



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:45 PM UTC

PDB ID : 1C3C / pdb_00001c3c
Title : T. MARITIMA ADENYLOSUCCINATE LYASE
Authors : Toth, E.A.; Yeates, T.O.
Deposited on : 1999-07-27
Resolution : 1.80 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

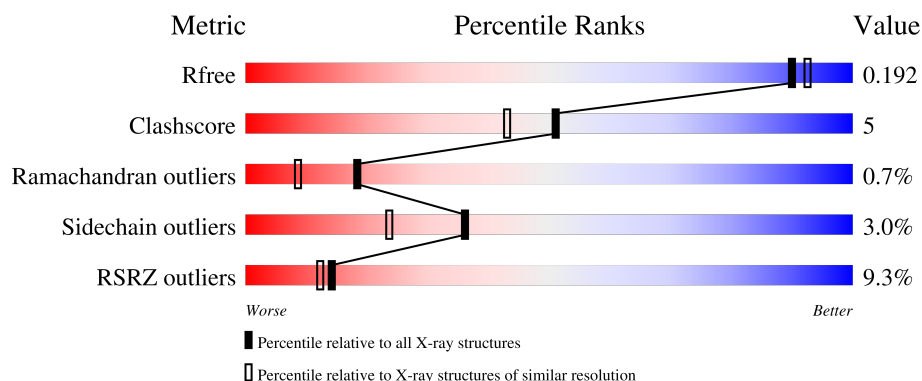
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	429	8% 83% 13% ..
1	B	429	10% 83% 13% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7427 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 PROTEIN (ADENYLOSUCCINATE LYASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	0	0
			3367	2154	577	624	12			
1	B	424	Total	C	N	O	S	0	0	0
			3367	2154	577	624	12			

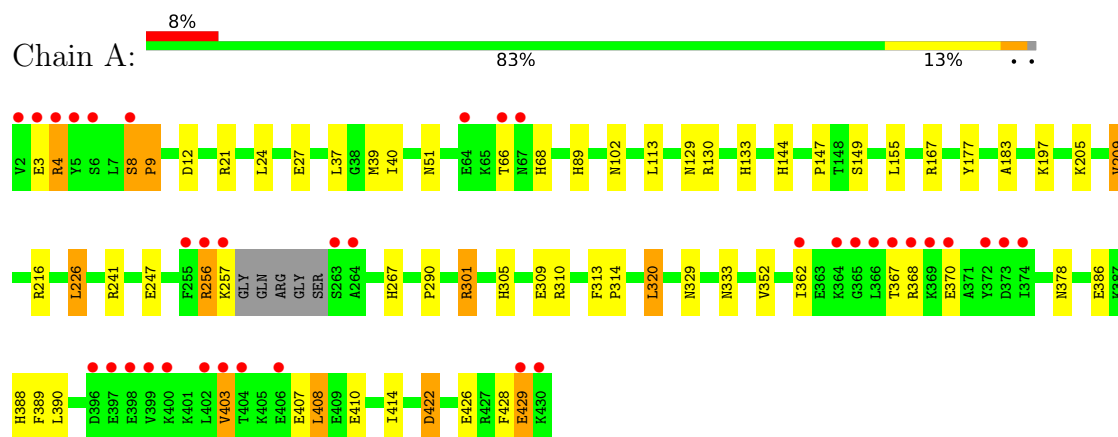
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	347	Total	O	0	0
			347	347		
2	B	346	Total	O	0	0
			346	346		

