



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:41 PM UTC

PDB ID : 6KHR / pdb_00006khr
Title : Structure of glycynamide-RNase-transformylase T from Mycobacterium tuberculosis
Authors : Chen, C.; Wang, J.
Deposited on : 2019-07-16
Resolution : 2.79 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

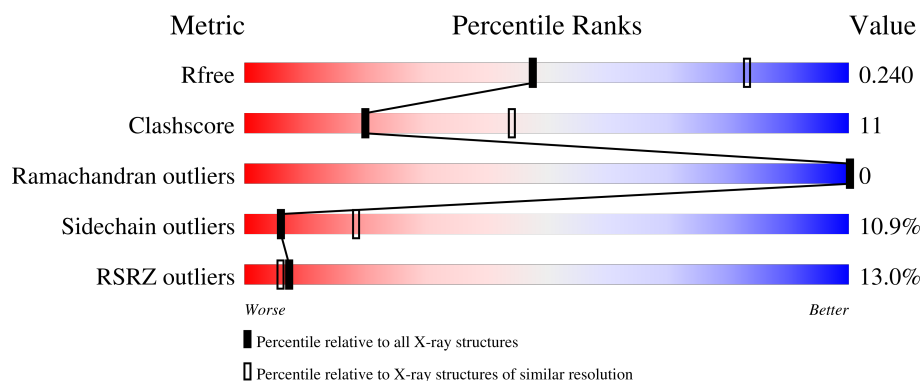
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	420	
1	B	420	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4920 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 Formate-dependent phosphoribosylglycinamide formyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	349	Total	C	N	O	S	0	0	0
			2543	1604	450	478	11			
1	B	320	Total	C	N	O	S	0	0	0
			2343	1476	414	442	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P95197
A	2	VAL	-	expression tag	UNP P95197
B	1	MET	-	initiating methionine	UNP P95197
B	2	VAL	-	expression tag	UNP P95197

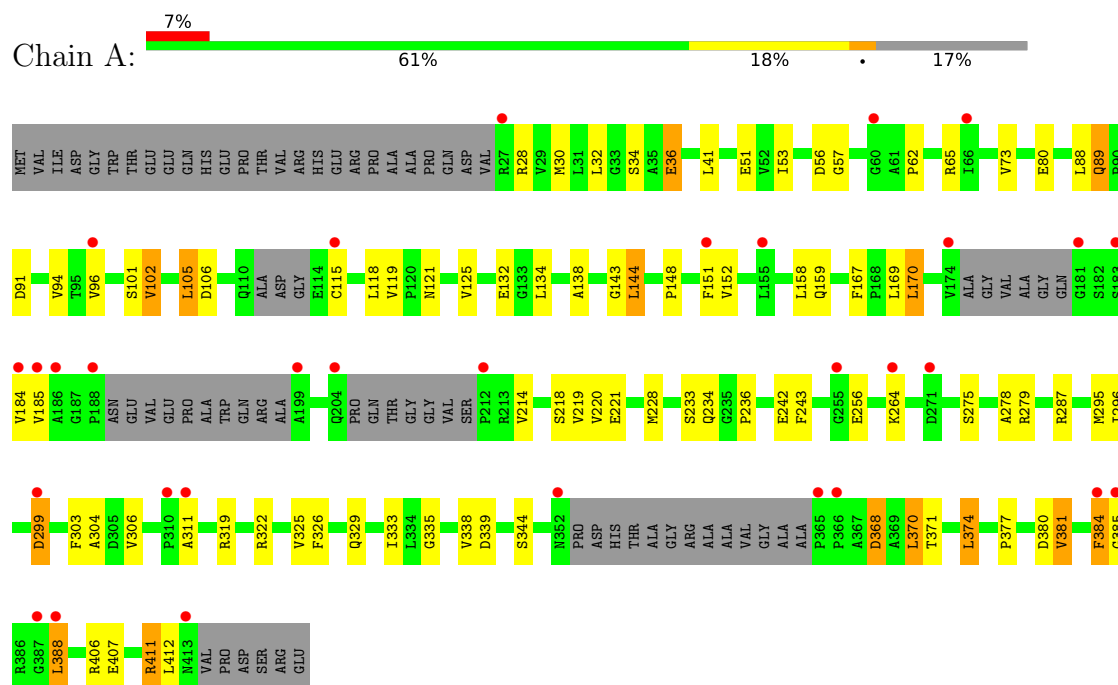
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	15	Total	O	0	0
			15	15		
2	B	19	Total	O	0	0
			19	19		

