



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 24, 2026 – 12:53 AM UTC

PDB ID : 2D2R / pdb_00002d2r
Title : Crystal structure of Helicobacter pylori Undecaprenyl Pyrophosphate Synthase
Authors : Kuo, C.J.; Guo, R.T.; Chen, C.L.; Ko, T.P.; Cheng, Y.S.; Cheng, Y.L.; Liang, P.H.; Wang, A.H.-J.
Deposited on : 2005-09-16
Resolution : 1.88 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

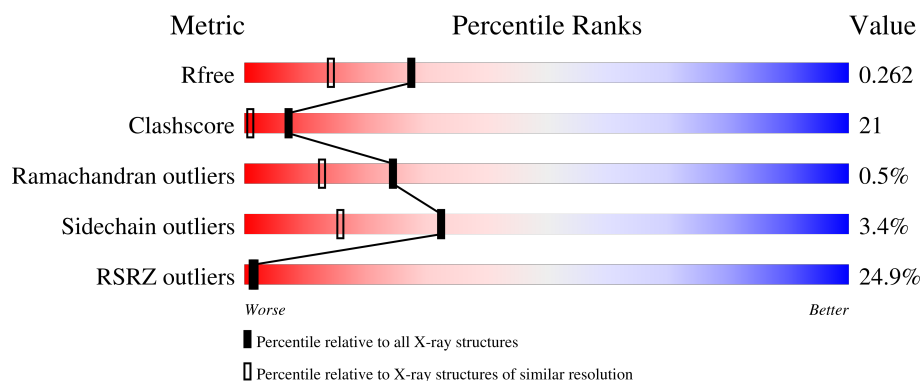
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	245	
1	B	245	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4035 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 Undecaprenyl Pyrophosphate Synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	0	0
			1772	1137	305	324	6			
1	B	203	Total	C	N	O	S	0	0	0
			1682	1080	293	303	6			

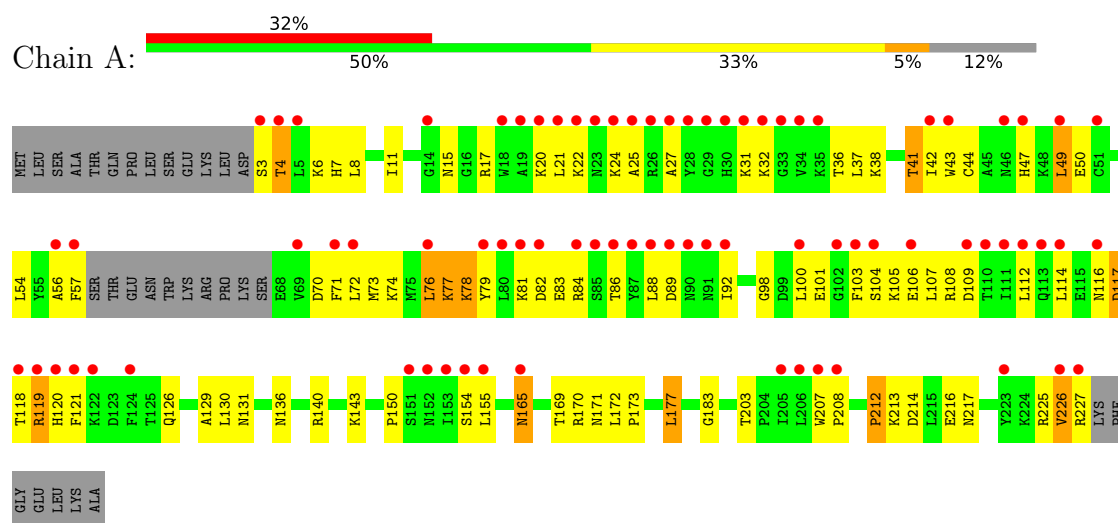
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	252	Total	O	0	0
			252	252		
2	B	329	Total	O	0	0
			329	329		

