



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2021 – 04:08 PM JST

PDB ID : 7DAM
Title : Adenosine triphosphate phosphoribosyltransferase from *Vibrio cholerae*
Authors : Pal, R.K.; Gourinath, S.; Biswal, B.K.
Deposited on : 2020-10-16
Resolution : 2.70 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.23.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

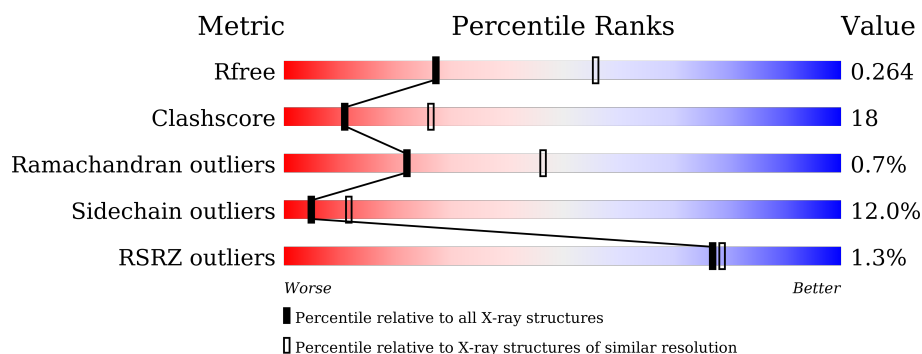
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	310	
1	B	310	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3764 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 ATP phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	1	0
			1973	1244	339	380	10			
1	B	272	Total	C	N	O	S	0	0	0
			1781	1119	302	353	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP C3LU29
A	-4	HIS	-	expression tag	UNP C3LU29
A	-3	HIS	-	expression tag	UNP C3LU29
A	-2	HIS	-	expression tag	UNP C3LU29
A	-1	HIS	-	expression tag	UNP C3LU29
A	0	HIS	-	expression tag	UNP C3LU29
B	-5	HIS	-	expression tag	UNP C3LU29
B	-4	HIS	-	expression tag	UNP C3LU29
B	-3	HIS	-	expression tag	UNP C3LU29
B	-2	HIS	-	expression tag	UNP C3LU29
B	-1	HIS	-	expression tag	UNP C3LU29
B	0	HIS	-	expression tag	UNP C3LU29

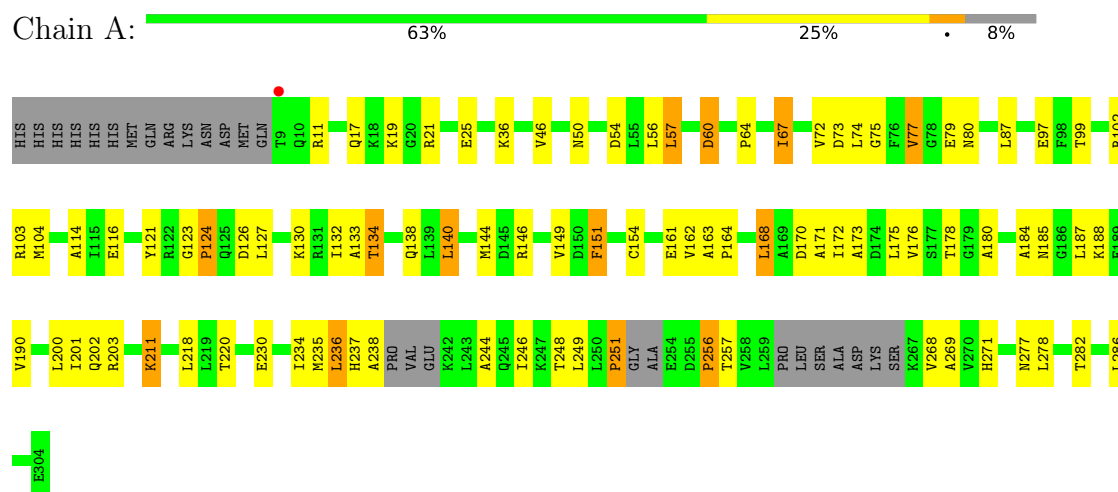
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	7	Total	O	0	0
			7	7		
2	B	3	Total	O	0	0
			3	3		