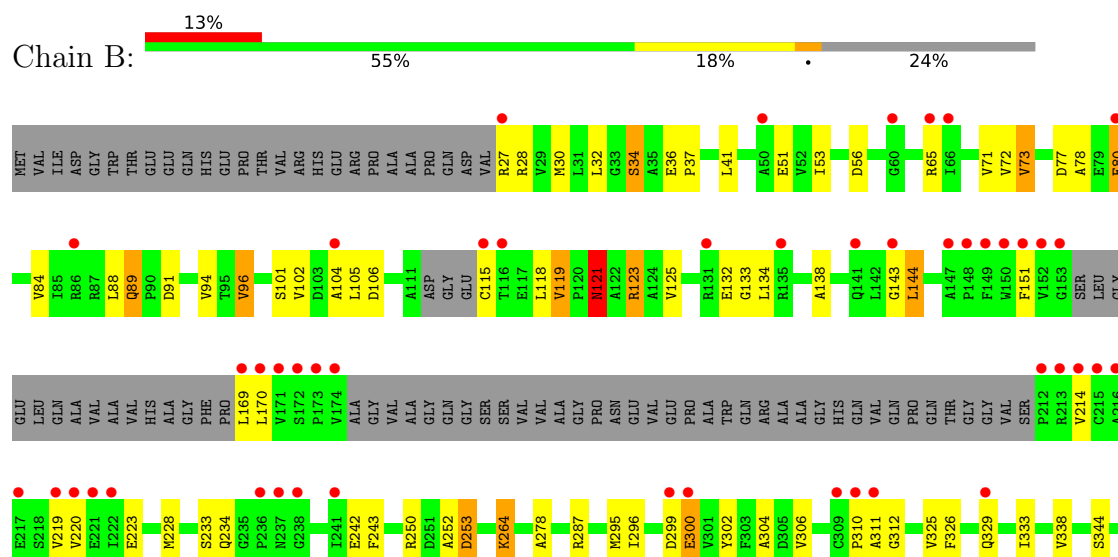
3 Residue-property plots

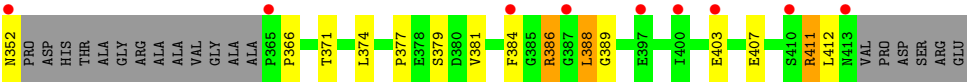
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Formate-dependent phosphoribosylglycinamide formyltransferase



- Molecule 1: Formate-dependent phosphoribosylglycinamide formyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	103.13Å 103.13Å 148.69Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.72 – 2.79 48.72 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.72-2.79) 99.2 (48.72-2.79)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.18 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.14rc3_3206	Depositor
R, R_{free}	0.217 , 0.242 0.216 , 0.240	Depositor DCC
R_{free} test set	1137 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	57.1	Xtriage
Anisotropy	0.291	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 62.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4920	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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	0.73	3/2579 (0.1%)	0.92	11/3513 (0.3%)
1	B	0.83	9/2374 (0.4%)	0.89	1/3234 (0.0%)
All	All	0.78	12/4953 (0.2%)	0.91	12/6747 (0.2%)

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 (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	36	GLU	C-N	12.29	1.49	1.33
1	B	37	PRO	C-O	-6.75	1.16	1.24
1	B	119	VAL	C-O	-6.59	1.16	1.24
1	B	377	PRO	C-O	-6.34	1.16	1.23
1	B	118	LEU	C-O	-5.93	1.16	1.24
1	B	379	SER	C-O	-5.82	1.16	1.23
1	A	377	PRO	C-O	-5.64	1.17	1.23
1	B	96	VAL	C-O	-5.59	1.17	1.24
1	B	34	SER	C-O	-5.50	1.16	1.23
1	A	381	VAL	C-O	-5.35	1.17	1.23
1	B	366	PRO	C-O	-5.15	1.17	1.23
1	B	106	ASP	C-O	-5.06	1.17	1.24

