-9	SER	-	expression tag	UNP Q2N4T5
A	-8	SER	-	expression tag	UNP Q2N4T5
A	-7	GLY	-	expression tag	UNP Q2N4T5
A	-6	LEU	-	expression tag	UNP Q2N4T5
A	-5	VAL	-	expression tag	UNP Q2N4T5
A	-4	PRO	-	expression tag	UNP Q2N4T5
A	-3	ARG	-	expression tag	UNP Q2N4T5
A	-2	GLY	-	expression tag	UNP Q2N4T5
A	-1	SER	-	expression tag	UNP Q2N4T5
A	0	HIS	-	expression tag	UNP Q2N4T5
B	-19	MET	-	expression tag	UNP Q2N4T5
B	-18	GLY	-	expression tag	UNP Q2N4T5
B	-17	SER	-	expression tag	UNP Q2N4T5
B	-16	SER	-	expression tag	UNP Q2N4T5
B	-15	HIS	-	expression tag	UNP Q2N4T5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP Q2N4T5
B	-13	HIS	-	expression tag	UNP Q2N4T5
B	-12	HIS	-	expression tag	UNP Q2N4T5
B	-11	HIS	-	expression tag	UNP Q2N4T5
B	-10	HIS	-	expression tag	UNP Q2N4T5
B	-9	SER	-	expression tag	UNP Q2N4T5
B	-8	SER	-	expression tag	UNP Q2N4T5
B	-7	GLY	-	expression tag	UNP Q2N4T5
B	-6	LEU	-	expression tag	UNP Q2N4T5
B	-5	VAL	-	expression tag	UNP Q2N4T5
B	-4	PRO	-	expression tag	UNP Q2N4T5
B	-3	ARG	-	expression tag	UNP Q2N4T5
B	-2	GLY	-	expression tag	UNP Q2N4T5
B	-1	SER	-	expression tag	UNP Q2N4T5
B	0	HIS	-	expression tag	UNP Q2N4T5

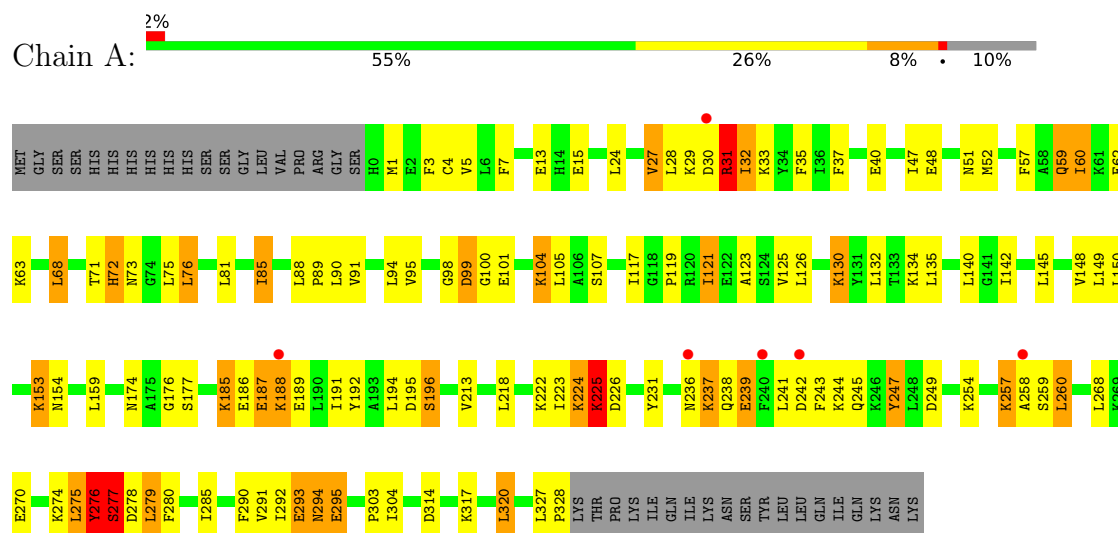
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	43	Total O 43 43	0	0
2	B	52	Total O 52 52	0	0