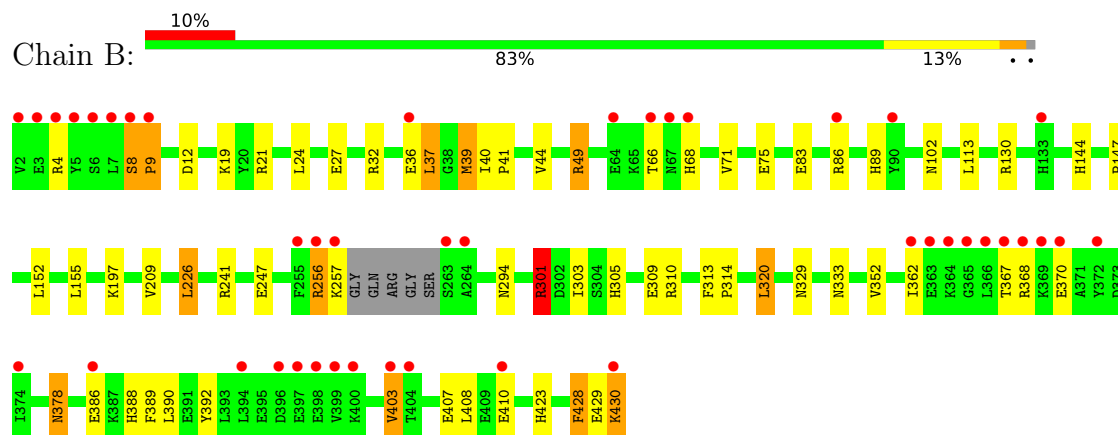
3 Residue-property plots [i](#)

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: PROTEIN (ADENYLOSUCCINATE LYASE)



• Molecule 1: PROTEIN (ADENYLOSUCCINATE LYASE)



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	120.81Å 126.75Å 169.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.80 50.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.8 (50.00-1.80) 97.9 (50.00-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.07 (at 1.79Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.179 , 0.207 0.166 , 0.192	Depositor DCC
R_{free} test set	5858 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	15.5	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.024 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7427	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% 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	0.86	2/3436 (0.1%)	1.61	42/4659 (0.9%)
1	B	0.90	1/3436 (0.0%)	1.61	41/4659 (0.9%)
All	All	0.88	3/6872 (0.0%)	1.61	83/9318 (0.9%)

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	3
1	B	0	1
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	39	MET	SD-CE	-10.04	1.54	1.79
1	A	429	GLU	CA-C	6.14	1.61	1.53
1	A	429	GLU	C-N	-5.70	1.25	1.33

