



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:37 PM UTC

PDB ID : 1WW8 / pdb_00001ww8
Title : Crystal Structure of malic enzyme from *Pyrococcus horikoshii* Ot3
Authors : Lokanath, N.K.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-01-04
Resolution : 2.50 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

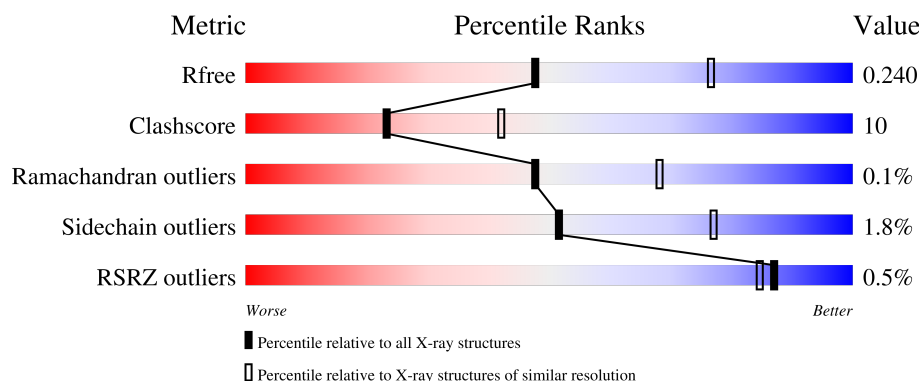
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	439	 72% 25% ..
1	B	439	 76% 21% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6908 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 malate oxidoreductase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	433	Total	C	N	O	S	Se	0	0	0
			3339	2143	570	618	2	6			
1	B	433	Total	C	N	O	S	Se	0	0	0
			3337	2141	571	618	2	5			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	GB 14591090
A	90	MSE	MET	modified residue	GB 14591090
A	109	MSE	MET	modified residue	GB 14591090
A	286	MSE	MET	modified residue	GB 14591090
A	351	MSE	MET	modified residue	GB 14591090
A	393	MSE	MET	modified residue	GB 14591090
B	1	MSE	MET	modified residue	GB 14591090
B	90	MSE	MET	modified residue	GB 14591090
B	109	MSE	MET	modified residue	GB 14591090
B	286	MSE	MET	modified residue	GB 14591090
B	351	MSE	MET	modified residue	GB 14591090
B	393	MSE	MET	modified residue	GB 14591090

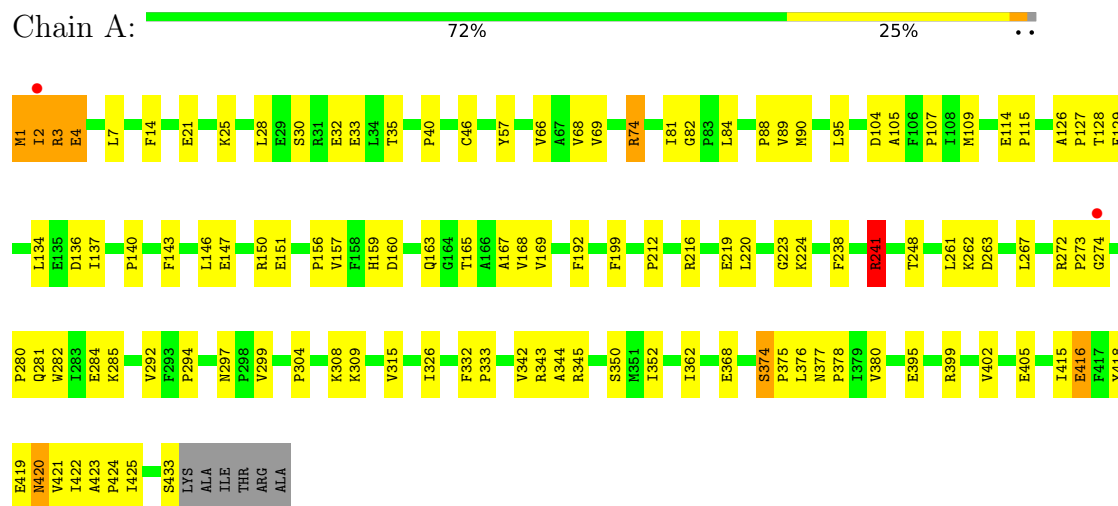
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	121	Total	O	0	0
			121	121		
2	B	111	Total	O	0	0
			111	111		