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	36	GLU	CA-C-N	11.12	131.51	119.28
1	A	36	GLU	C-N-CA	11.12	131.51	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	385	GLY	CA-C-N	-9.92	109.03	122.72
1	A	385	GLY	C-N-CA	-9.92	109.03	122.72
1	A	384	PHE	CA-C-N	-7.79	114.61	122.63
1	A	384	PHE	C-N-CA	-7.79	114.61	122.63
1	A	121	ASN	CB-CA-C	-7.62	95.98	109.62
1	A	299	ASP	CB-CA-C	-7.49	94.35	109.55
1	A	256	GLU	CB-CA-C	-5.79	99.28	109.71
1	B	121	ASN	CB-CA-C	5.61	121.18	109.68
1	A	36	GLU	O-C-N	-5.54	115.55	120.48
1	A	384	PHE	CA-C-O	-5.11	114.19	119.51

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	384	PHE	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	2543	0	2581	58	1
1	B	2343	0	2392	62	1
2	A	15	0	0	7	0
2	B	19	0	0	9	0
All	All	4920	0	4973	112	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 11.

All (112) 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:299:ASP:CG	1:B:253:ASP:OD1	1.95	1.10
1:A:169:LEU:HB2	1:A:185:VAL:HG13	1.34	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:LEU:CB	1:A:185:VAL:HG13	1.92	0.98
1:A:299:ASP:OD2	1:B:253:ASP:OD1	1.80	0.97
1:B:352:ASN:CG	2:B:501:HOH:O	2.10	0.93
1:B:352:ASN:OD1	2:B:501:HOH:O	1.85	0.93
1:B:133:GLY:O	2:B:502:HOH:O	1.89	0.88
1:A:299:ASP:OD1	1:B:253:ASP:OD1	1.93	0.86
1:B:310:PRO:O	2:B:503:HOH:O	1.95	0.84
1:B:352:ASN:O	2:B:504:HOH:O	1.97	0.82
1:A:170:LEU:HD23	1:A:219:VAL:HG22	1.59	0.82
1:A:299:ASP:OD2	1:B:252:ALA:HB1	1.82	0.79
1:B:123:ARG:HH11	1:B:123:ARG:HG3	1.47	0.79
1:B:41:LEU:HD13	1:B:329:GLN:HG3	1.65	0.79
1:A:41:LEU:HD13	1:A:329:GLN:HG3	1.65	0.78
1:A:279:ARG:NE	2:A:501:HOH:O	1.94	0.75
1:B:311:ALA:C	2:B:503:HOH:O	2.32	0.73
1:A:335:GLY:O	2:A:502:HOH:O	2.05	0.73
1:A:28:ARG:NH2	1:A:88:LEU:O	2.26	0.68
1:B:28:ARG:NH2	1:B:88:LEU:O	2.26	0.68
1:B:123:ARG:HH11	1:B:123:ARG:CG	2.08	0.67
1:B:121:ASN:C	1:B:121:ASN:HD22	2.02	0.67
1:A:303:PHE:O	2:A:505:HOH:O	2.12	0.67
1:B:388:LEU:HD22	1:B:389:GLY:N	2.10	0.66
