



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:31 PM UTC

PDB ID : 1M1Z / pdb_00001m1z
Title : BETA-LACTAM SYNTHETASE APO ENZYME
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Deposited on : 2002-06-20
Resolution : 1.95 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

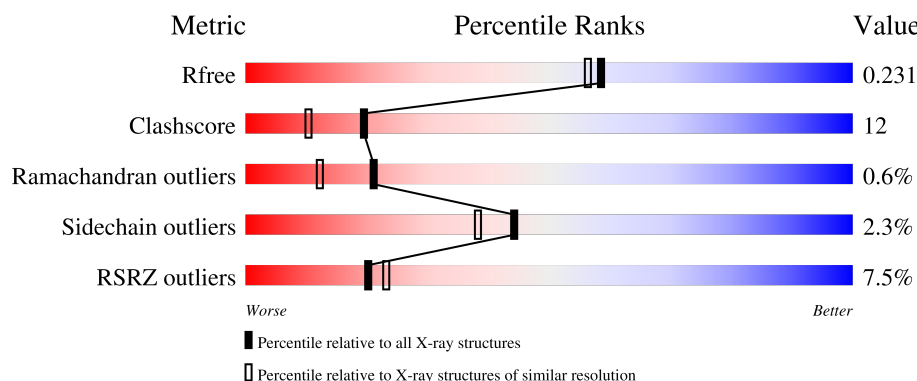
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	513	10% 71% 20% • 6%
1	B	513	5% 73% 20% • •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7883 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 BETA-LACTAM SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	480	Total	C	N	O	S	0	0	0
			3604	2254	659	683	8			
1	B	490	Total	C	N	O	S	0	0	0
			3676	2296	674	698	8			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	268	Total	O	0	0
			268	268		
2	B	335	Total	O	0	0
			335	335		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.18Å 97.47Å 80.97Å 90.00° 90.42° 90.00°	Depositor
Resolution (Å)	23.88 – 1.95 23.88 – 1.95	Depositor EDS
% Data completeness (in resolution range)	91.2 (23.88-1.95) 91.3 (23.88-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.00 (at 1.91Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.197 , 0.231 0.197 , 0.231	Depositor DCC
R_{free} test set	6156 reflections (8.23%)	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.036 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7883	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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	0.40	1/3674 (0.0%)	0.93	10/5012 (0.2%)
1	B	0.53	4/3749 (0.1%)	1.18	25/5116 (0.5%)
All	All	0.47	5/7423 (0.1%)	1.07	35/10128 (0.3%)

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	1
1	B	0	4
All	All	0	5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	93	PRO	N-CD	13.49	1.66	1.47
1	B	92	GLY	C-N	6.54	1.46	1.34
1	B	91	ALA	CA-C	6.15	1.60	1.52
1	A	455	SER	N-CA	5.85	1.53	1.46
1	B	91	ALA	C-N	-5.65	1.25	1.33