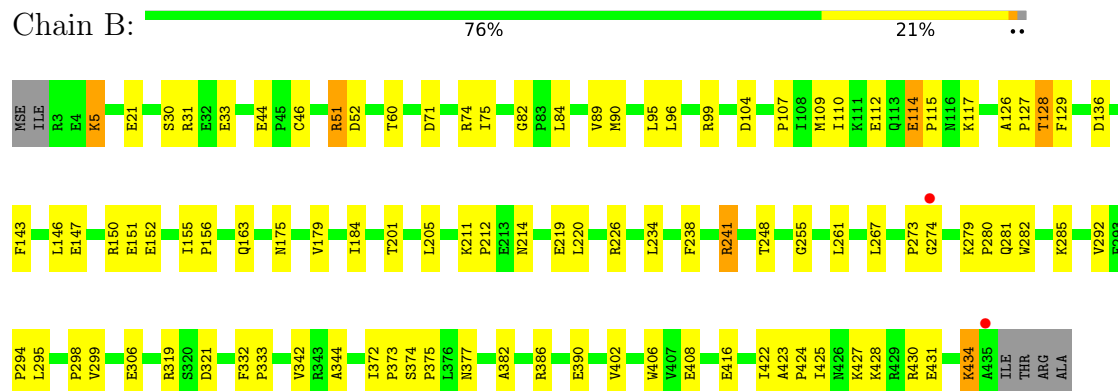
3 Residue-property plots

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: malate oxidoreductase



- Molecule 1: malate oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.46Å 85.68Å 172.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.4 (20.00-2.50) 98.3 (20.00-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.28 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.190 , 0.240 0.187 , 0.240	Depositor DCC
R_{free} test set	1474 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 31.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6908	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% 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	0.48	2/3398 (0.1%)	0.98	16/4595 (0.3%)
1	B	0.46	0/3396	1.03	20/4592 (0.4%)
All	All	0.47	2/6794 (0.0%)	1.01	36/9187 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	MSE	CG-SE	-5.13	1.80	1.95
1	A	1	MSE	SE-CE	-5.02	1.80	1.95

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	75	ILE	O-C-N	10.26	134.33	122.83
1	B	75	ILE	CA-C-O	-10.08	108.85	120.74
1	B	241	ARG	CD-NE-CZ	-8.13	113.02	124.40
1	B	84	LEU	N-CA-CB	8.05	122.05	110.13
1	B	184	ILE	N-CA-C	6.88	118.46	110.62
1	B	163	GLN	N-CA-C	6.78	118.67	111.28
1	B	5	LYS	N-CA-CB	-6.72	99.87	110.28
1	B	427	LYS	N-CA-C	-6.51	103.87	110.97
1	B	295	LEU	N-CA-C	6.51	120.22	112.93
1	A	163	GLN	N-CA-C	6.30	118.23	111.36
1	B	152	GLU	N-CA-C	6.13	120.83	113.23
1	B	82	GLY	CA-C-N	6.00	125.88	119.28
1	B	82	GLY	C-N-CA	6.00	125.88	119.28
1	B	129	PHE	N-CA-C	5.99	119.70	110.42
1	A	377	ASN	CA-C-N	5.90	125.77	119.28
1	A	377	ASN	C-N-CA	5.90	125.77	119.28
1	B	84	LEU	CB-CG-CD2	5.86	128.27	110.70
1	A	374	SER	CA-C-N	5.74	125.42	119.56
1	A	374	SER	C-N-CA	5.74	125.42	119.56
1	A	69	VAL	N-CA-C	5.55	116.58	108.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	PHE	N-CA-C	5.43	118.56	110.14
1	B	234	LEU	N-CA-C	5.38	117.84	111.33
1	B	99	ARG	N-CA-C	5.30	117.14	111.36
1	A	160	ASP	N-CA-C	5.27	117.44	111.11
1	A	81	ILE	N-CA-C	5.16	117.18	112.43
1	A	82	GLY	CA-C-N	5.08	124.58	119.19
1	A	82	GLY	C-N-CA	5.08	124.58	119.19
1	A	421	VAL	N-CA-C	5.06	116.00	111.90
1	A	241	ARG	CD-NE-CZ	-5.05	117.33	124.40
1	A	74	ARG	CD-NE-CZ	-5.04	117.34	124.40
1	A	68	VAL	N-CA-C	-5.04	99.87	107.73
1	A	3	ARG	CD-NE-CZ	-5.03	117.36	124.40
1	B	377	ASN	CA-C-N	5.03	124.63	119.05
1	B	377	ASN	C-N-CA	5.03	124.63	119.05
1	B	114	GLU	CA-C-N	5.00	124.44	119.24
1	B	114	GLU	C-N-CA	5.00	124.44	119.24

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	3339	0	3448	85	0
1	B	3337	0	3443	66	0
2	A	121	0	0	3	0
2	B	111	0	0	1	0
All	All	6908	0	6891	139	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 10.