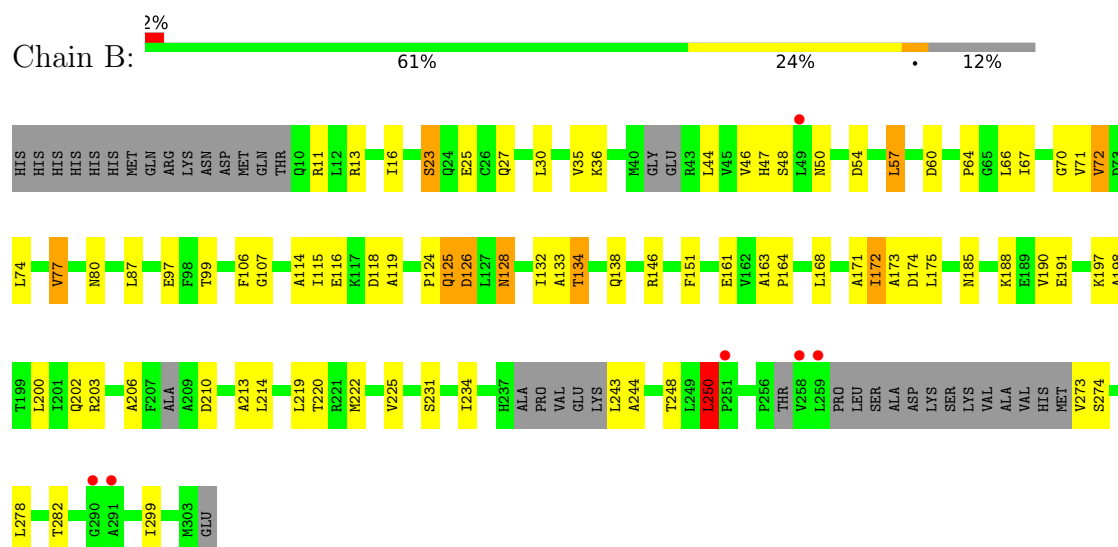
3 Residue-property plots

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: ATP phosphoribosyltransferase



• Molecule 1: ATP phosphoribosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	117.51Å 117.51Å 121.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.67 – 2.70 36.67 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (36.67-2.70) 99.7 (36.67-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.210 , 0.265 0.211 , 0.264	Depositor DCC
R_{free} test set	1745 reflections (10.26%)	wwPDB-VP
Wilson B-factor (Å ²)	69.9	Xtriage
Anisotropy	0.235	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.19 , 17.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.448 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.543 for H, K, L 0.457 for K, H, -L	Depositor
Outliers	0 of 17005 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3764	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.92% 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	0.69	0/1995	0.81	0/2713
1	B	0.71	0/1788	0.80	0/2447
All	All	0.70	0/3783	0.80	0/5160

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	172	ILE	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	1973	0	1817	81	0
1	B	1781	0	1578	53	0
2	A	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	3	0	0	0	0
All	All	3764	0	3395	130	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 18.