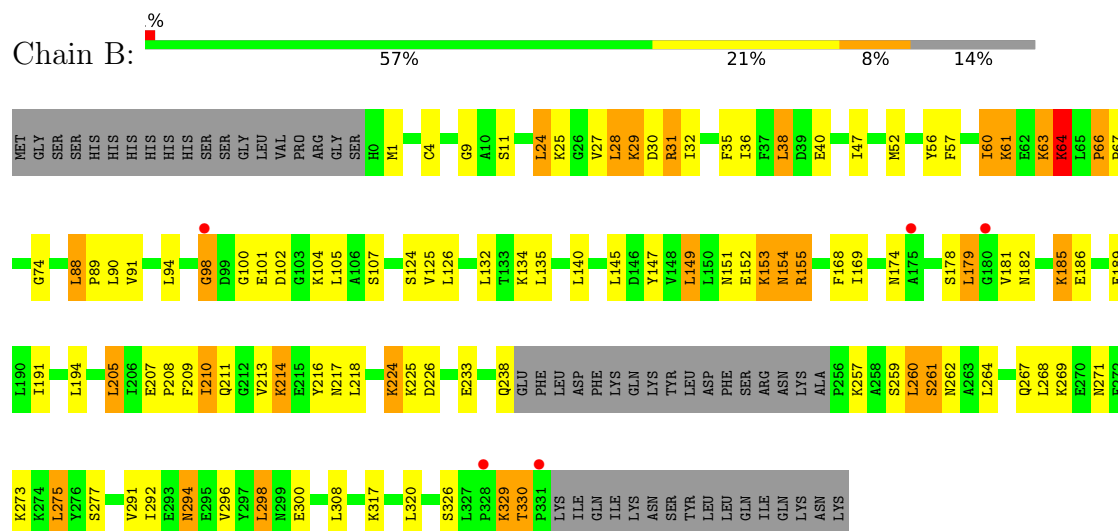
3 Residue-property plots [i](#)

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: D-alanine-D-alanine ligase



• Molecule 1: D-alanine-D-alanine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.97Å 89.02Å 121.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.40 15.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.5 (15.00-2.40) 97.1 (15.00-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.39Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.229 , 0.283 0.232 , 0.270	Depositor DCC
R_{free} test set	1521 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	48.1	Xtriage
Anisotropy	0.637	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5286	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.08	18/2720 (0.7%)	1.28	31/3667 (0.8%)
1	B	0.87	7/2587 (0.3%)	1.17	18/3489 (0.5%)
All	All	0.98	25/5307 (0.5%)	1.23	49/7156 (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	A	0	2
1	B	0	1
All	All	0	3

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	279	LEU	CB-CG	-19.24	1.15	1.53
1	A	277	SER	CB-OG	17.10	1.76	1.42
1	A	32	ILE	CB-CG1	-15.38	1.22	1.53
1	B	30	ASP	CB-CG	-11.24	1.24	1.52
1	B	225	LYS	CB-CG	-11.09	1.19	1.52
1	A	27	VAL	CB-CG2	-11.00	1.16	1.52
1	B	29	LYS	CA-CB	-10.83	1.32	1.52
1	B	31	ARG	CB-CG	10.18	1.82	1.52
1	B	178	SER	CB-OG	9.94	1.62	1.42
1	A	226	ASP	C-N	9.88	1.47	1.33
1	A	225	LYS	CB-CG	9.31	1.80	1.52
1	A	27	VAL	CB-CG1	8.66	1.81	1.52
1	B	27	VAL	C-N	8.21	1.46	1.33
1	B	28	LEU	CB-CG	7.83	1.69	1.53
1	A	295	GLU	C-N	-7.76	1.23	1.33
1	A	278	ASP	CB-CG	7.59	1.71	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	28	LEU	CB-CG	-7.52	1.38	1.53
1	A	72	HIS	CB-CG	7.37	1.60	1.50
1	A	280	PHE	CA-CB	6.67	1.64	1.53
1	A	257	LYS	C-N	-6.44	1.24	1.33
1	A	280	PHE	C-N	-5.89	1.25	1.33
1	A	275	LEU	C-N	5.80	1.41	1.33
1	A	292	ILE	C-N	5.69	1.41	1.33
1	A	260	LEU	C-N	5.43	1.40	1.33
1	A	226	ASP	CB-CG	-5.42	1.38	1.52

