



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 28, 2024 – 07:24 pm GMT

PDB ID : 2WE7
Title : Crystal structure of Mycobacterium tuberculosis Rv0376c homologue from Mycobacterium smegmatis
Authors : Cho, H.J.; Kang, B.S.
Deposited on : 2009-03-29
Resolution : 2.90 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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

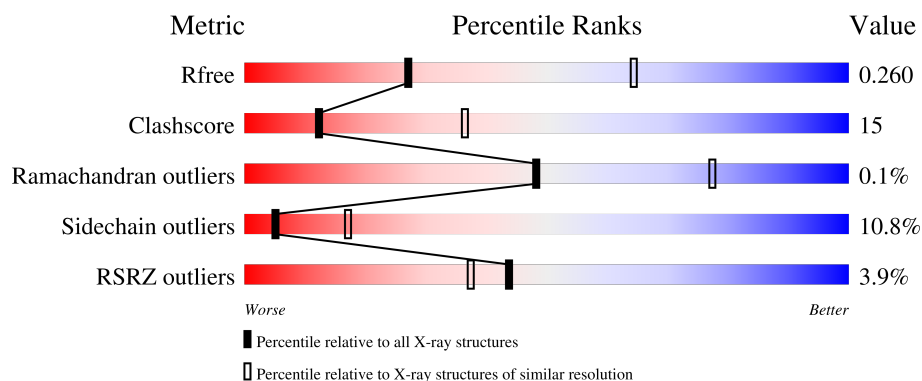
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.90 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}	164625	2335 (2.90-2.90)
Clashscore	180529	2564 (2.90-2.90)
Ramachandran outliers	177936	2514 (2.90-2.90)
Sidechain outliers	177891	2516 (2.90-2.90)
RSRZ outliers	164620	2337 (2.90-2.90)

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	386	
1	B	386	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5046 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 Xanthine dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	Se	0	3	0
			2562	1604	462	489	3	4			
1	B	335	Total	C	N	O	S	Se	0	0	0
			2463	1545	436	475	3	4			

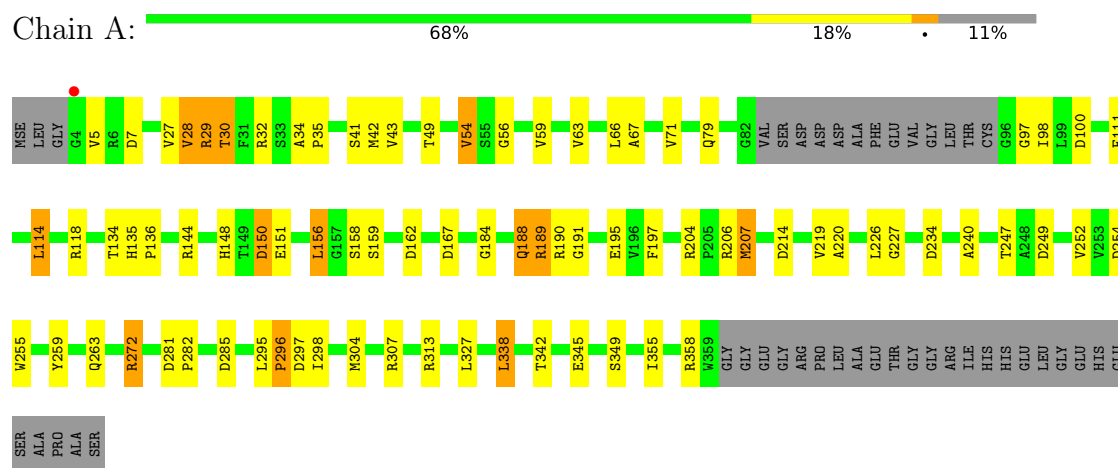
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	20	Total	O	0	0
			20	20		
2	B	1	Total	O	0	0
			1	1		