All (139) 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:292:VAL:HG12	1:B:294:PRO:HD3	1.48	0.95
1:A:261:LEU:HD21	1:A:267:LEU:HD13	1.55	0.86
1:A:422:ILE:HA	1:A:425:ILE:HD12	1.57	0.84
1:A:280:PRO:HG2	1:A:281:GLN:NE2	2.00	0.76
1:A:74:ARG:O	1:A:136:ASP:HB3	1.88	0.73
1:A:168:VAL:HG12	1:A:326:ILE:HG23	1.70	0.72
1:A:292:VAL:HG12	1:A:294:PRO:HD3	1.70	0.72
1:B:319:ARG:HG2	1:B:319:ARG:HH11	1.55	0.71
1:B:422:ILE:HA	1:B:425:ILE:HD12	1.75	0.68
1:A:285:LYS:HA	1:A:285:LYS:HE2	1.75	0.68
1:B:219:GLU:HG2	1:B:238:PHE:CE1	2.29	0.67
1:B:280:PRO:HG3	1:B:306:GLU:HG3	1.76	0.66
1:B:143:PHE:CE2	1:B:241:ARG:HD3	2.31	0.66
1:B:274:GLY:H	1:B:299:VAL:HB	1.60	0.65
1:A:30:SER:OG	1:A:33:GLU:HG3	1.97	0.64
1:B:279:LYS:HE2	1:B:282:TRP:CZ2	2.33	0.64
1:A:1:MSE:HG2	1:A:4:GLU:HG2	1.80	0.63
1:A:89:VAL:HG23	1:B:46:CYS:SG	2.40	0.62
1:A:219:GLU:HG3	1:A:220:LEU:H	1.64	0.62
1:A:308:LYS:HD2	2:A:549:HOH:O	2.00	0.61
1:B:423:ALA:HB3	1:B:424:PRO:HD3	1.80	0.61
1:A:261:LEU:CD2	1:A:267:LEU:HD13	2.30	0.61
1:A:199:PHE:CE1	1:A:241:ARG:HB3	2.36	0.61
1:A:147:GLU:O	1:A:151:GLU:HG2	2.00	0.60
1:B:434:LYS:HE3	1:B:434:LYS:HA	1.83	0.60
1:A:405:GLU:HG3	2:A:456:HOH:O	2.01	0.60
1:A:1:MSE:CG	1:A:4:GLU:HG2	2.32	0.60
1:A:219:GLU:HG3	1:A:220:LEU:N	2.17	0.60
1:A:3:ARG:HA	1:B:44:GLU:HG3	1.83	0.60
1:B:428:LYS:HG3	1:B:431:GLU:OE1	2.01	0.59
1:A:261:LEU:HD12	1:A:282:TRP:HB3	1.83	0.59
1:B:143:PHE:O	1:B:147:GLU:HG2	2.03	0.58
1:B:430:ARG:CZ	1:B:434:LYS:HG2	2.34	0.58
1:A:342:VAL:HG12	1:A:402:VAL:HG22	1.85	0.58
1:A:423:ALA:HB3	1:A:424:PRO:HD3	1.85	0.57
1:B:372:ILE:HD12	1:B:373:PRO:HD2	1.86	0.57
1:A:212:PRO:O	1:A:248:THR:HA	2.05	0.56
1:B:95:LEU:HD12	1:B:96:LEU:N	2.19	0.56
1:B:31:ARG:HG2	1:B:31:ARG:HH11	1.71	0.56
1:A:150:ARG:HA	1:A:157:VAL:HG11	1.87	0.56
1:A:90:MSE:HG3	1:A:109:MSE:SE	2.56	0.56
1:B:402:VAL:HG21	1:B:406:TRP:CE3	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MSE:HG2	1:A:4:GLU:CG	2.34	0.55
1:B:212:PRO:O	1:B:248:THR:HA	2.05	0.55
1:A:415:ILE:O	1:A:419:GLU:HG2	2.08	0.54
1:A:46:CYS:SG	1:B:89:VAL:HG23	2.48	0.53
1:A:219:GLU:HG2	1:A:238:PHE:CE1	2.43	0.53
1:B:211:LYS:HB2	1:B:214:ASN:HD22	1.72	0.53
1:A:150:ARG:HA	1:A:157:VAL:CG1	2.39	0.53
1:B:319:ARG:HG2	1:B:319:ARG:NH1	2.22	0.53
1:A:107:PRO:O	1:B:128:THR:HG21	2.08	0.53
1:A:273:PRO:HB3	1:A:297:ASN:O	2.07	0.53
1:B:112:GLU:HG2	1:B:117:LYS:HB3	1.91	0.53
1:A:332:PHE:CG	1:A:333:PRO:HD3	2.44	0.53
1:B:31:ARG:HG2	1:B:31:ARG:NH1	2.24	0.52
1:B:342:VAL:HG12	1:B:402:VAL:HG12	1.90	0.52
1:A:156:PRO:HD3	1:A:344:ALA:O	2.08	0.52
1:A:140:PRO:HB3	1:A:241:ARG:NH2	2.24	0.51
1:A:90:MSE:HE1	1:A:136:ASP:OD2	2.11	0.51
1:A:223:GLY:O	1:A:224:LYS:HD3	2.10	0.51
1:B:226:ARG:HH21	1:B:255:GLY:HA2	1.75	0.51
1:A:14:PHE:H	1:A:433:SER:HB3	1.75	0.51
1:B:382:ALA:HB1	1:B:408:GLU:HG3	1.94	0.50
1:B:279:LYS:HE2	1:B:282:TRP:CE2	2.47	0.50
1:B:156:PRO:HD3	1:B:344:ALA:O	2.12	0.50
1:B:30:SER:OG	1:B:33:GLU:HG3	2.12	0.50
1:B:201:THR:O	1:B:205:LEU:HB2	2.12	0.50
1:B:126:ALA:N	1:B:127:PRO:CD	2.75	0.49
1:B:89:VAL:HB	1:B:90:MSE:HE2	1.95	0.49