All (83) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	429	GLU	CA-C-N	19.31	156.46	121.70
1	A	429	GLU	C-N-CA	19.31	156.46	121.70
1	A	21	ARG	CD-NE-CZ	19.17	151.24	124.40
1	B	310	ARG	CD-NE-CZ	16.09	146.93	124.40
1	B	21	ARG	CD-NE-CZ	15.98	146.77	124.40
1	A	310	ARG	CD-NE-CZ	14.33	144.47	124.40
1	B	378	ASN	OD1-CG-ND2	-14.31	108.29	122.60
1	A	8	SER	CA-C-N	13.89	137.21	119.84
1	A	8	SER	C-N-CA	13.89	137.21	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	8	SER	CA-C-N	13.89	137.21	119.84
1	B	8	SER	C-N-CA	13.89	137.21	119.84
1	A	378	ASN	OD1-CG-ND2	-13.19	109.41	122.60
1	A	429	GLU	O-C-N	-13.15	106.47	122.65
1	B	429	GLU	CA-C-N	12.54	144.27	121.70
1	B	429	GLU	C-N-CA	12.54	144.27	121.70
1	A	209	VAL	CB-CA-C	-10.54	95.79	110.12
1	A	8	SER	O-C-N	9.38	132.11	121.32
1	A	8	SER	CA-C-O	-9.09	107.71	120.16
1	A	9	PRO	CA-N-CD	-8.96	99.45	112.00
1	B	209	VAL	CB-CA-C	-8.88	98.04	110.12
1	A	389	PHE	CA-CB-CG	8.82	122.62	113.80
1	B	9	PRO	CA-N-CD	-8.50	100.10	112.00
1	A	428	PHE	CA-CB-CG	-8.02	105.78	113.80
1	B	8	SER	O-C-N	7.77	130.26	121.32
1	B	12	ASP	CA-CB-CG	-7.71	104.89	112.60
1	A	9	PRO	N-CA-C	7.67	128.27	112.47
1	B	241	ARG	NE-CZ-NH1	-7.52	113.98	121.50
1	A	422	ASP	CB-CG-OD2	7.40	135.42	118.40
1	A	408	LEU	CD1-CG-CD2	-7.33	94.67	110.80
1	B	386	GLU	CA-CB-CG	-7.32	99.47	114.10
1	B	9	PRO	N-CA-CB	7.26	110.88	103.25
1	B	21	ARG	NE-CZ-NH1	-7.19	114.31	121.50
1	B	408	LEU	CD1-CG-CD2	-7.11	95.16	110.80
1	B	389	PHE	CA-CB-CG	7.00	120.80	113.80
1	A	9	PRO	N-CA-CB	6.93	110.52	103.25
1	A	130	ARG	NE-CZ-NH1	-6.56	114.94	121.50
1	A	310	ARG	NE-CZ-NH2	-6.48	113.37	119.20
1	B	21	ARG	NE-CZ-NH2	6.48	125.03	119.20
1	B	386	GLU	OE1-CD-OE2	6.42	138.31	122.90
1	B	310	ARG	NE-CZ-NH2	-6.40	113.44	119.20
1	B	430	LYS	CA-C-O	-6.39	109.93	120.80
1	B	39	MET	CG-SD-CE	6.25	114.65	100.90
1	A	12	ASP	CA-CB-CG	-6.22	106.38	112.60
1	A	426	GLU	N-CA-CB	-6.06	100.89	110.22
1	B	130	ARG	NE-CZ-NH1	-6.05	115.45	121.50
1	A	241	ARG	NE-CZ-NH1	-6.03	115.47	121.50
1	B	378	ASN	CA-CB-CG	-5.96	106.64	112.60
1	B	386	GLU	CG-CD-OE2	-5.87	104.90	118.40
1	A	21	ARG	NE-CZ-NH2	5.86	124.47	119.20
1	A	183	ALA	CA-C-O	-5.83	114.66	120.90
1	B	44	VAL	CA-C-O	-5.83	114.99	121.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	378	ASN	CA-CB-CG	-5.76	106.84	112.60
1	A	4	ARG	CA-C-O	-5.73	113.78	120.20
1	B	241	ARG	NE-CZ-NH2	5.68	124.31	119.20
1	A	216	ARG	NE-CZ-NH1	-5.62	115.88	121.50
1	B	44	VAL	O-C-N	5.54	127.33	121.91
1	B	378	ASN	CB-CG-OD1	5.52	131.85	120.80
1	A	24	LEU	CD1-CG-CD2	-5.51	98.68	110.80
1	A	209	VAL	CA-CB-CG2	5.51	119.76	110.40
1	A	51	ASN	OD1-CG-ND2	5.49	128.09	122.60
1	B	310	ARG	NE-CZ-NH1	5.47	126.97	121.50
1	B	428	PHE	CA-CB-CG	-5.46	108.34	113.80
1	A	3	GLU	CA-C-O	5.40	126.46	120.63
1	B	423	HIS	O-C-N	5.38	127.83	122.12
1	A	102	ASN	CA-CB-CG	-5.38	107.22	112.60
1	A	9	PRO	O-C-N	5.35	129.86	122.64
1	B	152	LEU	CD1-CG-CD2	-5.33	99.08	110.80
1	B	423	HIS	CA-CB-CG	5.31	119.11	113.80
1	A	290	PRO	N-CA-C	-5.29	106.64	113.57
1	A	428	PHE	CB-CA-C	-5.24	101.32	110.17
1	B	49	ARG	CD-NE-CZ	-5.22	117.09	124.40
1	B	9	PRO	N-CA-C	5.21	123.21	112.47
1	B	301	ARG	N-CA-CB	5.17	119.23	110.49
1	A	155	LEU	CD1-CG-CD2	-5.17	99.43	110.80
1	A	167	ARG	NE-CZ-NH2	-5.13	114.58	119.20
1	A	426	GLU	OE1-CD-OE2	-5.12	110.60	122.90
1	B	294	ASN	CA-CB-CG	5.11	117.71	112.60
1	A	40	ILE	N-CA-CB	5.10	116.01	111.67
1	B	429	GLU	N-CA-C	-5.08	105.34	112.25
1	A	129	ASN	CA-CB-CG	-5.05	107.55	112.60
1	A	414	ILE	CA-CB-CG1	-5.05	101.82	110.40
1	B	8	SER	CA-C-O	-5.01	113.30	120.16
1	B	102	ASN	CA-CB-CG	-5.00	107.60	112.60