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	30	ASP	CA-CB-CG	29.66	142.26	112.60
1	A	276	TYR	CB-CG-CD1	-16.30	96.35	120.80
1	A	278	ASP	CA-CB-CG	-14.76	97.84	112.60
1	A	31	ARG	CB-CG-CD	14.45	144.53	111.30
1	A	276	TYR	CB-CG-CD2	13.29	140.73	120.80
1	A	27	VAL	CA-CB-CG1	-11.01	91.69	110.40
1	A	27	VAL	CA-CB-CG2	10.55	128.34	110.40
1	A	294	ASN	CB-CA-C	10.06	119.79	109.83
1	A	225	LYS	CA-CB-CG	-9.94	94.23	114.10
1	A	226	ASP	CA-CB-CG	9.74	122.34	112.60
1	B	28	LEU	CB-CG-CD2	9.35	138.75	110.70
1	A	257	LYS	O-C-N	-8.83	111.71	123.10
1	B	224	LYS	CB-CA-C	8.57	127.48	110.42
1	B	225	LYS	CA-CB-CG	8.50	131.09	114.10
1	B	30	ASP	CB-CG-OD1	8.45	137.84	118.40
1	A	280	PHE	N-CA-C	8.13	122.92	112.92
1	B	30	ASP	CB-CG-OD2	-8.03	99.94	118.40
1	A	32	ILE	CA-CB-CG2	-7.91	97.05	110.50
1	B	31	ARG	CA-CB-CG	-7.88	98.33	114.10
1	A	257	LYS	CA-C-N	7.87	133.72	121.99
1	A	257	LYS	C-N-CA	7.87	133.72	121.99
1	A	72	HIS	CB-CG-CD2	-7.78	121.08	131.20
1	A	99	ASP	CB-CA-C	-7.68	95.13	110.42
1	A	277	SER	N-CA-C	7.49	126.76	110.80
1	A	260	LEU	CB-CG-CD2	-6.84	90.19	110.70
1	A	276	TYR	O-C-N	6.81	130.85	122.34
1	A	72	HIS	CB-CG-ND1	6.68	132.73	122.70
1	A	99	ASP	N-CA-C	6.62	124.89	110.80
1	A	293	GLU	CA-C-N	-6.47	116.70	126.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	293	GLU	C-N-CA	-6.47	116.70	126.86
1	B	225	LYS	O-C-N	-6.41	113.52	122.37
1	A	278	ASP	CB-CG-OD1	6.38	133.08	118.40
1	B	178	SER	CA-CB-OG	-6.28	98.54	111.10
1	B	29	LYS	CB-CA-C	6.10	123.56	116.54
1	B	27	VAL	O-C-N	5.92	133.16	121.91
1	B	225	LYS	N-CA-CB	-5.77	101.61	110.44
1	A	224	LYS	CB-CA-C	-5.71	109.47	117.23
1	A	275	LEU	CA-C-N	-5.70	113.14	122.54
1	A	275	LEU	C-N-CA	-5.70	113.14	122.54
1	A	280	PHE	N-CA-CB	-5.54	102.47	110.56
1	B	300	GLU	N-CA-C	5.33	116.39	108.60
1	A	225	LYS	CB-CG-CD	-5.32	99.07	111.30
1	B	28	LEU	CA-CB-CG	-5.25	97.92	116.30
1	B	225	LYS	CB-CG-CD	-5.23	99.28	111.30
1	A	88	LEU	CA-C-N	-5.14	115.43	120.52
1	A	88	LEU	C-N-CA	-5.14	115.43	120.52
1	B	224	LYS	N-CA-CB	5.14	119.18	110.49
1	B	66	PRO	CA-C-N	5.09	125.33	119.83
1	B	66	PRO	C-N-CA	5.09	125.33	119.83