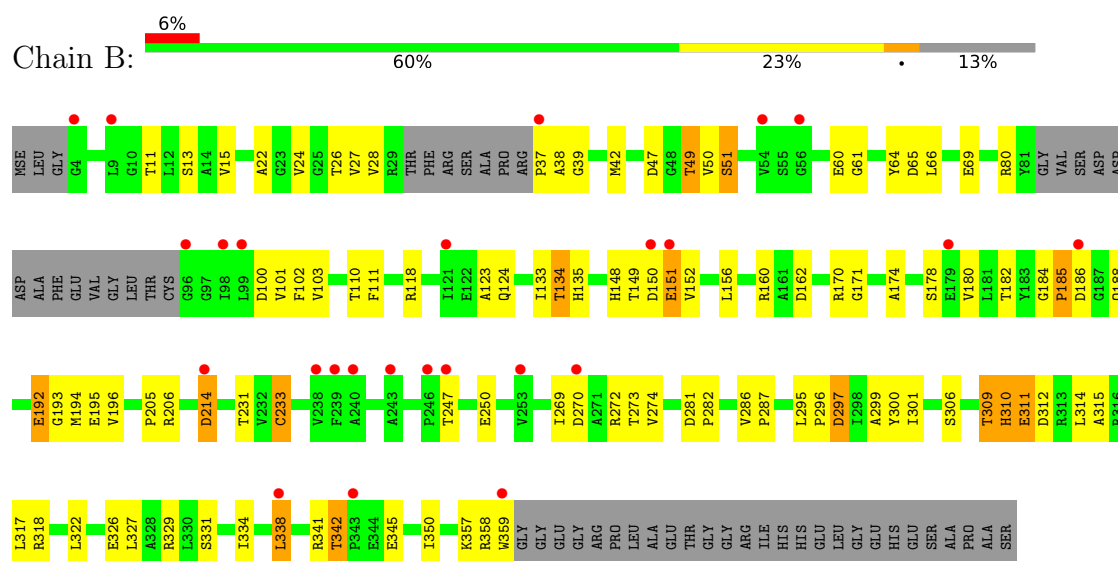
3 Residue-property plots

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: Xanthine dehydrogenase



- Molecule 1: Xanthine dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	101.04Å 101.04Å 316.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.29 – 2.90 38.29 – 2.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (38.29-2.90) 98.5 (38.29-2.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	45.39 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.195 , 0.252 0.207 , 0.260	Depositor DCC
R_{free} test set	1135 reflections (5.21%)	wwPDB-VP
Wilson B-factor (Å ²)	54.5	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 50.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5046	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.31% 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	0.95	2/2614 (0.1%)	0.78	2/3557 (0.1%)
1	B	0.63	1/2503 (0.0%)	0.67	0/3408
All	All	0.81	3/5117 (0.1%)	0.73	2/6965 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	252	VAL	CB-CG1	-5.52	1.41	1.52
1	B	233	CYS	CB-SG	-5.38	1.73	1.81
1	A	252	VAL	CB-CG2	-5.12	1.42	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	191	GLY	N-CA-C	-5.29	99.88	113.10
1	A	296	PRO	N-CA-C	-5.26	98.42	112.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2562	0	2555	51	0
1	B	2463	0	2437	101	0
2	A	20	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
All	All	5046	0	4992	152	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 15.