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	92	GLY	O-C-N	-29.74	92.03	121.77
1	B	91	ALA	CA-C-N	23.02	158.01	121.87
1	B	91	ALA	C-N-CA	23.02	158.01	121.87
1	B	93	PRO	CA-N-CD	-13.68	92.34	111.50
1	B	91	ALA	CA-C-O	-13.04	106.90	120.98
1	A	454	SER	CA-C-N	-9.48	107.57	120.28
1	A	454	SER	C-N-CA	-9.48	107.57	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	453	THR	CB-CA-C	-8.91	96.90	110.26
1	B	93	PRO	N-CA-CB	8.78	112.26	102.60
1	B	173	ALA	CA-C-N	8.48	135.85	122.61
1	B	173	ALA	C-N-CA	8.48	135.85	122.61
1	A	454	SER	O-C-N	-7.72	110.64	123.00
1	B	92	GLY	C-N-CD	7.61	137.34	120.60
1	A	455	SER	CB-CA-C	7.58	123.38	110.79
1	B	92	GLY	N-CA-C	7.50	127.64	112.34
1	A	117	VAL	N-CA-C	7.50	119.64	108.54
1	B	117	VAL	N-CA-C	7.28	119.31	108.54
1	B	401	ASP	N-CA-C	-6.62	101.16	110.50
1	A	29	ALA	N-CA-C	-6.50	104.05	112.23
1	A	401	ASP	N-CA-C	-6.39	100.97	110.23
1	B	29	ALA	N-CA-C	-6.32	104.62	112.90
1	B	343	ARG	N-CA-C	-6.29	98.48	108.73
1	B	13	LEU	N-CA-C	-5.99	98.31	108.26
1	B	390	LEU	N-CA-C	-5.90	105.93	113.01
1	B	93	PRO	N-CD-CG	-5.65	97.02	103.80
1	B	195	ILE	N-CA-C	5.61	116.38	108.36
1	A	13	LEU	N-CA-C	-5.52	99.09	108.26
1	B	174	ASP	OD1-CG-OD2	-5.49	109.72	122.90
1	B	93	PRO	CB-CA-C	5.32	125.30	112.00
1	A	110	ASP	CB-CA-C	-5.29	108.93	116.34
1	A	195	ILE	N-CA-C	5.25	115.52	108.17
1	B	57	ALA	CA-C-N	5.13	126.25	119.84
1	B	57	ALA	C-N-CA	5.13	126.25	119.84
1	B	122	ALA	N-CA-C	-5.09	100.41	109.06
1	B	248	LEU	N-CA-C	5.03	117.16	109.07

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	454	SER	Mainchain
1	B	174	ASP	Sidechain
1	B	452	THR	Mainchain
1	B	453	THR	Mainchain
1	B	91	ALA	Mainchain

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	3604	0	3595	81	0
1	B	3676	0	3663	93	0
2	A	268	0	0	4	0
2	B	335	0	0	4	0
All	All	7883	0	7258	169	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 12.