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	276	TYR	Sidechain
1	A	30	ASP	Peptide
1	B	63	LYS	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2660	0	2679	113	1
1	B	2531	0	2558	62	1
2	A	43	0	0	3	0
2	B	52	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5286	0	5237	172	1

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 16.

All (172) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:ARG:CB	1:B:31:ARG:CG	1.83	1.56
1:A:225:LYS:CB	1:A:225:LYS:CG	1.80	1.54
1:A:27:VAL:CB	1:A:27:VAL:CG1	1.81	1.54
1:A:231:TYR:CE1	1:A:242:ASP:OD1	1.81	1.32
1:A:277:SER:CB	1:A:277:SER:OG	1.76	1.32
1:A:244:LYS:NZ	1:A:259:SER:OG	1.74	1.20
1:B:179:LEU:CD1	2:B:386:HOH:O	1.97	1.11
1:B:64:LYS:O	1:B:64:LYS:HG3	1.45	1.07
1:A:35:PHE:HB3	1:A:52:MET:HE3	1.34	1.07
1:A:35:PHE:CB	1:A:52:MET:HE3	1.86	1.06
1:A:231:TYR:CD1	1:A:242:ASP:HB2	1.92	1.05
1:A:27:VAL:CG1	1:A:27:VAL:CA	2.41	0.99
1:A:225:LYS:CG	1:A:225:LYS:CA	2.44	0.94
1:A:231:TYR:CE1	1:A:242:ASP:HB2	2.02	0.94
1:A:35:PHE:HB3	1:A:52:MET:CE	1.99	0.92
1:A:32:ILE:HG22	1:A:32:ILE:O	1.68	0.92
1:B:179:LEU:HD13	2:B:386:HOH:O	1.64	0.91
1:A:27:VAL:CG1	1:A:27:VAL:CG2	2.46	0.90
1:A:231:TYR:CD1	1:A:242:ASP:CB	2.57	0.88
1:A:123:ALA:HB2	1:A:279:LEU:O	1.73	0.87
1:A:231:TYR:CD1	1:A:242:ASP:OD1	2.28	0.86
1:A:225:LYS:CB	1:A:225:LYS:CD	2.54	0.85
1:A:241:LEU:HD22	1:A:258:ALA:O	1.77	0.84
1:A:231:TYR:CE1	1:A:242:ASP:CG	2.55	0.84
1:B:31:ARG:CG	1:B:31:ARG:CA	2.55	0.84
1:A:35:PHE:CG	1:A:52:MET:HE3	2.14	0.83
1:A:231:TYR:HD1	1:A:242:ASP:HB2	1.45	0.82
1:A:231:TYR:CE1	1:A:242:ASP:CB	2.65	0.80
1:B:64:LYS:O	1:B:64:LYS:CG	2.30	0.80
1:B:179:LEU:HD11	2:B:386:HOH:O	1.72	0.79
1:B:213:VAL:HG21	1:B:292:ILE:HG12	1.64	0.79
1:A:98:GLY:O	1:A:99:ASP:HB2	1.84	0.77
1:B:329:LYS:C	1:B:330:THR:HG22	2.09	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:CYS:SG	1:A:89:PRO:HD2	2.25	0.76
1:B:4:CYS:SG	1:B:36:ILE:CD1	2.74	0.76
1:B:28:LEU:HB2	1:B:32:ILE:HD11	1.69	0.76