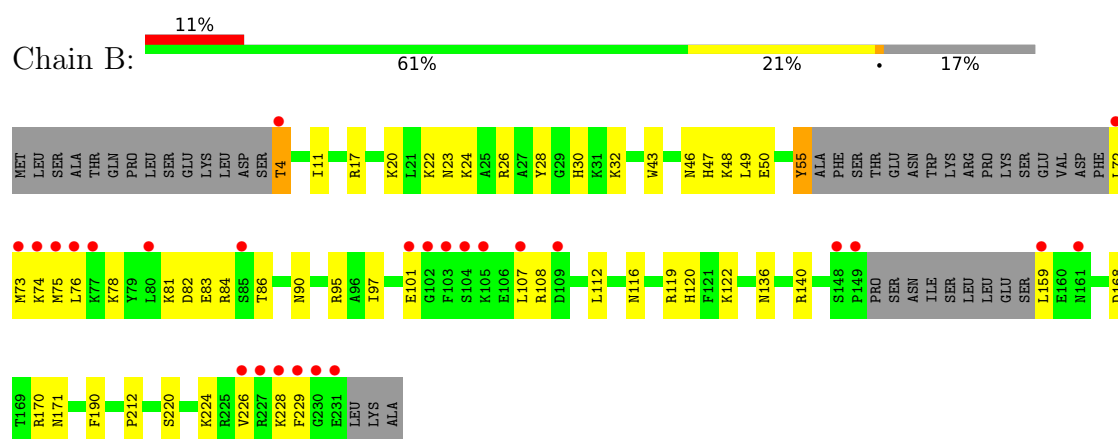
3 Residue-property plots [i](#)

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: Undecaprenyl Pyrophosphate Synthase



• Molecule 1: Undecaprenyl Pyrophosphate Synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.63Å 58.91Å 153.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 1.88 25.00 – 1.88	Depositor EDS
% Data completeness (in resolution range)	92.4 (25.00-1.88) 92.4 (25.00-1.88)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.19 (at 1.88Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.193 , 0.240 0.242 , 0.262	Depositor DCC
R_{free} test set	1719 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	9.0	Xtriage
Anisotropy	0.411	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 69.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.29$, $\langle L^2 \rangle = 0.13$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	4035	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.88	1/1807 (0.1%)	1.18	14/2432 (0.6%)
1	B	1.15	0/1714	1.22	7/2301 (0.3%)
All	All	1.02	1/3521 (0.0%)	1.20	21/4733 (0.4%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	165	ASN	CG-OD1	5.66	1.34	1.23

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	170	ARG	CA-C-N	-11.46	104.27	122.73
1	B	170	ARG	C-N-CA	-11.46	104.27	122.73
1	A	22	LYS	N-CA-C	-7.88	103.23	112.92
1	A	78	LYS	N-CA-C	-7.61	102.93	111.07
1	A	203	THR	CA-C-N	6.95	126.58	119.56
1	A	203	THR	C-N-CA	6.95	126.58	119.56
1	B	73	MET	N-CA-C	-6.60	103.92	112.23
1	A	117	ASP	N-CA-C	6.20	118.84	111.71
1	A	49	LEU	N-CA-C	-6.16	99.61	109.96
1	A	150	PRO	N-CA-C	-5.99	101.86	111.38
1	A	98	GLY	N-CA-C	-5.84	104.92	111.63
1	B	48	LYS	N-CA-C	5.68	119.51	112.24
1	A	172	LEU	N-CA-C	-5.62	102.82	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	11	ILE	N-CA-C	-5.60	99.17	107.51
1	A	165	ASN	CB-CA-C	5.59	120.07	110.01
1	A	183	GLY	N-CA-C	5.45	121.33	114.25
1	A	131	ASN	N-CA-C	-5.39	103.95	111.39
1	A	4	THR	N-CA-C	-5.19	106.46	112.89
1	B	49	LEU	N-CA-C	-5.08	103.82	110.53
1	B	11	ILE	N-CA-C	-5.07	99.96	107.51
1	B	82	ASP	N-CA-C	5.04	116.77	111.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	55	TYR	Sidechain

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	1772	0	1795	86	0
1	B	1682	0	1711	57	0
2	A	252	0	0	22	0
2	B	329	0	0	30	0
All	All	4035	0	3506	143	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 21.