All (169) 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:41:GLN:HE21	1:B:45:SER:HB3	1.28	0.93
1:A:92:GLY:H	1:A:93:PRO:HD2	1.34	0.92
1:B:308:GLN:NE2	1:B:335:ARG:HH11	1.73	0.86
1:A:112:HIS:HD2	1:A:115:ARG:HE	1.23	0.85
1:B:174:ASP:O	1:B:188:PRO:HG3	1.79	0.82
1:B:31:ARG:HH11	1:B:31:ARG:HB2	1.45	0.81
1:A:41:GLN:HG2	1:A:45:SER:OG	1.86	0.75
1:A:216:ILE:HD11	1:B:116:LEU:HD22	1.70	0.74
1:A:148:ALA:HB3	1:A:151:GLU:HG3	1.70	0.73
1:A:216:ILE:HD12	1:B:88:VAL:HG11	1.72	0.72
1:B:186:GLN:HE22	1:B:381:ASN:HD21	1.37	0.72
1:B:174:ASP:O	1:B:188:PRO:CG	2.38	0.71
1:B:383:MET:HG2	1:B:397:HIS:CD2	2.25	0.71
1:B:41:GLN:NE2	1:B:45:SER:HB3	2.06	0.69
1:B:348:TYR:CE2	1:B:383:MET:HG3	2.28	0.69
1:B:308:GLN:HE21	1:B:335:ARG:HD3	1.60	0.66
1:B:328:LEU:HB2	1:B:329:PRO:HD3	1.79	0.64
1:A:95:PRO:HB3	1:A:100:GLU:HB3	1.79	0.64
1:A:314:TRP:CD1	1:A:485:THR:HB	2.34	0.62
1:B:214:ARG:NH2	1:B:370:LEU:HD12	2.14	0.62
1:B:27:VAL:O	1:B:27:VAL:HG23	1.99	0.61
1:A:13:LEU:HD23	1:A:26:PRO:HG3	1.83	0.60
1:B:169:GLY:HA3	1:B:183:GLY:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:LEU:O	1:B:107:GLU:HG3	2.03	0.59
1:A:456:PHE:CE2	1:A:474:LYS:HD3	2.38	0.59
1:B:308:GLN:NE2	1:B:335:ARG:NH1	2.49	0.59
1:A:239:THR:HG22	1:A:240:PRO:HD2	1.84	0.59
1:A:273:MET:HB2	1:A:333:LEU:HD13	1.83	0.59
1:B:176:ARG:HD2	2:B:657:HOH:O	2.01	0.58
1:B:140:SER:HB2	1:B:383:MET:HE2	1.85	0.58
1:B:41:GLN:HE21	1:B:45:SER:CB	2.08	0.58
1:B:437:GLU:HG2	1:B:441:ARG:NH2	2.19	0.58
1:A:437:GLU:O	1:A:441:ARG:HB2	2.04	0.57
1:B:314:TRP:CD1	1:B:485:THR:HB	2.39	0.57
1:A:456:PHE:CD2	1:A:474:LYS:HD3	2.40	0.57
1:B:26:PRO:HG3	1:B:49:THR:HG21	1.87	0.57
1:A:52:HIS:NE2	1:A:58:PRO:HG3	2.19	0.56
1:B:352:ILE:O	1:B:370:LEU:HD23	2.05	0.56
1:A:279:ASN:HD22	1:A:281:PHE:H	1.53	0.56
1:B:297:GLU:C	1:B:298:ILE:HD12	2.30	0.56
1:A:250:GLY:O	1:A:441:ARG:HD2	2.06	0.56
1:B:325:GLU:HG2	1:B:456:PHE:HD2	1.72	0.55
1:B:298:ILE:HD12	1:B:298:ILE:N	2.22	0.55
1:A:309:LEU:HD11	1:A:456:PHE:CE1	2.41	0.55
1:B:437:GLU:HG2	1:B:441:ARG:HH22	1.71	0.55
1:B:308:GLN:NE2	1:B:335:ARG:HD3	2.22	0.55
1:B:354:LEU:HD21	1:B:408:LEU:HD22	1.88	0.54
1:B:490:ARG:HG2	1:B:494:GLU:HG3	1.90	0.54
1:B:327:LEU:HD13	1:B:387:LEU:HD21	1.87	0.54
1:B:386:VAL:HA	1:B:389:THR:OG1	2.07	0.54
1:A:92:GLY:N	1:A:93:PRO:HD2	2.14	0.53
1:A:112:HIS:CD2	1:A:115:ARG:HE	2.14	0.53
1:B:348:TYR:OH	1:B:382:GLU:HG3	2.08	0.53
1:B:65:SER:HB3	1:B:71:THR:HB	1.90	0.53
1:B:453:THR:CG2	1:B:457:SER:HB2	2.39	0.53
1:B:472:GLU:HG3	1:B:476:GLN:HE21	1.73	0.53
1:B:215:ARG:HH11	1:B:215:ARG:HG3	1.74	0.52