1:A:239:GLU:OE2	2:A:376:HOH:O	2.04	0.75
1:A:231:TYR:CD1	1:A:242:ASP:CG	2.65	0.74
1:A:247:TYR:H	1:A:247:TYR:HD1	1.37	0.73
1:A:222:LYS:NZ	1:A:277:SER:O	2.24	0.70
1:B:9:GLY:HA2	1:B:38:LEU:HD13	1.74	0.70
1:B:260:LEU:HD22	1:B:264:LEU:HD23	1.75	0.68
1:A:231:TYR:HD1	1:A:242:ASP:CB	2.03	0.68
1:B:40:GLU:H	1:B:40:GLU:CD	2.02	0.68
1:A:121:ILE:HD12	1:A:121:ILE:N	2.09	0.68
1:B:4:CYS:SG	1:B:36:ILE:HD11	2.33	0.67
1:B:74:GLY:HA2	1:B:88:LEU:HG	1.77	0.67
1:B:4:CYS:SG	1:B:36:ILE:HD12	2.33	0.66
1:A:276:TYR:CD2	1:A:276:TYR:O	2.49	0.66
1:A:32:ILE:HG21	1:A:35:PHE:CE1	2.32	0.64
1:A:225:LYS:CG	1:A:225:LYS:HA	2.28	0.64
1:A:285:ILE:HG22	1:A:303:PRO:HA	1.80	0.64
1:A:270:GLU:O	1:A:274:LYS:HG2	1.99	0.62
1:A:35:PHE:CB	1:A:52:MET:CE	2.67	0.62
1:B:329:LYS:C	1:B:330:THR:CG2	2.73	0.61
1:A:153:LYS:O	1:A:153:LYS:HG2	2.01	0.60
1:A:191:ILE:HD12	1:A:191:ILE:H	1.67	0.60
1:A:231:TYR:CZ	1:A:242:ASP:OD1	2.50	0.60
1:A:37:PHE:HB2	1:A:47:ILE:HD13	1.84	0.60
1:A:48:GLU:HB2	1:A:51:ASN:HD22	1.66	0.60
1:B:224:LYS:HD2	1:B:326:SER:HB2	1.84	0.59
1:A:31:ARG:NH1	1:A:317:LYS:CD	2.66	0.59
1:A:37:PHE:HB3	1:A:47:ILE:HD11	1.85	0.59
1:A:27:VAL:CG1	1:A:27:VAL:HA	2.33	0.58
1:A:31:ARG:NH1	1:A:317:LYS:HD2	2.19	0.58
1:A:125:VAL:HG21	1:B:125:VAL:HG21	1.86	0.58
1:A:192:TYR:O	1:A:196:SER:HB3	2.03	0.58
1:B:31:ARG:HG3	1:B:317:LYS:NZ	2.19	0.57
1:A:95:VAL:O	1:A:100:GLY:HA3	2.04	0.57
1:B:28:LEU:HB2	1:B:32:ILE:CD1	2.34	0.57
1:B:56:TYR:CE1	1:B:63:LYS:HE2	2.39	0.57
1:A:291:VAL:HA	1:A:295:GLU:O	2.05	0.57
1:B:216:TYR:CZ	1:B:260:LEU:HD21	2.40	0.56
1:B:214:LYS:HB2	1:B:291:VAL:HB	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:ARG:CB	1:B:31:ARG:CD	2.80	0.56
1:A:32:ILE:HG21	1:A:35:PHE:CZ	2.41	0.56
1:B:98:GLY:O	1:B:102:ASP:CG	2.49	0.55
1:B:275:LEU:HD21	1:B:298:LEU:HG	1.89	0.55
1:A:35:PHE:CD2	1:A:52:MET:HE3	2.41	0.55
1:B:61:LYS:C	1:B:63:LYS:H	2.15	0.54
1:A:13:GLU:OE1	1:A:177:SER:HB2	2.08	0.54
1:A:24:LEU:HA	1:A:27:VAL:HG22	1.89	0.54