All (143) 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:73:MET:HE1	1:A:104:SER:N	1.69	1.06
1:A:73:MET:HE1	1:A:104:SER:H	0.85	1.01
1:A:165:ASN:HB3	2:A:347:HOH:O	1.64	0.96
1:A:37:LEU:O	1:A:41:THR:HG23	1.64	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:MET:CE	1:A:104:SER:H	1.80	0.92
1:A:105:LYS:HD3	1:A:108:ARG:HH21	1.39	0.86
1:A:136:ASN:HD21	1:A:140:ARG:HE	1.23	0.85
1:A:112:LEU:HD22	2:A:409:HOH:O	1.79	0.82
1:A:83:GLU:O	1:A:86:THR:HG22	1.81	0.80
1:A:20:LYS:HD2	1:A:21:LEU:HD23	1.62	0.79
1:A:136:ASN:ND2	1:A:140:ARG:HE	1.80	0.78
1:A:31:LYS:HD3	2:A:358:HOH:O	1.84	0.76
1:B:23:ASN:HB2	2:B:535:HOH:O	1.84	0.76
1:B:20:LYS:HD2	2:B:375:HOH:O	1.86	0.74
1:A:41:THR:HG21	1:A:126:GLN:HE21	1.51	0.73
1:B:159:LEU:N	2:B:339:HOH:O	2.21	0.73
1:B:17:ARG:NH2	2:B:305:HOH:O	2.22	0.72
1:A:77:LYS:HZ3	1:A:81:LYS:HD2	1.54	0.72
1:B:76:LEU:HD23	1:B:107:LEU:HD21	1.72	0.70
1:A:83:GLU:HG3	2:A:448:HOH:O	1.91	0.69
1:B:4:THR:N	2:B:525:HOH:O	2.24	0.69
1:A:73:MET:HE3	1:A:107:LEU:HD23	1.74	0.68
1:A:15:ASN:HD22	1:A:208:PRO:HD3	1.57	0.68
1:B:30:HIS:HB3	1:B:75:MET:HE1	1.76	0.68
1:B:78:LYS:HG2	2:B:236:HOH:O	1.95	0.66
1:A:106:GLU:HG2	2:A:438:HOH:O	1.94	0.66
1:B:136:ASN:ND2	1:B:140:ARG:HE	1.93	0.66
1:A:77:LYS:HZ3	1:A:81:LYS:CD	2.08	0.65
1:A:43:TRP:O	1:A:47:HIS:HD2	1.80	0.65
1:A:101:GLU:HA	1:A:108:ARG:HH11	1.61	0.64
1:B:122:LYS:O	1:B:122:LYS:HG3	1.97	0.64
1:B:136:ASN:HD21	1:B:140:ARG:HE	1.43	0.64
1:A:116:ASN:ND2	1:A:119:ARG:HH22	1.96	0.63
2:A:458:HOH:O	1:B:228:LYS:HD3	1.99	0.61
1:A:56:ALA:HB3	1:A:129:ALA:O	1.99	0.61
1:A:89:ASP:OD2	2:A:362:HOH:O	2.16	0.61
1:A:116:ASN:O	1:A:119:ARG:HG2	2.01	0.61
1:B:26:ARG:HD2	2:B:524:HOH:O	2.02	0.60
1:B:50:GLU:HG2	2:B:540:HOH:O	2.01	0.60
1:B:95:ARG:HH22	1:B:171:ASN:H	1.48	0.60
1:A:37:LEU:HG	1:A:54:LEU:HD13	1.82	0.59
1:A:105:LYS:O	1:A:109:ASP:HB2	2.03	0.59
1:A:76:LEU:O	1:A:79:TYR:HB3	2.03	0.58
1:B:24:LYS:HD3	1:B:28:TYR:CD2	2.39	0.58
1:B:72:LEU:HA	2:B:561:HOH:O	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:SER:HB2	2:A:342:HOH:O	2.04	0.58
1:A:44:CYS:HB3	1:A:49:LEU:HD22	1.86	0.57