All (130) 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:282:THR:O	1:A:286:LEU:HD12	1.75	0.85
1:B:116:GLU:O	1:B:119:ALA:HB3	1.77	0.83
1:A:134:THR:CG2	1:A:173:ALA:HB3	2.20	0.72
1:B:203:ARG:CB	1:B:206:ALA:HB3	2.20	0.72
1:A:11:ARG:O	1:A:211:LYS:NZ	2.23	0.71
1:A:73:ASP:HA	1:A:203:ARG:HG3	1.73	0.71
1:A:236:LEU:O	1:A:269:ALA:HB1	1.90	0.70
1:B:134:THR:CG2	1:B:173:ALA:HB3	2.22	0.70
1:A:151:PHE:H	1:A:151:PHE:HD1	1.38	0.70
1:A:237:HIS:HA	1:A:269:ALA:HB2	1.73	0.69
1:B:11:ARG:NH2	1:B:54:ASP:OD1	2.28	0.66
1:B:161:GLU:CB	1:B:185:ASN:ND2	2.59	0.66
1:A:230:GLU:O	1:A:277:ASN:N	2.29	0.66
1:A:21:ARG:O	1:A:25:GLU:HG3	1.97	0.63
1:B:30:LEU:O	1:B:35:VAL:HG12	1.99	0.63
1:A:163:ALA:HB3	1:A:164:PRO:HD3	1.82	0.61
1:A:180:ALA:O	1:A:184:ALA:N	2.33	0.60
1:A:246:ILE:O	1:A:249:LEU:HB2	2.00	0.60
1:B:163:ALA:HB3	1:B:164:PRO:HD3	1.84	0.60
1:B:80:ASN:HD22	1:B:175:LEU:HD11	1.67	0.59
1:A:235:MET:HG2	1:A:271[A]:HIS:ND1	2.18	0.58
1:A:56:LEU:HD12	1:A:72:VAL:HG11	1.86	0.57
1:A:257:THR:HG21	1:A:271[B]:HIS:HB3	1.86	0.57
1:A:73:ASP:HA	1:A:203:ARG:CG	2.35	0.57
1:A:257:THR:HG21	1:A:271[A]:HIS:H	1.70	0.56
1:A:235:MET:CG	1:A:271[A]:HIS:ND1	2.69	0.55
1:A:236:LEU:O	1:A:236:LEU:HD12	2.07	0.55
1:A:257:THR:HG21	1:A:271[B]:HIS:H	1.71	0.54
1:A:235:MET:SD	1:A:271[A]:HIS:ND1	2.81	0.54
1:A:57:LEU:HD23	1:A:57:LEU:N	2.23	0.54
1:A:168:LEU:HD23	1:B:71:VAL:HG11	1.89	0.53
1:B:74:LEU:HD21	1:B:200:LEU:HD11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ASP:HA	1:A:203:ARG:CD	2.38	0.53
1:A:236:LEU:O	1:A:269:ALA:CB	2.56	0.53
1:B:106:PHE:O	1:B:198:ALA:HB3	2.09	0.52
1:B:80:ASN:ND2	1:B:175:LEU:HD11	2.25	0.52
1:A:102:ARG:HG2	1:A:104:MET:HE2	1.91	0.52
1:B:46:VAL:O	1:B:54:ASP:HA	2.09	0.52
1:A:244:ALA:O	1:A:248:THR:HG23	2.11	0.51
1:A:46:VAL:O	1:A:54:ASP:HA	2.10	0.51
1:B:116:GLU:O	1:B:119:ALA:CB	2.55	0.51
1:B:74:LEU:CD2	1:B:200:LEU:HD11	2.40	0.51
1:B:273:VAL:HG23	1:B:299:ILE:HD11	1.91	0.51
1:A:56:LEU:HD12	1:A:72:VAL:CG1	2.40	0.51
1:A:161:GLU:CB	1:A:185:ASN:OD1	2.59	0.50
1:B:57:LEU:HD23	1:B:57:LEU:N	2.26	0.50
1:A:162:VAL:HG21	1:B:44:LEU:HD12	1.93	0.50
1:A:237:HIS:HA	1:A:269:ALA:CB	2.42	0.50
1:A:151:PHE:CD1	1:A:151:PHE:N	2.77	0.50
1:A:154:CYS:SG	1:A:168:LEU:HD23	2.53	0.49
1:A:80:ASN:HD22	1:A:175:LEU:HD11	1.77	0.49
1:B:35:VAL:O	1:B:35:VAL:HG13	2.12	0.49
1:A:168:LEU:CD2	1:B:71:VAL:HG11	2.42	0.49
1:A:132:ILE:HA	1:A:171:ALA:O	2.13	0.48
1:A:102:ARG:HB3	1:A:104:MET:HE1	1.96	0.48
1:A:123:GLY:O	1:A:124:PRO:CB	2.62	0.48
1:A:127:LEU:HD23	1:A:127:LEU:N	2.28	0.48
1:B:234:ILE:N	1:B:234:ILE:HD12	2.28	0.48
1:A:74:LEU:HD21	1:A:200:LEU:HD11	1.96	0.48
1:A:134:THR:HG23	1:A:173:ALA:HB3	1.92	0.48
1:B:99:THR:O	1:B:202:GLN:N	2.47	0.48
1:A:123:GLY:HA2	1:A:190:VAL:CG1	2.44	0.47
1:A:278:LEU:C	1:A:278:LEU:HD12	2.35	0.47
1:B:11:ARG:HH22	1:B:47:HIS:HB3	1.79	0.47
1:B:134:THR:HG23	1:B:173:ALA:HB3	1.95	0.47
1:A:36:LYS:CB	1:A:50:ASN:CB	2.92	0.47
1:A:67:ILE:HD11	1:A:74:LEU:C	2.35	0.47