1:A:48:GLU:HB2	1:A:51:ASN:ND2	2.22	0.54
1:A:121:ILE:N	1:A:121:ILE:CD1	2.71	0.54
1:B:154:ASN:HD22	1:B:154:ASN:C	2.15	0.54
1:A:57:PHE:HA	1:A:60:ILE:HB	1.91	0.53
1:A:72:HIS:H	1:A:72:HIS:CD2	2.24	0.53
1:A:225:LYS:CB	1:A:225:LYS:HD3	2.35	0.53
1:B:217:ASN:HB2	1:B:233:GLU:HB3	1.90	0.53
1:A:159:LEU:HD11	1:A:187:GLU:HB2	1.90	0.53
1:A:37:PHE:HB2	1:A:47:ILE:CD1	2.38	0.53
1:A:231:TYR:HE1	1:A:242:ASP:HB2	1.67	0.52
1:A:121:ILE:CD1	1:A:121:ILE:H	2.22	0.52
1:A:231:TYR:HE1	1:A:242:ASP:CB	2.20	0.52
1:B:169:ILE:HD11	1:B:181:VAL:HG12	1.90	0.52
1:A:37:PHE:CB	1:A:47:ILE:HD11	2.40	0.52
1:A:237:LYS:O	1:A:238:GLN:HB2	2.09	0.51
1:B:24:LEU:HG	1:B:94:LEU:HD11	1.91	0.51
1:A:5:VAL:HG13	1:A:94:LEU:HD12	1.93	0.51
1:A:31:ARG:NH1	1:A:317:LYS:HD3	2.26	0.51
1:B:149:LEU:HD23	1:B:205:LEU:HD13	1.92	0.51
1:A:130:LYS:HE3	1:A:174:ASN:HA	1.94	0.50
1:A:37:PHE:CB	1:A:47:ILE:CD1	2.89	0.50
1:A:72:HIS:O	1:A:73:ASN:HB2	2.11	0.50
1:A:134:LYS:HD2	2:A:360:HOH:O	2.12	0.49
1:A:121:ILE:HD12	1:A:121:ILE:H	1.75	0.49
1:A:213:VAL:HG21	1:A:290:PHE:HB3	1.93	0.49
1:A:89:PRO:O	1:A:90:LEU:C	2.55	0.49
1:A:4:CYS:HG	1:A:89:PRO:HD2	1.78	0.48
1:A:47:ILE:CG2	1:A:52:MET:HE2	2.44	0.48
1:A:186:GLU:O	1:A:188:LYS:N	2.46	0.48
1:B:134:LYS:HD3	1:B:147:TYR:CD2	2.49	0.47
1:B:168:PHE:HB2	1:B:207:GLU:O	2.14	0.47
1:A:98:GLY:O	1:A:99:ASP:CB	2.58	0.47
1:B:269:LYS:O	1:B:273:LYS:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:ASN:H	1:B:154:ASN:HD21	1.63	0.47
1:A:99:ASP:O	1:A:105:LEU:HD12	2.15	0.47
1:A:7:PHE:CZ	1:A:52:MET:HE1	2.51	0.46
1:A:185:LYS:N	1:A:189:GLU:OE2	2.33	0.46
1:B:35:PHE:CD2	1:B:52:MET:HE2	2.51	0.46
1:B:291:VAL:CG1	1:B:294:ASN:HA	2.45	0.46
1:A:119:PRO:HG3	1:A:303:PRO:HB2	1.97	0.46
1:A:247:TYR:CD1	1:A:247:TYR:N	2.82	0.46
1:B:57:PHE:CD2	1:B:60:ILE:HG13	2.51	0.46
1:A:71:THR:HG21	1:A:85:ILE:HD11	1.98	0.45
1:B:330:THR:O	1:B:330:THR:OG1	2.30	0.45
1:A:223:ILE:O	1:A:224:LYS:HB2	2.17	0.44
1:B:271:ASN:HD22	1:B:296:VAL:HG11	1.81	0.44