1:B:26:ARG:CD	2:B:524:HOH:O	2.51	0.57
1:A:76:LEU:HD11	1:A:130:LEU:HD22	1.87	0.57
1:A:81:LYS:HE2	2:A:474:HOH:O	2.05	0.56
1:A:101:GLU:HA	1:A:108:ARG:NH1	2.19	0.56
1:A:121:PHE:N	1:A:121:PHE:CD1	2.69	0.56
1:B:108:ARG:O	1:B:112:LEU:HG	2.05	0.56
1:A:77:LYS:HG3	1:A:78:LYS:N	2.21	0.56
1:A:72:LEU:HB2	2:A:395:HOH:O	2.05	0.56
1:B:226:VAL:HG13	1:B:229:PHE:HE2	1.72	0.55
1:A:108:ARG:O	1:A:112:LEU:HG	2.06	0.55
1:A:43:TRP:CD1	1:A:212:PRO:HB3	2.42	0.55
1:B:22:LYS:HB2	1:B:24:LYS:HG3	1.89	0.55
1:B:101:GLU:HG2	2:B:314:HOH:O	2.06	0.54
1:A:3:SER:HB3	2:A:266:HOH:O	2.06	0.53
1:A:41:THR:HG21	1:A:126:GLN:NE2	2.23	0.53
1:A:118:THR:C	1:A:120:HIS:H	2.16	0.53
1:A:105:LYS:CD	1:A:108:ARG:HH21	2.18	0.52
1:A:77:LYS:HD3	1:A:81:LYS:HD3	1.89	0.52
1:A:15:ASN:ND2	1:A:208:PRO:HD3	2.24	0.52
1:A:114:LEU:O	1:A:118:THR:HG23	2.09	0.52
1:B:43:TRP:O	1:B:47:HIS:HD2	1.93	0.52
1:A:82:ASP:OD2	2:A:449:HOH:O	2.19	0.52
1:A:84:ARG:NH2	1:A:117:ASP:OD2	2.43	0.51
1:A:17:ARG:NH2	2:A:294:HOH:O	2.44	0.51
1:B:24:LYS:O	2:B:376:HOH:O	2.19	0.51
1:A:43:TRP:O	1:A:47:HIS:CD2	2.62	0.50
1:B:55:TYR:HB2	1:B:190:PHE:CE1	2.46	0.50
1:A:84:ARG:NH1	1:A:88:LEU:HD11	2.27	0.50
1:A:49:LEU:HD12	2:A:364:HOH:O	2.12	0.50
1:A:119:ARG:NH1	2:A:400:HOH:O	2.43	0.50
1:B:72:LEU:CA	2:B:561:HOH:O	2.60	0.50
1:B:226:VAL:HG13	1:B:229:PHE:CE2	2.47	0.49
1:A:37:LEU:HG	1:A:54:LEU:CD1	2.43	0.49
1:B:30:HIS:HB3	1:B:75:MET:CE	2.43	0.49
1:B:32:LYS:NZ	2:B:403:HOH:O	2.46	0.49
1:B:30:HIS:CB	1:B:75:MET:HE1	2.42	0.49
1:A:207:TRP:HB3	1:A:208:PRO:HD3	1.94	0.48
1:A:213:LYS:O	1:A:217:ASN:ND2	2.46	0.48
1:A:83:GLU:HG2	1:A:86:THR:HG21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:HIS:H	1:B:120:HIS:CD2	2.31	0.48
1:A:38:LYS:O	1:A:42:ILE:HG13	2.14	0.47
1:B:83:GLU:OE2	1:B:86:THR:HG21	2.14	0.47
1:A:73:MET:HE2	1:A:73:MET:HB3	1.67	0.47
1:B:229:PHE:CE2	2:B:263:HOH:O	2.56	0.47
1:A:100:LEU:HD22	1:A:103:PHE:HE2	1.80	0.47
1:A:8:LEU:HD13	1:A:177:LEU:HG	1.95	0.47
1:B:220:SER:O	1:B:224:LYS:HG3	2.15	0.47
1:B:95:ARG:HG3	2:B:541:HOH:O	2.15	0.46
1:A:216:GLU:HG2	2:A:388:HOH:O	2.16	0.46
1:B:46:ASN:ND2	2:B:539:HOH:O	2.47	0.46