1:A:311:ALA:O	2:A:504:HOH:O	2.12	0.66
1:A:169:LEU:CB	1:A:185:VAL:CG1	2.72	0.65
1:A:170:LEU:HD23	1:A:219:VAL:CG2	2.26	0.64
1:A:36:GLU:HG2	1:A:384:PHE:CG	2.33	0.63
1:A:368:ASP:OD1	1:A:368:ASP:N	2.31	0.62
1:A:295:MET:HE2	1:A:304:ALA:HB2	1.82	0.62
1:B:295:MET:HE2	1:B:304:ALA:HB2	1.82	0.61
1:A:152:VAL:HG21	1:A:158:LEU:HG	1.81	0.60
1:A:89:GLN:NE2	1:A:115:CYS:SG	2.74	0.60
1:A:233:SER:HA	1:A:287:ARG:HG2	1.84	0.60
1:B:89:GLN:NE2	1:B:115:CYS:SG	2.74	0.59
1:B:233:SER:HA	1:B:287:ARG:HG2	1.84	0.58
1:A:119:VAL:HG11	1:A:333:ILE:HA	1.85	0.58
1:A:370:LEU:HD22	1:A:374:LEU:HD22	1.87	0.57
1:B:170:LEU:HG	1:B:219:VAL:HG22	1.87	0.57
1:A:169:LEU:HB3	1:A:185:VAL:CG1	2.34	0.56
1:B:300:GLU:OE2	1:B:302:TYR:OH	2.15	0.56
1:B:78:ALA:HA	1:B:104:ALA:HB2	1.88	0.55
1:B:94:VAL:HG13	1:B:119:VAL:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:PHE:CD2	1:A:388:LEU:HD12	2.43	0.54
1:B:28:ARG:NH1	1:B:53:ILE:HD11	2.24	0.53
1:A:299:ASP:HB3	1:B:250:ARG:HH22	1.73	0.53
1:A:28:ARG:N	1:A:91:ASP:OD2	2.39	0.53
1:B:371:THR:O	2:B:505:HOH:O	2.19	0.53
1:A:221:GLU:HB3	1:B:403:GLU:OE1	2.09	0.52
1:A:132:GLU:HG2	1:A:151:PHE:CE2	2.45	0.51
1:A:28:ARG:NH1	1:A:53:ILE:HD11	2.24	0.51
1:B:132:GLU:HG2	1:B:151:PHE:CE2	2.45	0.51
1:A:279:ARG:CD	2:A:501:HOH:O	2.53	0.50
1:A:148:PRO:HB2	1:A:218:SER:HB2	1.94	0.50
1:B:253:ASP:OD1	1:B:253:ASP:N	2.44	0.49
1:B:170:LEU:HG	1:B:219:VAL:CG2	2.43	0.49
1:B:386:ARG:O	1:B:386:ARG:HG3	2.13	0.49
1:B:134:LEU:HD21	1:B:306:VAL:HG12	1.95	0.48
1:A:134:LEU:HD21	1:A:306:VAL:HG12	1.95	0.48
1:B:27:ARG:N	2:B:508:HOH:O	2.45	0.48
1:A:101:SER:O	1:A:105:LEU:HD22	2.14	0.48
1:B:119:VAL:HG11	1:B:333:ILE:HA	1.94	0.48
1:A:325:VAL:O	1:A:329:GLN:HG2	2.14	0.47
1:B:102:VAL:HG12	1:B:125:VAL:HG12	1.96	0.47
1:B:77:ASP:HB3	1:B:80:GLU:HB2	1.95	0.47
1:A:228:MET:HE3	1:A:228:MET:HB3	1.77	0.47
1:B:28:ARG:N	1:B:91:ASP:OD2	2.39	0.47
1:A:30:MET:HE2	1:A:88:LEU:HD22	1.97	0.47
1:B:325:VAL:O	1:B:329:GLN:HG2	2.14	0.47