1:B:261:SER:HB2	1:B:264:LEU:H	1.82	0.44
1:A:32:ILE:CG2	1:A:35:PHE:CE1	2.98	0.44
1:A:187:GLU:O	1:A:187:GLU:CG	2.62	0.44
1:B:185:LYS:HG2	1:B:189:GLU:OE2	2.17	0.44
1:A:59:GLN:O	1:A:62:GLU:HG2	2.16	0.44
1:B:89:PRO:O	1:B:90:LEU:C	2.61	0.44
1:B:154:ASN:HD22	1:B:155:ARG:N	2.16	0.44
1:A:40:GLU:H	1:A:40:GLU:CD	2.26	0.44
1:A:104:LYS:HA	1:B:107:SER:HB3	1.99	0.44
1:A:225:LYS:CA	1:A:225:LYS:HG3	2.44	0.44
1:A:242:ASP:OD2	1:A:243:PHE:N	2.39	0.44
1:A:68:LEU:HD23	1:A:75:LEU:HD13	2.00	0.43
1:A:293:GLU:O	1:A:294:ASN:HB2	2.17	0.43
1:B:66:PRO:HA	1:B:67:PRO:HD3	1.80	0.43
1:B:174:ASN:OD1	1:B:174:ASN:O	2.37	0.43
1:B:291:VAL:HG13	1:B:294:ASN:HA	2.00	0.43
1:A:3:PHE:CZ	1:A:320:LEU:HD13	2.53	0.43
1:A:188:LYS:HA	1:A:188:LYS:HD2	1.64	0.43
1:A:99:ASP:O	1:A:105:LEU:CG	2.67	0.43
1:B:207:GLU:HB3	1:B:208:PRO:HD2	2.01	0.43
1:A:99:ASP:O	1:A:105:LEU:HG	2.19	0.42
1:B:11:SER:HB2	1:B:98:GLY:HA3	2.02	0.42
1:B:31:ARG:HG3	1:B:317:LYS:HZ2	1.84	0.42
1:A:213:VAL:CG2	1:A:290:PHE:HB3	2.50	0.42
1:B:74:GLY:HA2	1:B:88:LEU:CG	2.47	0.42
1:A:254:LYS:HG2	2:A:387:HOH:O	2.18	0.42
1:A:148:VAL:HG12	1:A:150:LEU:CD2	2.49	0.42
1:B:61:LYS:C	1:B:63:LYS:N	2.77	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:LEU:HA	1:A:328:PRO:HD3	1.92	0.42
1:A:225:LYS:HD3	1:A:225:LYS:HB2	2.02	0.41
1:A:153:LYS:O	1:A:153:LYS:CG	2.68	0.41
1:B:31:ARG:CA	1:B:31:ARG:HG2	2.48	0.41
1:B:210:ILE:HG22	2:B:399:HOH:O	2.20	0.41
1:A:31:ARG:HH12	1:A:317:LYS:CD	2.32	0.41
1:A:101:GLU:CG	1:A:304:ILE:HD12	2.50	0.41
1:A:244:LYS:NZ	1:A:259:SER:HG	2.06	0.41
1:A:107:SER:HB3	1:B:104:LYS:HA	2.03	0.41
1:A:142:ILE:HD11	1:A:274:LYS:HG3	2.03	0.41
1:A:191:ILE:HD12	1:A:191:ILE:N	2.33	0.41
1:A:31:ARG:HH12	1:A:317:LYS:HD2	1.86	0.41
1:B:101:GLU:O	1:B:124:SER:HB3	2.20	0.40
1:B:209:PHE:CE2	1:B:211:GLN:HB3	2.57	0.40
1:A:76:LEU:HD22	1:A:85:ILE:CD1	2.52	0.40
1:B:100:GLY:HA2	1:B:105:LEU:HD12	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:ASP:OD1	1:B:261:SER:OG[1_655]	1.93	0.27