1:B:370:LEU:C	1:B:370:LEU:HD13	2.35	0.52
1:A:313:VAL:HG11	1:A:478:VAL:HG13	1.91	0.52
1:A:454:SER:O	1:A:455:SER:C	2.48	0.52
1:A:110:ASP:CG	1:A:111:LEU:H	2.18	0.51
1:B:353:PRO:HA	1:B:370:LEU:HD21	1.92	0.51
1:B:422:ASP:OD2	1:B:423:LYS:HE2	2.09	0.51
1:B:305:LEU:C	1:B:305:LEU:HD23	2.36	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:LEU:O	1:B:89:LEU:HD13	2.11	0.51
1:A:302:THR:CG2	1:A:455:SER:HB2	2.41	0.51
1:B:480:GLU:OE1	1:B:500:VAL:HG13	2.11	0.50
1:B:341:GLU:HG2	2:B:542:HOH:O	2.10	0.50
1:B:382:GLU:HG2	2:B:827:HOH:O	2.10	0.50
1:B:453:THR:HG22	1:B:457:SER:HB2	1.93	0.50
1:A:349:GLY:O	1:A:353:PRO:HD2	2.11	0.50
1:A:212:LEU:HD13	1:B:206:ARG:HB2	1.93	0.49
1:B:110:ASP:CG	1:B:111:LEU:H	2.21	0.49
1:B:333:LEU:C	1:B:333:LEU:HD23	2.37	0.49
1:A:52:HIS:CD2	1:A:58:PRO:HG3	2.48	0.49
1:B:327:LEU:HD22	1:B:387:LEU:HD13	1.95	0.49
1:B:390:LEU:C	1:B:390:LEU:HD12	2.38	0.49
1:A:386:VAL:HA	1:A:389:THR:OG1	2.12	0.48
1:B:85:LEU:O	1:B:88:VAL:HG22	2.13	0.48
1:B:348:TYR:HH	1:B:382:GLU:CD	2.20	0.48
1:A:30:THR:HG23	1:A:389:THR:HB	1.96	0.48
1:A:112:HIS:HE1	1:B:368:THR:OG1	1.97	0.48
1:B:348:TYR:OH	1:B:382:GLU:CD	2.56	0.48
1:A:345:LEU:HD22	1:A:396:THR:CG2	2.43	0.48
1:B:358:HIS:HB3	1:B:417:ARG:NH2	2.28	0.48
1:A:17:ARG:NH1	1:A:151:GLU:HB3	2.29	0.48
1:A:148:ALA:HB3	1:A:151:GLU:CG	2.41	0.47
1:B:353:PRO:HA	1:B:370:LEU:CD2	2.44	0.47
1:A:29:ALA:HA	1:A:314:TRP:CH2	2.50	0.47
1:B:166:ASP:O	1:B:168:LYS:HD2	2.14	0.47
1:B:325:GLU:OE2	1:B:456:PHE:HB3	2.14	0.47
1:B:422:ASP:C	1:B:423:LYS:HG2	2.39	0.47
1:B:29:ALA:HA	1:B:314:TRP:CH2	2.49	0.47
1:A:4:PRO:HB3	2:A:717:HOH:O	2.15	0.47
1:B:170:PHE:HB3	1:B:171:PRO:HD2	1.95	0.47
1:B:195:ILE:HG12	1:B:202:ALA:HB2	1.97	0.47
1:B:41:GLN:HB2	1:B:45:SER:HB3	1.97	0.47
1:A:73:ALA:HB2	1:A:103:LEU:HD13	1.97	0.46
1:B:308:GLN:HE22	1:B:335:ARG:NH1	2.12	0.46
1:B:473:ALA:O	1:B:477:VAL:HG23	2.15	0.46
1:A:435:PRO:HG2	1:A:438:THR:OG1	2.16	0.46
1:A:249:SER:HB3	2:A:519:HOH:O	2.16	0.46
1:B:431:ALA:HA	1:B:439:VAL:HG11	1.98	0.46
1:A:303:THR:HG23	1:A:304:GLU:N	2.30	0.46
1:B:308:GLN:HE21	1:B:335:ARG:HH11	1.57	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:308:GLN:HE22	1:B:335:ARG:HH11	1.60	0.46
1:A:127:THR:HG23	1:A:127:THR:O	2.15	0.46
1:A:496:ASP:O	1:A:500:VAL:HG23	2.16	0.46
1:A:223:VAL:HG13	1:A:429:ALA:HA	1.97	0.46
1:A:308:GLN:HB2	1:A:328:LEU:HD22	1.97	0.45
1:A:348:TYR:HH	1:A:382:GLU:CD	2.24	0.45
1:B:273:MET:HE2	1:B:333:LEU:HA	1.97	0.45
1:A:214:ARG:HB2	1:B:116:LEU:HD23	1.97	0.45
1:A:132:LEU:HD23	1:A:194:ASP:HA	1.99	0.45