All (152) 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:114:LEU:O	1:A:114:LEU:HD12	1.34	1.21
1:A:189:ARG:CG	1:A:189:ARG:HH11	1.56	1.17
1:B:27:VAL:HG23	1:B:38:ALA:HA	1.21	1.14
1:A:189:ARG:HH11	1:A:189:ARG:HG3	1.16	1.05
1:B:342:THR:HG22	1:B:345:GLU:HG3	1.37	1.04
1:B:28:VAL:CG1	1:B:180:VAL:CG2	2.40	0.99
1:B:47:ASP:OD1	1:B:49:THR:HG23	1.66	0.95
1:B:27:VAL:CG2	1:B:38:ALA:HA	1.96	0.94
1:B:28:VAL:HG12	1:B:180:VAL:HG21	1.49	0.93
1:A:134:THR:HB	1:A:195:GLU:HG2	1.51	0.90
1:B:28:VAL:CG1	1:B:180:VAL:HG21	2.03	0.87
1:A:114:LEU:HD12	1:A:114:LEU:C	1.87	0.84
1:B:342:THR:HG22	1:B:345:GLU:CG	2.07	0.84
1:A:189:ARG:O	1:A:190:ARG:HB2	1.74	0.83
1:A:189:ARG:CG	1:A:189:ARG:NH1	2.30	0.82
1:B:27:VAL:HG23	1:B:27:VAL:O	1.79	0.82
1:A:189:ARG:HG3	1:A:189:ARG:NH1	1.87	0.81
1:B:28:VAL:HG12	1:B:180:VAL:CG2	2.05	0.81
1:B:310:HIS:CE1	1:B:314:LEU:HD11	2.16	0.80
1:B:27:VAL:CG2	1:B:27:VAL:O	2.32	0.78
1:A:189:ARG:HH11	1:A:189:ARG:HG2	1.45	0.78
1:A:204[B]:ARG:HG3	1:A:204[B]:ARG:HH11	1.48	0.78
1:B:28:VAL:CG1	1:B:180:VAL:HG22	2.13	0.77
1:A:30:THR:CG2	1:A:34:ALA:HB3	2.13	0.77
1:B:310:HIS:CE1	1:B:314:LEU:HG	2.20	0.76
1:B:269:ILE:O	1:B:269:ILE:HG13	1.83	0.76
1:B:27:VAL:HG21	1:B:37:PRO:O	1.85	0.76
1:A:30:THR:HG22	1:A:34:ALA:HB3	1.66	0.76
1:B:50:VAL:HG12	1:B:51:SER:N	2.00	0.74
1:B:310:HIS:CE1	1:B:314:LEU:CG	2.72	0.73
1:B:47:ASP:OD1	1:B:49:THR:CG2	2.36	0.72
1:B:269:ILE:O	1:B:269:ILE:CG1	2.38	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:310:HIS:CE1	1:B:314:LEU:CD1	2.72	0.71
1:A:59:VAL:O	1:A:63:VAL:HG23	1.90	0.71
1:B:182:THR:HG22	1:B:195:GLU:HG2	1.71	0.70
1:B:342:THR:CG2	1:B:345:GLU:HG3	2.18	0.70
1:A:35:PRO:HG3	1:A:56:GLY:H	1.57	0.70
1:B:28:VAL:HG13	1:B:180:VAL:HG22	1.74	0.69
1:B:192:GLU:HG2	1:B:193:GLY:N	2.07	0.69
1:B:42:MSE:HE2	1:B:50:VAL:HG11	1.73	0.69
1:A:207:MSE:HE1	1:A:219:VAL:O	1.91	0.69
1:A:184:GLY:HA3	1:A:188:GLN:HG2	1.77	0.67
1:A:358:ARG:HB3	2:A:2020:HOH:O	1.94	0.66
1:A:29:ARG:NH1	1:A:100:ASP:OD2	2.29	0.66
1:B:151:GLU:OE1	1:B:151:GLU:C	2.35	0.66
1:B:184:GLY:O	1:B:186:ASP:N	2.30	0.65
1:B:134:THR:O	1:B:194:MSE:HA	1.96	0.65
1:A:32:ARG:HB2	1:A:97:GLY:HA2	1.79	0.65
1:B:151:GLU:OE1	1:B:152:VAL:N	2.30	0.65
1:B:286:VAL:HB	1:B:287:PRO:HD3	1.79	0.65
1:B:341:ARG:N	1:B:345:GLU:OE1	2.31	0.63