1:A:90:MSE:CG	1:A:109:MSE:SE	3.11	0.49
1:B:281:GLN:H	1:B:281:GLN:CD	2.21	0.48
1:B:386:ARG:O	1:B:390:GLU:HG3	2.14	0.48
1:A:66:VAL:O	1:A:105:ALA:HA	2.12	0.48
1:A:272:ARG:NH1	2:A:557:HOH:O	2.46	0.48
1:A:146:LEU:O	1:A:150:ARG:HB2	2.13	0.48
1:B:90:MSE:HG3	1:B:109:MSE:SE	2.64	0.47
1:B:332:PHE:CG	1:B:333:PRO:HD3	2.49	0.47
1:B:155:ILE:HB	1:B:156:PRO:HD2	1.96	0.47
1:B:211:LYS:HB2	1:B:214:ASN:ND2	2.29	0.47
1:B:298:PRO:O	1:B:319:ARG:HD2	2.13	0.47
1:A:376:LEU:O	1:A:378:PRO:HD3	2.15	0.47
1:A:285:LYS:HE2	1:A:285:LYS:CA	2.43	0.47
1:B:146:LEU:O	1:B:150:ARG:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:GLU:CG	1:A:220:LEU:H	2.26	0.46
1:B:285:LYS:N	1:B:285:LYS:HD2	2.30	0.46
1:B:430:ARG:NH1	1:B:434:LYS:HG2	2.30	0.46
1:A:126:ALA:N	1:A:127:PRO:CD	2.79	0.46
1:A:167:ALA:HB2	1:A:352:ILE:HG23	1.98	0.46
1:B:95:LEU:HD12	1:B:95:LEU:C	2.40	0.46
1:A:280:PRO:HG2	1:A:281:GLN:HE21	1.79	0.46
1:B:51:ARG:HG3	1:B:52:ASP:N	2.31	0.46
1:A:21:GLU:HG3	1:B:104:ASP:OD1	2.16	0.45
1:A:280:PRO:O	1:A:284:GLU:HG3	2.16	0.45
1:A:32:GLU:CD	1:B:5:LYS:NZ	2.75	0.45
1:A:368:GLU:H	1:A:368:GLU:CD	2.25	0.45
1:A:374:SER:HA	1:A:375:PRO:HD3	1.83	0.45
1:A:350:SER:OG	1:A:395:GLU:OE2	2.33	0.45
1:B:219:GLU:HG3	1:B:220:LEU:N	2.32	0.45
1:A:274:GLY:H	1:A:299:VAL:HB	1.81	0.44
1:A:3:ARG:O	1:A:7:LEU:HG	2.18	0.44
1:A:345:ARG:HE	1:A:345:ARG:HB3	1.55	0.44
1:A:143:PHE:O	1:A:147:GLU:HG2	2.18	0.44
1:A:219:GLU:CG	1:A:220:LEU:N	2.80	0.44
1:A:343:ARG:HG2	1:A:399:ARG:HE	1.83	0.44
1:A:128:THR:HG21	1:B:107:PRO:O	2.17	0.44
1:A:28:LEU:HD11	1:B:95:LEU:HD13	1.99	0.43
1:B:112:GLU:HG2	1:B:117:LYS:CB	2.47	0.43
1:A:25:LYS:HD3	1:B:21:GLU:HB2	1.99	0.43
1:A:304:PRO:HA	1:A:315:VAL:HG11	2.01	0.43
1:A:422:ILE:HA	1:A:425:ILE:CD1	2.38	0.43
1:A:28:LEU:CD1	1:B:95:LEU:HD13	2.48	0.42
1:A:192:PHE:CD1	1:A:192:PHE:N	2.87	0.42
1:B:226:ARG:HD3	2:B:446:HOH:O	2.19	0.42
1:A:416:GLU:O	1:A:420:ASN:HB2	2.18	0.42
1:B:319:ARG:HB3	1:B:321:ASP:OD1	2.18	0.42
1:B:74:ARG:O	1:B:136:ASP:HB3	2.20	0.42
1:B:261:LEU:HD21	1:B:267:LEU:HD13	2.01	0.42
1:A:1:MSE:HG3	1:A:4:GLU:HG2	2.03	0.41
1:A:165:THR:O	1:A:169:VAL:HG23	2.19	0.41
1:B:114:GLU:HA	1:B:115:PRO:HD2	1.92	0.41
1:A:134:LEU:HB3	1:A:137:ILE:HD12	2.03	0.41
1:A:423:ALA:N	1:A:424:PRO:CD	2.83	0.41
1:A:418:TYR:CD2	1:A:418:TYR:C	2.98	0.41
1:B:71:ASP:HA	1:B:110:ILE:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:423:ALA:N	1:B:424:PRO:CD	2.84	0.41
1:A:362:ILE:HG13	1:A:380:VAL:HG22	2.02	0.41
1:A:57:TYR:CE2	1:A:343:ARG:HD3	2.56	0.41
1:A:134:LEU:HB3	1:A:137:ILE:CD1	2.51	0.41
1:A:134:LEU:O	1:A:159:HIS:HD2	2.04	0.41
1:A:262:LYS:O	1:A:263:ASP:HB2	2.20	0.41
1:B:175:ASN:O	1:B:179:VAL:HG23	2.21	0.41
1:A:2:ILE:HG23	1:A:3:ARG:N	2.36	0.41
1:A:88:PRO:HG3	1:B:60:THR:HG22	2.03	0.41
1:A:309:LYS:HB2	1:A:309:LYS:HE3	1.88	0.40
1:B:374:SER:HA	1:B:375:PRO:HD3	1.92	0.40
1:A:35:THR:HB	1:A:40:PRO:HD2	2.04	0.40
1:A:104:ASP:OD1	1:B:21:GLU:HG3	2.22	0.40
1:A:192:PHE:N	1:A:192:PHE:HD1	2.19	0.40
1:A:114:GLU:HA	1:A:115:PRO:HD2	1.88	0.40