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	327/367 (89%)	307 (94%)	17 (5%)	3 (1%)	14	22
1	B	311/367 (85%)	293 (94%)	13 (4%)	5 (2%)	7	11
All	All	638/734 (87%)	600 (94%)	30 (5%)	8 (1%)	9	14

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	SER
1	A	154	ASN
1	B	64	LYS
1	B	98	GLY
1	B	153	LYS
1	B	277	SER
1	B	155	ARG
1	A	176	GLY

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	291/326 (89%)	247 (85%)	44 (15%)	3	3
1	B	278/326 (85%)	232 (84%)	46 (16%)	2	3
All	All	569/652 (87%)	479 (84%)	90 (16%)	2	3

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	15	GLU
1	A	29	LYS
1	A	31	ARG
1	A	33	LYS
1	A	59	GLN
1	A	60	ILE
1	A	63	LYS
1	A	68	LEU
1	A	76	LEU
1	A	81	LEU
1	A	85	ILE
1	A	91	VAL
1	A	104	LYS
1	A	117	ILE

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Mol	Chain	Res	Type
1	A	121	ILE
1	A	126	LEU
1	A	130	LYS
1	A	132	LEU
1	A	135	LEU
1	A	140	LEU
1	A	145	LEU
1	A	149	LEU
1	A	153	LYS
1	A	185	LYS
1	A	187	GLU
1	A	188	LYS
1	A	194	LEU
1	A	195	ASP
1	A	196	SER
1	A	218	LEU
1	A	225	LYS
1	A	236	ASN
1	A	237	LYS
1	A	239	GLU
1	A	245	GLN
1	A	247	TYR
1	A	257	LYS
1	A	260	LEU
1	A	268	LEU
1	A	275	LEU
1	A	276	TYR
1	A	314	ASP
1	A	320	LEU
1	B	1	MET
1	B	24	LEU
1	B	25	LYS
1	B	29	LYS
1	B	38	LEU
1	B	47	ILE
1	B	60	ILE
1	B	61	LYS
1	B	64	LYS
1	B	88	LEU
1	B	91	VAL
1	B	126	LEU
1	B	132	LEU

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Mol	Chain	Res	Type
1	B	135	LEU
1	B	140	LEU
1	B	145	LEU
1	B	149	LEU
1	B	152	GLU
1	B	153	LYS
1	B	154	ASN
1	B	179	LEU
1	B	182	ASN
1	B	185	LYS
1	B	186	GLU
1	B	191	ILE
1	B	194	LEU
1	B	205	LEU
1	B	210	ILE
1	B	214	LYS
1	B	218	LEU
1	B	226	ASP
1	B	238	GLN
1	B	257	LYS
1	B	259	SER
1	B	260	LEU
1	B	261	SER
1	B	262	ASN
1	B	267	GLN
1	B	268	LEU
1	B	275	LEU
1	B	294	ASN
1	B	298	LEU
1	B	308	LEU
1	B	320	LEU
1	B	329	LYS
1	B	330	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	HIS
1	A	51	ASN
1	A	53	HIS
1	A	72	HIS
1	A	163	ASN

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Mol	Chain	Res	Type
1	A	217	ASN
1	A	245	GLN
1	A	262	ASN
1	A	271	ASN
1	A	322	ASN
1	A	325	GLN
1	B	59	GLN
1	B	154	ASN
1	B	165	ASN
1	B	174	ASN
1	B	238	GLN
1	B	271	ASN
1	B	294	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	329/367 (89%)	-0.03	6 (1%) 67 63	20, 51, 82, 102	0
1	B	315/367 (85%)	-0.02	5 (1%) 70 66	20, 58, 86, 118	0
All	All	644/734 (87%)	-0.02	11 (1%) 69 65	20, 54, 85, 118	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	236	ASN	7.5
1	B	328	PRO	6.5
1	A	242	ASP	5.9
1	A	258	ALA	4.4
1	A	188	LYS	3.5
1	B	331	PRO	3.0
1	B	175	ALA	2.7
1	B	98	GLY	2.2
1	A	30	ASP	2.2
1	B	180	GLY	2.2
1	A	240	PHE	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.