1:B:123:ARG:CG	1:B:123:ARG:NH1	2.72	0.46
1:B:30:MET:HE2	1:B:88:LEU:HD22	1.97	0.46
1:A:34:SER:O	1:A:62:PRO:HD2	2.16	0.45
1:A:53:ILE:HD13	1:A:88:LEU:HD23	1.99	0.45
1:B:72:VAL:C	1:B:73:VAL:CG2	2.91	0.44
1:B:73:VAL:HG21	1:B:84:VAL:HG21	1.99	0.44
1:B:326:PHE:HA	1:B:329:GLN:HG2	1.99	0.44
1:A:242:GLU:HA	2:A:503:HOH:O	2.17	0.44
1:A:275:SER:O	1:A:279:ARG:HG3	2.18	0.44
1:A:143:GLY:CA	1:B:264:LYS:HE2	2.48	0.44
1:A:326:PHE:HA	1:A:329:GLN:HG2	1.99	0.44
1:A:138:ALA:O	1:A:144:LEU:HB2	2.19	0.43
1:A:384:PHE:HD2	1:A:388:LEU:HD12	1.83	0.43
1:B:53:ILE:HD13	1:B:88:LEU:HD23	1.99	0.43
1:A:242:GLU:HG3	1:A:338:VAL:HG23	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:PHE:N	2:A:503:HOH:O	2.10	0.43
1:B:121:ASN:C	1:B:121:ASN:ND2	2.74	0.43
1:B:36:GLU:HG2	1:B:388:LEU:HD12	2.00	0.43
1:B:71:VAL:HG12	1:B:73:VAL:CG2	2.49	0.43
1:A:167:PHE:HE1	1:A:185:VAL:HG22	1.83	0.43
1:B:138:ALA:O	1:B:144:LEU:HB2	2.18	0.43
1:B:94:VAL:HG21	1:B:333:ILE:HD13	2.01	0.43
1:B:388:LEU:HD22	1:B:388:LEU:C	2.44	0.42
1:B:101:SER:HB3	1:B:104:ALA:HB3	2.00	0.42
1:B:228:MET:HE3	1:B:228:MET:HB3	1.77	0.42
1:B:242:GLU:HG3	1:B:338:VAL:HG23	2.01	0.42
1:A:102:VAL:HG12	1:A:125:VAL:HG12	2.01	0.42
1:A:56:ASP:OD1	1:A:57:GLY:N	2.46	0.42
1:A:94:VAL:HG21	1:A:333:ILE:HD13	2.01	0.42
1:A:279:ARG:NH2	1:B:223:GLU:OE2	2.53	0.42
1:A:184:VAL:O	1:A:184:VAL:HG23	2.20	0.41
1:B:34:SER:HB2	1:B:56:ASP:HB2	2.01	0.41
1:B:312:GLY:N	2:B:503:HOH:O	2.51	0.41
1:B:72:VAL:C	1:B:73:VAL:HG22	2.45	0.41
1:B:243:PHE:HZ	1:B:278:ALA:HB2	1.86	0.41
1:B:169:LEU:O	1:B:170:LEU:HD23	2.21	0.41
1:B:388:LEU:HD22	1:B:389:GLY:H	1.85	0.41
1:A:169:LEU:HA	1:A:169:LEU:HD23	1.86	0.41
1:A:322:ARG:NH2	1:A:339:ASP:OD2	2.54	0.41
1:B:407:GLU:O	1:B:411:ARG:HG2	2.22	0.40
1:A:243:PHE:HZ	1:A:278:ALA:HB2	1.86	0.40
1:A:319:ARG:NH2	1:A:380:ASP:OD2	2.49	0.40
1:A:407:GLU:O	1:A:411:ARG:HG2	2.22	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:236:PRO:O	1:B:143:GLY:O[5_554]	1.86	0.34