1:B:17:ARG:O	1:B:20:LYS:HG2	2.15	0.46
1:B:72:LEU:N	2:B:562:HOH:O	2.48	0.46
1:A:226:VAL:O	1:A:227:ARG:C	2.58	0.45
1:B:32:LYS:HD3	2:B:404:HOH:O	2.15	0.45
1:B:224:LYS:NZ	2:B:484:HOH:O	2.49	0.45
1:A:105:LYS:NZ	1:A:108:ARG:NH2	2.64	0.45
1:A:112:LEU:HD11	2:A:261:HOH:O	2.16	0.45
1:B:81:LYS:HD2	2:B:417:HOH:O	2.15	0.45
1:B:90:ASN:OD1	2:B:507:HOH:O	2.21	0.45
1:A:92:ILE:N	1:A:121:PHE:HD2	2.15	0.45
1:A:170:ARG:O	1:A:171:ASN:HB2	2.17	0.45
1:A:25:ALA:HB3	2:A:316:HOH:O	2.16	0.45
1:B:26:ARG:HD3	2:B:524:HOH:O	2.17	0.45
1:A:24:LYS:HD3	1:A:24:LYS:HA	1.71	0.45
1:A:227:ARG:HD2	2:A:310:HOH:O	2.16	0.45
1:A:105:LYS:HD3	1:A:108:ARG:NH2	2.20	0.44
1:B:122:LYS:HD2	2:B:340:HOH:O	2.17	0.44
1:B:226:VAL:HA	1:B:229:PHE:CE2	2.52	0.44
1:A:77:LYS:NZ	1:A:81:LYS:NZ	2.66	0.44
1:B:97:ILE:HG22	1:B:168:ASP:HB3	2.00	0.44
1:A:169:THR:O	1:A:170:ARG:C	2.60	0.44
1:A:118:THR:C	1:A:120:HIS:N	2.75	0.43
1:B:116:ASN:O	1:B:119:ARG:HG2	2.19	0.43
1:A:36:THR:O	1:A:37:LEU:C	2.58	0.43
1:A:6:LYS:NZ	1:A:50:GLU:OE2	2.39	0.43
1:B:22:LYS:CB	1:B:24:LYS:HG3	2.50	0.42
1:A:84:ARG:HH11	1:A:84:ARG:HG2	1.85	0.42
1:B:26:ARG:NH1	2:B:421:HOH:O	2.52	0.42
1:B:83:GLU:O	1:B:84:ARG:C	2.62	0.42
1:B:26:ARG:NH2	2:B:421:HOH:O	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:LYS:HE2	1:A:77:LYS:HB2	1.91	0.42
1:A:7:HIS:HB2	1:A:173:PRO:HG2	2.01	0.42
1:A:225:ARG:O	1:A:227:ARG:N	2.53	0.42
1:A:83:GLU:O	1:A:84:ARG:C	2.63	0.42
1:A:27:ALA:HA	2:A:484:HOH:O	2.20	0.41
1:B:72:LEU:C	2:B:561:HOH:O	2.62	0.41
1:A:100:LEU:HD22	1:A:103:PHE:CE2	2.54	0.41
2:A:384:HOH:O	1:B:20:LYS:NZ	2.53	0.41
1:A:71:PHE:O	1:A:74:LYS:HB3	2.20	0.41
1:B:75:MET:HE3	2:B:316:HOH:O	2.21	0.41
1:A:143:LYS:NZ	2:A:368:HOH:O	2.54	0.41
1:B:43:TRP:O	1:B:47:HIS:CD2	2.74	0.41
1:B:20:LYS:HD3	2:B:304:HOH:O	2.20	0.40
1:B:26:ARG:CZ	2:B:421:HOH:O	2.70	0.40
1:A:83:GLU:HG2	1:A:86:THR:CG2	2.51	0.40
1:A:214:ASP:HA	1:A:217:ASN:HD22	1.86	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	211/245 (86%)	201 (95%)	8 (4%)	2 (1%)	14	4
1	B	197/245 (80%)	194 (98%)	3 (2%)	0	100	100
All	All	408/490 (83%)	395 (97%)	11 (3%)	2 (0%)	24	13