1:B:50:VAL:CG1	1:B:51:SER:N	2.62	0.62
1:B:27:VAL:HG22	1:B:39:GLY:H	1.65	0.62
1:B:160:ARG:NH2	1:B:188:GLN:HB3	2.15	0.60
1:A:5:VAL:HG12	1:A:41:SER:OG	2.03	0.59
1:B:102:PHE:CD1	1:B:133:ILE:HG22	2.36	0.59
1:B:50:VAL:CG2	1:B:64:TYR:CE2	2.86	0.59
1:B:270:ASP:OD1	1:B:272:ARG:HG2	2.03	0.58
1:A:234:ASP:HB3	1:A:240:ALA:HB2	1.84	0.58
1:B:331:SER:OG	1:B:357:LYS:HE3	2.03	0.57
1:B:28:VAL:HG13	1:B:180:VAL:CG2	2.29	0.57
1:B:13:SER:OG	1:B:118:ARG:NH1	2.38	0.56
1:B:27:VAL:CG2	1:B:38:ALA:CA	2.76	0.56
1:A:67:ALA:O	1:A:71:VAL:HG23	2.06	0.55
1:B:50:VAL:CG2	1:B:64:TYR:HE2	2.19	0.55
1:B:310:HIS:HE1	1:B:314:LEU:HD21	1.70	0.55
1:A:295:LEU:O	1:A:296:PRO:C	2.43	0.55
1:B:15:VAL:HG11	1:B:22:ALA:HB2	1.88	0.55
1:B:26:THR:O	1:B:101:VAL:HA	2.06	0.55
1:B:270:ASP:OD1	1:B:272:ARG:N	2.28	0.55
1:A:28:VAL:HG22	1:A:197:PHE:CG	2.43	0.54
1:A:156:LEU:HB2	1:A:162:ASP:OD1	2.07	0.54
1:A:226:LEU:HD21	1:A:355:ILE:HD11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:ARG:HA	1:B:299:ALA:HB2	1.90	0.53
1:B:295:LEU:O	1:B:296:PRO:C	2.45	0.53
1:B:50:VAL:HG21	1:B:64:TYR:CD2	2.44	0.53
1:B:273:THR:HG22	1:B:274:VAL:N	2.23	0.53
1:B:50:VAL:HG23	1:B:64:TYR:HE2	1.74	0.52
1:A:35:PRO:CG	1:A:56:GLY:H	2.20	0.52
1:B:27:VAL:HG22	1:B:39:GLY:N	2.24	0.52
1:A:27:VAL:HG23	1:A:54:VAL:HG22	1.91	0.52
1:A:134:THR:CB	1:A:195:GLU:HG2	2.33	0.52
1:B:27:VAL:HG23	1:B:38:ALA:CA	2.15	0.51
1:A:204[B]:ARG:HG3	1:A:204[B]:ARG:NH1	2.22	0.51
1:B:315:ALA:HA	1:B:318:ARG:HD2	1.91	0.51
1:B:151:GLU:N	1:B:170:ARG:HH21	2.09	0.50
1:B:273:THR:CG2	1:B:274:VAL:N	2.74	0.50
1:B:342:THR:HG22	1:B:345:GLU:CD	2.32	0.50
1:A:272:ARG:HG3	1:A:272:ARG:HH11	1.76	0.50
1:B:27:VAL:CG2	1:B:37:PRO:O	2.59	0.50
1:B:306:SER:OG	1:B:309:THR:CG2	2.59	0.50
1:B:186:ASP:OD2	1:B:186:ASP:O	2.30	0.50
1:B:309:THR:O	1:B:309:THR:OG1	2.30	0.50
1:B:214:ASP:OD1	1:B:214:ASP:N	2.37	0.49
1:B:306:SER:O	1:B:310:HIS:N	2.43	0.49
1:B:310:HIS:HE1	1:B:314:LEU:CG	2.24	0.49
1:B:148:HIS:O	1:B:170:ARG:NE	2.44	0.49
1:B:150:ASP:OD2	1:B:150:ASP:O	2.30	0.49
1:B:273:THR:O	1:B:299:ALA:HB3	2.13	0.48
1:A:148:HIS:HB2	1:A:151:GLU:HG2	1.95	0.48
1:B:311:GLU:O	1:B:312:ASP:C	2.48	0.48
1:A:189:ARG:O	1:A:190:ARG:CB	2.45	0.48
1:B:306:SER:OG	1:B:309:THR:HG22	2.13	0.48
1:B:310:HIS:ND1	1:B:314:LEU:HG	2.28	0.48
1:B:274:VAL:HG22	1:B:300:TYR:HB3	1.95	0.47