1:A:74:LEU:CD2	1:A:200:LEU:HD11	2.44	0.47
1:B:13:ARG:HB3	1:B:72:VAL:HG13	1.95	0.47
1:B:70:GLY:CA	1:B:206:ALA:HB2	2.45	0.47
1:B:210:ASP:O	1:B:213:ALA:HB3	2.14	0.47
1:B:70:GLY:HA2	1:B:206:ALA:HB2	1.96	0.47
1:B:132:ILE:HA	1:B:171:ALA:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:SER:HA	1:B:274:SER:O	2.15	0.47
1:B:243:LEU:C	1:B:243:LEU:HD23	2.36	0.46
1:A:67:ILE:HG12	1:A:72:VAL:HG23	1.96	0.46
1:B:222:MET:O	1:B:225:VAL:HG22	2.15	0.46
1:A:60:ASP:N	1:A:60:ASP:OD1	2.47	0.46
1:A:99:THR:O	1:A:202:GLN:N	2.48	0.46
1:A:234:ILE:HD12	1:A:234:ILE:N	2.30	0.46
1:B:124:PRO:C	1:B:126:ASP:H	2.19	0.46
1:B:190:VAL:HG12	1:B:191:GLU:HG2	1.98	0.46
1:A:56:LEU:C	1:A:57:LEU:HD23	2.36	0.45
1:B:87:LEU:HB2	1:B:138:GLN:HG2	1.98	0.45
1:A:132:ILE:CG2	1:A:140:LEU:HD11	2.46	0.45
1:B:67:ILE:HG23	1:B:203:ARG:HA	1.97	0.45
1:B:250:LEU:HD22	1:B:250:LEU:O	2.17	0.45
1:B:172:ILE:HD12	1:B:174:ASP:HB2	1.99	0.44
1:A:67:ILE:HD12	1:A:201:ILE:HG13	2.00	0.44
1:B:115:ILE:HG22	1:B:119:ALA:HB3	1.98	0.44
1:B:35:VAL:O	1:B:35:VAL:CG1	2.65	0.44
1:A:73:ASP:HA	1:A:203:ARG:HD3	1.99	0.43
1:A:87:LEU:HB2	1:A:138:GLN:HG2	1.99	0.43
1:B:11:ARG:NH1	1:B:48:SER:O	2.51	0.43
1:B:36:LYS:H	1:B:50:ASN:CB	2.31	0.43
1:A:56:LEU:CD1	1:A:72:VAL:CG1	2.96	0.43
1:A:79:GLU:CD	1:A:103:ARG:HE	2.22	0.43
1:A:175:LEU:HD23	1:A:175:LEU:HA	1.89	0.43
1:A:238:ALA:HB3	1:A:268:VAL:O	2.19	0.43
1:A:235:MET:SD	1:A:271[A]:HIS:CE1	3.12	0.43
1:A:36:LYS:H	1:A:50:ASN:CB	2.32	0.43
1:A:80:ASN:ND2	1:A:175:LEU:HD11	2.33	0.43
1:B:125:GLN:HA	1:B:128:ASN:HB2	2.00	0.43
1:A:67:ILE:HD11	1:A:75:GLY:N	2.34	0.42
1:B:64:PRO:HD3	1:B:77:VAL:HG11	2.02	0.42
1:A:257:THR:CG2	1:A:271[A]:HIS:H	2.32	0.42
1:A:116:GLU:HA	1:A:187:LEU:HD23	2.02	0.42
1:A:64:PRO:HD3	1:A:77:VAL:HG11	2.02	0.42
1:A:257:THR:CG2	1:A:271[B]:HIS:H	2.33	0.42
1:A:17:GLN:HE21	1:A:17:GLN:HB3	1.59	0.42
1:A:17:GLN:NE2	1:A:19:LYS:O	2.51	0.41
1:A:130:LYS:HD2	1:A:170:ASP:CG	2.40	0.41
1:A:130:LYS:HB3	1:A:170:ASP:HB2	2.00	0.41
1:B:114:ALA:HA	1:B:188:LYS:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:LEU:C	1:B:35:VAL:HG12	2.39	0.41
1:B:107:GLY:O	1:B:197:LYS:HA	2.21	0.41
1:B:60:ASP:N	1:B:60:ASP:OD1	2.53	0.41
1:A:114:ALA:HA	1:A:188:LYS:O	2.21	0.41
1:B:133:ALA:O	1:B:172:ILE:HA	2.21	0.41
1:A:237:HIS:O	1:A:237:HIS:CG	2.74	0.41
1:A:248:THR:O	1:A:251:PRO:HG3	2.20	0.41
1:A:132:ILE:HG21	1:A:140:LEU:HD11	2.02	0.41
1:A:133:ALA:O	1:A:172:ILE:HA	2.21	0.41
1:A:218:LEU:HD23	1:A:218:LEU:HA	1.92	0.41
1:B:23:SER:OG	1:B:27:GLN:NE2	2.54	0.41
1:B:244:ALA:O	1:B:248:THR:N	2.50	0.41
1:A:126:ASP:O	1:A:149:VAL:CG1	2.69	0.40
1:A:244:ALA:O	1:A:248:THR:N	2.50	0.40
1:A:168:LEU:HD21	1:B:66:LEU:HD22	2.02	0.40
1:B:74:LEU:HD21	1:B:200:LEU:CD1	2.51	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	277/310 (89%)	255 (92%)	19 (7%)	3 (1%)	14	34
1	B	260/310 (84%)	243 (94%)	16 (6%)	1 (0%)	34	60
All	All	537/620 (87%)	498 (93%)	35 (6%)	4 (1%)	22	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	PRO
1	B	250	LEU