There are no symmetry-related clashes.

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	431/439 (98%)	412 (96%)	19 (4%)	0	100	100
1	B	431/439 (98%)	410 (95%)	20 (5%)	1 (0%)	43	63
All	All	862/878 (98%)	822 (95%)	39 (4%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	273	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	354/352 (101%)	346 (98%)	8 (2%)	44	72
1	B	353/352 (100%)	348 (99%)	5 (1%)	59	81
All	All	707/704 (100%)	694 (98%)	13 (2%)	51	77

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

Mol	Chain	Res	Type
1	A	2	ILE
1	A	4	GLU
1	A	84	LEU
1	A	95	LEU
1	A	216	ARG
1	A	241	ARG
1	A	416	GLU
1	A	420	ASN
1	B	51	ARG
1	B	128	THR
1	B	151	GLU
1	B	416	GLU
1	B	434	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	ASN
1	A	133	ASN
1	A	214	ASN
1	A	258	GLN
1	B	116	ASN
1	B	214	ASN
1	B	252	ASN
1	B	258	GLN
1	B	370	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 [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	427/439 (97%)	-0.44	2 (0%) 87 85	18, 32, 51, 63	0
1	B	428/439 (97%)	-0.44	2 (0%) 87 85	18, 32, 51, 60	0
All	All	855/878 (97%)	-0.44	4 (0%) 87 85	18, 32, 51, 63	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	274	GLY	3.6
1	A	2	ILE	2.5
1	A	274	GLY	2.3
1	B	435	ALA	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.