1:A:254:ASP:OD1	1:A:255:TRP:N	2.47	0.47
1:B:151:GLU:H	1:B:170:ARG:HH21	1.63	0.47
1:A:184:GLY:HA3	1:A:188:GLN:CG	2.44	0.47
1:B:135:HIS:NE2	1:B:185:PRO:O	2.46	0.47
1:B:27:VAL:CG2	1:B:39:GLY:H	2.26	0.47
1:B:134:THR:OG1	1:B:195:GLU:HB2	2.15	0.46
1:B:334:ILE:HG12	1:B:350:ILE:HG12	1.96	0.46
1:B:24:VAL:O	1:B:103:VAL:HA	2.15	0.46
1:B:231:THR:HG23	1:B:250:GLU:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ARG:NH1	1:A:189:ARG:HG2	2.16	0.46
1:B:60:GLU:O	1:B:61:GLY:C	2.53	0.46
1:A:342:THR:HG23	1:A:345:GLU:OE2	2.15	0.46
1:A:135:HIS:CG	1:A:136:PRO:HD2	2.51	0.46
1:B:123:ALA:O	1:B:124:GLN:HB2	2.15	0.45
1:B:310:HIS:HE1	1:B:314:LEU:CD2	2.29	0.45
1:A:66:LEU:HD22	1:A:79:GLN:HG3	1.98	0.45
1:A:285:ASP:OD2	1:A:313:ARG:HD3	2.16	0.45
1:B:27:VAL:HA	1:B:100:ASP:O	2.16	0.45
1:B:297:ASP:N	1:B:297:ASP:OD1	2.46	0.45
1:B:358:ARG:O	1:B:358:ARG:HG2	2.18	0.44
1:A:156:LEU:HD12	1:A:156:LEU:HA	1.67	0.44
1:A:281:ASP:HA	1:A:282:PRO:HD3	1.86	0.44
1:A:304:MSE:HE3	2:A:2015:HOH:O	2.17	0.44
1:B:156:LEU:HB2	1:B:162:ASP:OD1	2.17	0.44
1:A:204[A]:ARG:HG2	1:A:227:GLY:O	2.18	0.43
1:B:322:LEU:HD21	1:B:326:GLU:HB2	1.99	0.43
1:A:259:TYR:O	1:A:263:GLN:HG2	2.18	0.43
1:A:338:LEU:HD22	1:A:349:SER:HB2	1.99	0.43
1:B:317:LEU:HD22	1:B:327:LEU:HD21	2.00	0.43
1:B:42:MSE:CE	1:B:50:VAL:HG11	2.44	0.43
1:B:326:GLU:HA	1:B:329:ARG:HH22	1.84	0.43
1:B:80:ARG:HG3	1:B:100:ASP:OD1	2.18	0.42
1:B:150:ASP:O	1:B:150:ASP:CG	2.55	0.42
1:B:171:GLY:O	1:B:174:ALA:HB3	2.20	0.42
1:A:298:ILE:HG21	1:A:298:ILE:HD13	1.85	0.41
1:B:338:LEU:HA	1:B:338:LEU:HD12	1.78	0.41
1:B:27:VAL:CG2	1:B:39:GLY:N	2.83	0.41
1:B:231:THR:HA	1:B:250:GLU:O	2.20	0.41
1:B:65:ASP:OD2	1:B:65:ASP:N	2.53	0.41
1:B:50:VAL:CG2	1:B:64:TYR:CD2	3.04	0.41
1:A:207:MSE:CE	1:A:220:ALA:HA	2.50	0.41
1:A:207:MSE:CE	1:A:219:VAL:HG12	2.51	0.41
1:A:114:LEU:O	1:A:114:LEU:CD1	2.30	0.41
1:A:150:ASP:OD1	1:A:150:ASP:N	2.52	0.40
1:B:184:GLY:O	1:B:185:PRO:C	2.59	0.40
1:B:205:PRO:HB3	1:B:272:ARG:HG3	2.03	0.40
1:B:281:ASP:HA	1:B:282:PRO:HD2	1.82	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	342/386 (89%)	335 (98%)	7 (2%)	0	100	100
1	B	329/386 (85%)	318 (97%)	10 (3%)	1 (0%)	37	66
All	All	671/772 (87%)	653 (97%)	17 (2%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	185	PRO