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	429	GLU	Peptide
1	A	8	SER	Peptide,Mainchain
1	B	8	SER	Peptide

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	3367	0	3311	34	2
1	B	3367	0	3312	38	2
2	A	347	0	0	9	0
2	B	346	0	0	6	0
All	All	7427	0	6623	70	2

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 5.

All (70) 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:39:MET:SD	2:A:775:HOH:O	2.27	0.93
1:A:301:ARG:HH21	1:A:305:HIS:CE1	2.06	0.73
1:A:39:MET:CG	2:A:775:HOH:O	2.36	0.73
1:B:301:ARG:HH21	1:B:305:HIS:CE1	2.08	0.72
1:A:301:ARG:HH21	1:A:305:HIS:HE1	1.38	0.70
1:B:301:ARG:HH21	1:B:305:HIS:HE1	1.39	0.70
1:A:329:ASN:ND2	1:A:333:ASN:HD22	1.92	0.67
1:B:329:ASN:ND2	1:B:333:ASN:HD22	1.94	0.66
1:A:27:GLU:OE2	1:A:89:HIS:HD2	1.80	0.64
1:B:27:GLU:OE2	1:B:89:HIS:HD2	1.81	0.62
1:B:37:LEU:HD22	1:B:197:LYS:HD2	1.81	0.62
1:A:247:GLU:OE1	1:B:144:HIS:HE1	1.84	0.60
1:A:329:ASN:HD21	1:A:333:ASN:HD22	1.50	0.60
1:B:226:LEU:HG	1:B:320:LEU:HD12	1.83	0.59
1:B:144:HIS:HD2	2:B:457:HOH:O	1.86	0.57
1:A:144:HIS:HD2	2:A:448:HOH:O	1.87	0.57
1:A:144:HIS:HE1	1:B:247:GLU:OE1	1.89	0.56
1:B:428:PHE:HA	1:B:430:LYS:HE3	1.87	0.56
1:A:329:ASN:ND2	1:A:333:ASN:ND2	2.53	0.55
1:A:66:THR:HG22	1:A:68:HIS:N	2.21	0.55
1:A:226:LEU:HG	1:A:320:LEU:HD12	1.88	0.55
1:B:367:THR:HG22	1:B:370:GLU:OE2	2.07	0.55
1:B:362:ILE:HD11	1:B:368:ARG:HA	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:367:THR:HG22	1:A:370:GLU:OE2	2.06	0.54
1:A:39:MET:HG2	2:A:775:HOH:O	2.06	0.53
1:A:66:THR:HG22	1:A:68:HIS:H	1.73	0.53
1:B:388:HIS:HD2	2:B:520:HOH:O	1.91	0.53
1:A:256:ARG:O	1:A:257:LYS:CB	2.56	0.53
1:B:256:ARG:O	1:B:257:LYS:CB	2.56	0.53
1:B:83:GLU:O	1:B:86:ARG:HG3	2.09	0.52
1:B:32:ARG:O	1:B:36:GLU:HG3	2.10	0.52
1:A:329:ASN:HD21	1:A:333:ASN:ND2	2.06	0.52
1:B:27:GLU:OE2	1:B:89:HIS:CD2	2.61	0.52
1:A:388:HIS:HD2	2:A:509:HOH:O	1.93	0.51