1:A:268:LEU:O	1:A:293:THR:HB	2.17	0.45
1:A:497:THR:O	1:A:501:VAL:HG23	2.17	0.45
1:B:17:ARG:NH2	1:B:20:GLY:HA2	2.32	0.45
1:A:313:VAL:CG1	1:A:478:VAL:HG13	2.47	0.45
1:A:498:ASP:HB3	1:A:502:ARG:NH2	2.32	0.44
1:A:96:GLU:HG2	1:A:100:GLU:OE1	2.17	0.44
1:B:89:LEU:HD11	1:B:105:LEU:HD23	1.99	0.44
1:B:281:PHE:O	1:B:285:ARG:HG3	2.18	0.44
1:A:500:VAL:O	1:A:504:VAL:HG23	2.17	0.44
1:A:95:PRO:HG3	1:A:104:ARG:HG3	2.00	0.43
1:A:178:VAL:HG11	1:A:181:LEU:HD12	2.00	0.43
1:B:108:ARG:NH1	1:B:109:TYR:OH	2.52	0.43
1:A:290:HIS:CG	1:A:435:PRO:HD3	2.54	0.43
1:A:421:ARG:HH21	1:A:440:ASN:HA	1.84	0.43
1:A:465:VAL:HG12	1:A:467:GLU:H	1.84	0.43
1:A:60:ARG:HG3	2:A:668:HOH:O	2.18	0.43
1:A:475:ARG:HG2	1:A:475:ARG:HH11	1.83	0.43
1:B:89:LEU:HD11	1:B:105:LEU:CD2	2.49	0.43
1:B:172:LEU:HD13	1:B:193:MET:HE1	2.01	0.43
1:B:237:ARG:O	1:B:343:ARG:HD3	2.18	0.43
1:A:92:GLY:H	1:A:93:PRO:CD	2.17	0.42
1:A:491:HIS:HD2	1:A:493:SER:H	1.68	0.42
1:A:153:ARG:HE	1:A:164:HIS:CD2	2.38	0.42
1:A:390:LEU:HD12	1:A:390:LEU:C	2.45	0.42
1:B:352:ILE:HB	1:B:353:PRO:CD	2.49	0.42
1:A:366:LEU:C	1:A:366:LEU:HD23	2.45	0.42
1:B:43:GLU:H	1:B:43:GLU:HG2	1.70	0.42
1:B:341:GLU:OE1	1:B:341:GLU:N	2.53	0.42
1:A:173:ALA:O	1:A:174:ASP:HB2	2.18	0.42
1:A:300:ILE:HD13	1:A:332:ALA:HB1	2.02	0.42
1:A:352:ILE:HB	1:A:353:PRO:CD	2.50	0.42
1:B:320:ASP:HB3	1:B:323:ILE:HD13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:TYR:CG	1:A:349:GLY:N	2.89	0.41
1:A:199:SER:OG	1:A:201:THR:HG22	2.20	0.41
1:A:341:GLU:H	1:A:341:GLU:HG3	1.64	0.41
1:B:172:LEU:HB2	1:B:175:ALA:HB3	2.01	0.41
1:B:234:VAL:O	1:B:238:VAL:HG23	2.20	0.41
1:A:26:PRO:HD3	2:A:657:HOH:O	2.21	0.41
1:A:308:GLN:HG3	1:A:332:ALA:HB2	2.02	0.41
1:B:237:ARG:HB3	1:B:396:THR:HG21	2.02	0.41
1:A:216:ILE:HD13	1:A:412:GLU:HG2	2.01	0.41
1:B:174:ASP:O	1:B:188:PRO:HG2	2.17	0.41
1:A:483:ASP:O	1:A:488:GLY:N	2.54	0.41
1:A:274:GLY:O	1:A:299:THR:HA	2.20	0.41
1:A:421:ARG:HH22	1:A:439:VAL:HG12	1.86	0.41
1:A:352:ILE:HB	1:A:353:PRO:HD3	2.02	0.41
1:B:274:GLY:O	1:B:299:THR:HA	2.20	0.41
1:A:309:LEU:N	1:A:310:PRO:HD2	2.36	0.40
1:A:328:LEU:N	1:A:329:PRO:HD2	2.36	0.40
1:A:490:ARG:HH11	1:A:490:ARG:HG2	1.85	0.40
1:B:13:LEU:HD23	1:B:26:PRO:HB3	2.03	0.40
1:B:215:ARG:HG3	1:B:215:ARG:NH1	2.36	0.40
1:B:490:ARG:CG	1:B:494:GLU:HG3	2.49	0.40
1:A:363:LEU:N	1:A:364:PRO:CD	2.85	0.40
1:B:95:PRO:HB3	2:B:810:HOH:O	2.21	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	470/513 (92%)	452 (96%)	15 (3%)	3 (1%)	21	12
1	B	482/513 (94%)	463 (96%)	16 (3%)	3 (1%)	21	12