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	262/285 (92%)	231 (88%)	31 (12%)	4	13
1	B	250/285 (88%)	225 (90%)	25 (10%)	6	20
All	All	512/570 (90%)	456 (89%)	56 (11%)	5	17

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	28	VAL
1	A	29	ARG
1	A	30	THR
1	A	42	MSE

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Mol	Chain	Res	Type
1	A	43	VAL
1	A	49	THR
1	A	54	VAL
1	A	98	ILE
1	A	111	PHE
1	A	114	LEU
1	A	118	ARG
1	A	144	ARG
1	A	150	ASP
1	A	156	LEU
1	A	158	SER
1	A	159	SER
1	A	167	ASP
1	A	188	GLN
1	A	189	ARG
1	A	206	ARG
1	A	207	MSE
1	A	214	ASP
1	A	247	THR
1	A	249[A]	ASP
1	A	249[B]	ASP
1	A	272	ARG
1	A	297	ASP
1	A	307	ARG
1	A	327	LEU
1	A	338	LEU
1	B	11	THR
1	B	49	THR
1	B	51	SER
1	B	66	LEU
1	B	69	GLU
1	B	110	THR
1	B	111	PHE
1	B	134	THR
1	B	149	THR
1	B	151	GLU
1	B	178	SER
1	B	192	GLU
1	B	196	VAL
1	B	206	ARG
1	B	214	ASP
1	B	233	CYS

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Mol	Chain	Res	Type
1	B	247	THR
1	B	297	ASP
1	B	301	ILE
1	B	309	THR
1	B	310	HIS
1	B	311	GLU
1	B	338	LEU
1	B	342	THR
1	B	359	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	GLN
1	B	310	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 ⓘ

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	339/386 (87%)	-0.77	1 (0%) 90 88	5, 21, 47, 73	3 (0%)
1	B	331/386 (85%)	0.62	25 (7%) 21 18	20, 73, 100, 106	0
All	All	670/772 (86%)	-0.08	26 (3%) 44 37	5, 47, 96, 106	3 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	37	PRO	4.6
1	B	247	THR	3.9
1	B	238	VAL	3.7
1	B	186	ASP	3.4
1	B	359	TRP	3.2
1	A	4	GLY	3.2
1	B	179	GLU	2.9
1	B	214	ASP	2.9
1	B	150	ASP	2.9
1	B	239	PHE	2.9
1	B	270	ASP	2.8
1	B	4	GLY	2.8
1	B	243	ALA	2.7
1	B	240	ALA	2.6
1	B	246	PRO	2.5
1	B	96	GLY	2.4
1	B	9	LEU	2.2
1	B	98	ILE	2.2
1	B	253	VAL	2.2
1	B	151	GLU	2.2
1	B	338	LEU	2.2
1	B	343	PRO	2.1
1	B	54	VAL	2.1
1	B	56	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	121	ILE	2.0
1	B	99	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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.