1:B:378:ASN:HD22	1:B:392:TYR:HB3	1.76	0.50
1:B:329:ASN:ND2	1:B:333:ASN:ND2	2.57	0.50
1:B:329:ASN:HD21	1:B:333:ASN:HD22	1.59	0.50
1:A:39:MET:HE1	2:A:641:HOH:O	2.12	0.48
1:A:410:GLU:OE2	2:A:585:HOH:O	2.20	0.48
1:A:27:GLU:OE2	1:A:89:HIS:CD2	2.65	0.47
1:A:267:HIS:H	1:A:267:HIS:HD1	1.62	0.47
1:A:403:VAL:HG23	1:A:407:GLU:HB2	1.96	0.47
1:B:197:LYS:HG3	2:B:774:HOH:O	2.13	0.47
1:B:313:PHE:HB2	1:B:314:PRO:HD3	1.96	0.47
1:B:403:VAL:HG23	1:B:407:GLU:HB2	1.96	0.47
1:B:147:PRO:HD3	1:B:352:VAL:CG1	2.44	0.47
1:A:407:GLU:O	1:A:410:GLU:HG2	2.15	0.47
1:B:71:VAL:O	1:B:75:GLU:HG3	2.15	0.47
1:B:89:HIS:HE1	2:B:573:HOH:O	1.98	0.47
1:A:89:HIS:HE1	2:A:562:HOH:O	1.98	0.47
1:A:147:PRO:HD3	1:A:352:VAL:CG1	2.44	0.47
1:B:37:LEU:HD13	1:B:197:LYS:HE2	1.97	0.46
1:A:362:ILE:HD11	1:A:368:ARG:HA	1.97	0.46
1:B:49:ARG:HH11	1:B:49:ARG:HD2	1.57	0.45
1:B:40:ILE:HB	1:B:41:PRO:HD2	1.99	0.45
1:B:40:ILE:HB	1:B:41:PRO:CD	2.46	0.45
1:A:305:HIS:HD2	1:A:309:GLU:OE1	1.99	0.44
1:B:66:THR:HG22	1:B:68:HIS:H	1.82	0.44
1:B:305:HIS:HD2	1:B:309:GLU:OE1	2.00	0.44
1:A:313:PHE:HB2	1:A:314:PRO:HD3	1.99	0.44
1:B:66:THR:HG22	1:B:68:HIS:N	2.32	0.44
1:B:39:MET:HE1	2:B:652:HOH:O	2.17	0.42
1:B:19:LYS:HB3	1:B:19:LYS:HE3	1.80	0.42
1:B:37:LEU:HG	1:B:39:MET:HE3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:HIS:HE1	1:A:422:ASP:OD1	2.01	0.42
1:A:197:LYS:HG3	2:A:766:HOH:O	2.19	0.41
1:B:407:GLU:O	1:B:410:GLU:HG2	2.20	0.41
1:B:86:ARG:NH2	2:B:534:HOH:O	2.53	0.41
1:A:367:THR:HG22	1:A:370:GLU:HG3	2.03	0.41
1:A:177:TYR:CZ	1:A:205:LYS:HE2	2.56	0.40

All (2) 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:410:GLU:OE1	1:B:36:GLU:OE2[5_455]	1.90	0.30
1:A:410:GLU:OE2	1:B:36:GLU:OE2[5_455]	2.19	0.01

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	420/429 (98%)	410 (98%)	7 (2%)	3 (1%)	18	8
1	B	420/429 (98%)	409 (97%)	8 (2%)	3 (1%)	18	8
All	All	840/858 (98%)	819 (98%)	15 (2%)	6 (1%)	18	8