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	952/1026 (93%)	915 (96%)	31 (3%)	6 (1%)	21	12

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	92	GLY
1	B	339	GLY
1	B	19	GLY
1	A	92	GLY
1	A	489	GLY
1	A	240	PRO

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	373/395 (94%)	363 (97%)	10 (3%)	39	32
1	B	381/395 (96%)	374 (98%)	7 (2%)	51	47
All	All	754/790 (95%)	737 (98%)	17 (2%)	44	38

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	SER
1	A	103	LEU
1	A	239	THR
1	A	279	ASN
1	A	282	ARG
1	A	337	LEU
1	A	345	LEU
1	A	360	GLU
1	A	387	LEU
1	A	454	SER
1	B	31	ARG
1	B	41	GLN

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Mol	Chain	Res	Type
1	B	96	GLU
1	B	174	ASP
1	B	308	GLN
1	B	370	LEU
1	B	452	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	HIS
1	A	118	ASN
1	A	164	HIS
1	A	236	GLN
1	A	279	ASN
1	A	358	HIS
1	A	491	HIS
1	B	41	GLN
1	B	236	GLN
1	B	262	HIS
1	B	290	HIS
1	B	308	GLN
1	B	381	ASN
1	B	440	ASN
1	B	476	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 ⓘ

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	480/513 (93%)	0.47	49 (10%)	12 13	13, 26, 63, 76	0
1	B	490/513 (95%)	0.17	24 (4%)	35 41	14, 23, 45, 68	0
All	All	970/1026 (94%)	0.32	73 (7%)	20 23	13, 24, 55, 76	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	473	ALA	6.9
1	A	3	ALA	6.5
1	A	4	PRO	5.7
1	A	18	THR	5.3
1	B	24	PRO	4.9
1	A	454	SER	4.5
1	A	500	VAL	4.4
1	A	456	PHE	4.3
1	A	461	LEU	4.3
1	A	460	LEU	4.3
1	A	170	PHE	4.2
1	A	465	VAL	4.2
1	B	92	GLY	4.1
1	A	250	GLY	4.0
1	B	508	THR	3.9
1	B	93	PRO	3.6
1	B	358	HIS	3.6
1	B	91	ALA	3.5
1	B	19	GLY	3.5
1	B	174	ASP	3.5
1	A	164	HIS	3.2
1	A	455	SER	3.2
1	A	501	VAL	3.2
1	B	18	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	163	ALA	3.1
1	A	504	VAL	2.9
1	A	464	GLY	2.9
1	A	506	ASP	2.8
1	A	443	LYS	2.8
1	A	488	GLY	2.8
1	A	505	ALA	2.8
1	A	321	PRO	2.8
1	A	91	ALA	2.7
1	B	94	ALA	2.7
1	B	25	GLY	2.7
1	B	326	TYR	2.7
1	B	164	HIS	2.6
1	A	93	PRO	2.6
1	B	2	GLY	2.5
1	B	453	THR	2.5
1	A	463	HIS	2.5
1	B	452	THR	2.5
1	A	358	HIS	2.5
1	B	168	LYS	2.5
1	A	503	SER	2.4
1	A	502	ARG	2.4
1	A	495	VAL	2.4
1	A	241	GLY	2.4
1	A	361	ASP	2.3
1	A	17	ARG	2.3
1	B	348	TYR	2.3
1	A	5	VAL	2.3
1	B	165	ARG	2.3
1	A	457	SER	2.3
1	A	459	LEU	2.3
1	A	477	VAL	2.2
1	B	439	VAL	2.2
1	A	474	LYS	2.2
1	A	25	GLY	2.2
1	A	496	ASP	2.2
1	A	95	PRO	2.1
1	B	166	ASP	2.1
1	A	94	ALA	2.1
1	A	338	ASP	2.1
1	B	338	ASP	2.1
1	B	503	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	92	GLY	2.1
1	A	360	GLU	2.0
1	A	19	GLY	2.0
1	B	3	ALA	2.0
1	A	41	GLN	2.0
1	A	497	THR	2.0
1	A	240	PRO	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.