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	ARG
1	A	226	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	196/223 (88%)	186 (95%)	10 (5%)	21	7
1	B	184/223 (82%)	181 (98%)	3 (2%)	55	43
All	All	380/446 (85%)	367 (97%)	13 (3%)	32	16

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	32	LYS
1	A	41	THR
1	A	57	PHE
1	A	70	ASP
1	A	76	LEU
1	A	77	LYS
1	A	155	LEU
1	A	177	LEU
1	A	212	PRO
1	B	4	THR
1	B	74	LYS
1	B	212	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	46	ASN
1	A	47	HIS
1	A	91	ASN
1	A	116	ASN
1	A	126	GLN
1	A	136	ASN
1	A	217	ASN
1	B	46	ASN

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Mol	Chain	Res	Type
1	B	47	HIS
1	B	91	ASN
1	B	120	HIS
1	B	136	ASN
1	B	171	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	215/245 (87%)	1.71	78 (36%) 1 0	19, 41, 74, 93	0
1	B	203/245 (82%)	0.64	26 (12%) 7 8	13, 24, 55, 79	0
All	All	418/490 (85%)	1.19	104 (24%) 2 1	13, 32, 70, 93	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	28	TYR	6.1
1	B	72	LEU	5.7
1	A	121	PHE	5.4
1	A	69	VAL	5.4
1	A	29	GLY	5.1
1	A	27	ALA	5.1
1	A	21	LEU	4.9
1	A	57	PHE	4.8
1	A	155	LEU	4.7
1	B	76	LEU	4.6
1	A	88	LEU	4.4
1	A	226	VAL	4.4
1	A	87	TYR	4.3
1	A	112	LEU	4.3
1	A	103	PHE	4.2
1	A	14	GLY	4.2
1	A	120	HIS	4.2
1	A	18	TRP	4.2
1	A	71	PHE	4.1
1	A	19	ALA	4.1
1	A	25	ALA	4.1
1	A	86	THR	3.9
1	B	74	LYS	3.7
1	A	79	TYR	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	118	THR	3.7
1	B	107	LEU	3.6
1	B	230	GLY	3.5
1	A	20	LYS	3.5
1	A	31	LYS	3.5
1	B	105	LYS	3.4
1	B	103	PHE	3.4
1	A	24	LYS	3.3
1	B	149	PRO	3.3
1	A	154	SER	3.3
1	A	42	ILE	3.2
1	A	26	ARG	3.2
1	B	75	MET	3.2
1	A	56	ALA	3.2
1	B	231	GLU	3.2
1	A	43	TRP	3.2
1	A	124	PHE	3.2
1	A	47	HIS	3.2
1	A	85	SER	3.1
1	A	32	LYS	3.1
1	A	3	SER	3.1
1	A	23	ASN	3.1
1	B	77	LYS	3.1
1	A	30	HIS	3.1
1	A	227	ARG	3.0
1	A	92	ILE	3.0
1	A	80	LEU	3.0
1	A	100	LEU	3.0
1	A	116	ASN	3.0
1	A	33	GLY	3.0
1	A	72	LEU	3.0
1	A	76	LEU	3.0
1	B	73	MET	2.9
1	A	89	ASP	2.8
1	A	205	ILE	2.8
1	B	229	PHE	2.8
1	A	35	LYS	2.8
1	A	153	ILE	2.7
1	A	109	ASP	2.7
1	B	228	LYS	2.7
1	A	114	LEU	2.6
1	A	22	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	51	CYS	2.6
1	A	207	TRP	2.6
1	A	4	THR	2.6
1	A	119	ARG	2.6
1	B	80	LEU	2.6
1	B	159	LEU	2.6
1	B	85	SER	2.6
1	B	104	SER	2.6
1	B	148	SER	2.5
1	A	206	LEU	2.5
1	B	227	ARG	2.5
1	B	226	VAL	2.5
1	A	223	TYR	2.5
1	B	102	GLY	2.5
1	A	91	ASN	2.5
1	A	49	LEU	2.4
1	B	4	THR	2.4
1	A	113	GLN	2.3
1	A	84	ARG	2.3
1	B	109	ASP	2.3
1	A	90	ASN	2.3
1	A	151	SER	2.3
1	A	34	VAL	2.3
1	A	165	ASN	2.3
1	A	5	LEU	2.3
1	A	110	THR	2.2
1	A	106	GLU	2.2
1	B	161	ASN	2.2
1	A	208	PRO	2.2
1	A	152	ASN	2.1
1	A	111	ILE	2.1
1	A	122	LYS	2.1
1	A	104	SER	2.1
1	A	46	ASN	2.1
1	A	102	GLY	2.1
1	A	82	ASP	2.0
1	B	101	GLU	2.0
1	A	81	LYS	2.0

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

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.