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	337/420 (80%)	321 (95%)	16 (5%)	0	100	100
1	B	310/420 (74%)	299 (96%)	11 (4%)	0	100	100
All	All	647/840 (77%)	620 (96%)	27 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	262/316 (83%)	233 (89%)	29 (11%)	6	18
1	B	243/316 (77%)	217 (89%)	26 (11%)	6	19
All	All	505/632 (80%)	450 (89%)	55 (11%)	6	18

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LEU
1	A	51	GLU
1	A	65	ARG
1	A	73	VAL
1	A	80	GLU
1	A	89	GLN
1	A	96	VAL
1	A	102	VAL

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Mol	Chain	Res	Type
1	A	105	LEU
1	A	106	ASP
1	A	118	LEU
1	A	144	LEU
1	A	159	GLN
1	A	170	LEU
1	A	214	VAL
1	A	220	VAL
1	A	234	GLN
1	A	264	LYS
1	A	296	ILE
1	A	344	SER
1	A	368	ASP
1	A	370	LEU
1	A	371	THR
1	A	374	LEU
1	A	381	VAL
1	A	388	LEU
1	A	406	ARG
1	A	411	ARG
1	A	412	LEU
1	B	32	LEU
1	B	51	GLU
1	B	65	ARG
1	B	73	VAL
1	B	80	GLU
1	B	89	GLN
1	B	96	VAL
1	B	105	LEU
1	B	121	ASN
1	B	123	ARG
1	B	144	LEU
1	B	214	VAL
1	B	220	VAL
1	B	234	GLN
1	B	253	ASP
1	B	264	LYS
1	B	296	ILE
1	B	299	ASP
1	B	300	GLU
1	B	344	SER
1	B	374	LEU

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Mol	Chain	Res	Type
1	B	381	VAL
1	B	386	ARG
1	B	388	LEU
1	B	411	ARG
1	B	412	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	121	ASN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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	349/420 (83%)	0.56	31 (8%) 15 12	27, 58, 111, 184	0
1	B	320/420 (76%)	0.97	56 (17%) 4 3	28, 51, 142, 197	0
All	All	669/840 (79%)	0.76	87 (13%) 7 6	27, 54, 119, 197	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	170	LEU	8.9
1	B	171	VAL	8.7
1	B	174	VAL	7.0
1	B	216	ALA	6.6
1	B	311	ALA	6.5
1	A	311	ALA	6.0
1	B	169	LEU	6.0
1	B	152	VAL	5.7
1	B	115	CYS	5.5
1	B	173	PRO	5.4
1	A	185	VAL	5.2
1	A	212	PRO	5.1
1	B	153	GLY	5.0
1	B	214	VAL	4.7
1	A	174	VAL	4.7
1	B	149	PHE	4.5
1	B	310	PRO	4.5
1	A	181	GLY	4.3
1	A	384	PHE	4.3
1	A	310	PRO	4.2
1	A	60	GLY	4.2
1	A	365	PRO	4.1
1	B	212	PRO	4.0
1	B	215	CYS	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	150	TRP	4.0
1	B	397	GLU	3.9
1	A	388	LEU	3.9
1	B	217	GLU	3.8
1	B	413	ASN	3.7
1	B	238	GLY	3.7
1	A	188	PRO	3.6
1	B	151	PHE	3.6
1	B	147	ALA	3.5
1	B	365	PRO	3.5
1	A	255	GLY	3.4
1	B	172	SER	3.3
1	B	148	PRO	3.2
1	A	199	ALA	3.2
1	A	183	SER	3.2
1	B	143	GLY	3.2
1	B	299	ASP	3.2
1	A	387	GLY	3.1
1	A	366	PRO	3.0
1	B	236	PRO	3.0
1	B	387	GLY	3.0
1	A	27	ARG	3.0
1	A	352	ASN	3.0
1	B	309	CYS	2.9
1	A	151	PHE	2.9
1	B	27	ARG	2.9
1	B	410	SER	2.9
1	B	104	ALA	2.8
1	B	86	ARG	2.8
1	B	237	ASN	2.8
1	A	115	CYS	2.7
1	A	184	VAL	2.6
1	B	116	THR	2.6
1	B	60	GLY	2.6
1	B	219	VAL	2.5
1	B	300	GLU	2.5
1	B	131	ARG	2.4
1	B	213	ARG	2.4
1	B	400	ILE	2.4
1	B	66	ILE	2.4
1	B	403	GLU	2.4
1	B	220	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	135	ARG	2.4
1	A	204	GLN	2.4
1	B	80	GLU	2.4
1	B	222	ILE	2.3
1	B	241	ILE	2.3
1	A	299	ASP	2.3
1	B	329	GLN	2.3
1	A	271	ASP	2.2
1	B	50	ALA	2.2
1	B	141	GLN	2.2
1	A	186	ALA	2.2
1	B	65	ARG	2.1
1	A	96	VAL	2.1
1	B	384	PHE	2.1
1	B	352	ASN	2.1
1	A	155	LEU	2.1
1	A	264	LYS	2.0
1	A	413	ASN	2.0
1	A	66	ILE	2.0
1	B	221	GLU	2.0
1	A	385	GLY	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 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.