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Mol	Chain	Res	Type
1	A	256	PRO
1	A	121	TYR

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	176/263 (67%)	158 (90%)	18 (10%)	7	17
1	B	149/263 (57%)	128 (86%)	21 (14%)	3	8
All	All	325/526 (62%)	286 (88%)	39 (12%)	5	11

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	LEU
1	A	60	ASP
1	A	67	ILE
1	A	77	VAL
1	A	97	GLU
1	A	134	THR
1	A	140	LEU
1	A	144	MET
1	A	146	ARG
1	A	151	PHE
1	A	168	LEU
1	A	176	VAL
1	A	178	THR
1	A	211	LYS
1	A	220	THR
1	A	236	LEU
1	A	251	PRO
1	A	256	PRO
1	B	16	ILE
1	B	23	SER
1	B	25	GLU

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Mol	Chain	Res	Type
1	B	57	LEU
1	B	72	VAL
1	B	77	VAL
1	B	97	GLU
1	B	118	ASP
1	B	125	GLN
1	B	126	ASP
1	B	128	ASN
1	B	134	THR
1	B	146	ARG
1	B	151	PHE
1	B	168	LEU
1	B	214	LEU
1	B	219	LEU
1	B	220	THR
1	B	250	LEU
1	B	278	LEU
1	B	282	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	24	GLN
1	A	227	GLN
1	A	237	HIS
1	B	27	GLN
1	B	125	GLN
1	B	185	ASN
1	B	227	GLN

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 [i](#)

There are no monosaccharides 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	284/310 (91%)	-0.50	1 (0%) 92 93	40, 67, 98, 123	0
1	B	272/310 (87%)	-0.46	6 (2%) 62 63	42, 69, 98, 115	0
All	All	556/620 (89%)	-0.48	7 (1%) 77 78	40, 68, 98, 123	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	291	ALA	3.2
1	B	259	LEU	3.0
1	B	251	PRO	2.6
1	B	258	VAL	2.5
1	B	290	GLY	2.3
1	B	49	LEU	2.2
1	A	9	THR	2.2

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 monosaccharides 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.