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	PRO
1	A	256	ARG
1	A	301	ARG
1	B	9	PRO
1	B	256	ARG
1	B	301	ARG

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	355/382 (93%)	344 (97%)	11 (3%)	35	23
1	B	355/382 (93%)	345 (97%)	10 (3%)	38	26
All	All	710/764 (93%)	689 (97%)	21 (3%)	36	24

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	37	LEU
1	A	113	LEU
1	A	149	SER
1	A	209	VAL
1	A	226	LEU
1	A	320	LEU
1	A	386	GLU
1	A	390	LEU
1	A	403	VAL
1	A	408	LEU
1	B	4	ARG
1	B	24	LEU
1	B	37	LEU
1	B	113	LEU
1	B	155	LEU
1	B	226	LEU
1	B	303	ILE
1	B	320	LEU
1	B	390	LEU
1	B	403	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	HIS

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Mol	Chain	Res	Type
1	A	89	HIS
1	A	133	HIS
1	A	144	HIS
1	A	189	ASN
1	A	270	ASN
1	A	305	HIS
1	A	329	ASN
1	A	388	HIS
1	B	68	HIS
1	B	89	HIS
1	B	144	HIS
1	B	189	ASN
1	B	270	ASN
1	B	305	HIS
1	B	329	ASN
1	B	378	ASN
1	B	388	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	424/429 (98%)	0.00	36 (8%) 16 14	8, 15, 45, 57	0
1	B	424/429 (98%)	0.05	43 (10%) 12 11	8, 15, 45, 57	0
All	All	848/858 (98%)	0.03	79 (9%) 14 12	8, 15, 45, 57	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	367	THR	7.5
1	B	430	LYS	6.4
1	B	366	LEU	6.3
1	B	256	ARG	6.1
1	A	366	LEU	5.6
1	A	263	SER	5.5
1	B	3	GLU	5.4
1	B	368	ARG	5.3
1	A	256	ARG	5.1
1	A	3	GLU	5.0
1	A	66	THR	4.8
1	A	367	THR	4.8
1	A	257	LYS	4.7
1	B	365	GLY	4.7
1	A	430	LYS	4.5
1	A	67	ASN	4.3
1	B	6	SER	4.3
1	A	429	GLU	4.1
1	B	257	LYS	4.0
1	B	403	VAL	4.0
1	A	399	VAL	4.0
1	A	404	THR	3.9
1	A	6	SER	3.9
1	B	263	SER	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	403	VAL	3.9
1	A	368	ARG	3.8
1	B	66	THR	3.8
1	A	365	GLY	3.7
1	A	255	PHE	3.6
1	B	86	ARG	3.6
1	B	36	GLU	3.5
1	B	370	GLU	3.4
1	A	2	VAL	3.4
1	B	397	GLU	3.4
1	B	398	GLU	3.2
1	B	404	THR	3.1
1	B	7	LEU	3.1
1	A	4	ARG	3.1
1	B	255	PHE	3.1
1	A	369	LYS	3.0
1	B	264	ALA	3.0
1	A	64	GLU	2.9
1	B	369	LYS	2.9
1	A	398	GLU	2.9
1	A	362	ILE	2.9
1	B	90	TYR	2.8
1	B	364	LYS	2.8
1	A	8	SER	2.8
1	B	374	ILE	2.8
1	B	399	VAL	2.8
1	A	5	TYR	2.8
1	B	362	ILE	2.8
1	B	9	PRO	2.8
1	B	68	HIS	2.7
1	B	386	GLU	2.7
1	B	396	ASP	2.7
1	B	64	GLU	2.7
1	B	4	ARG	2.7
1	B	400	LYS	2.7
1	B	2	VAL	2.6
1	A	402	LEU	2.6
1	B	67	ASN	2.6
1	B	8	SER	2.5
1	B	410	GLU	2.5
1	A	370	GLU	2.4
1	A	396	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	374	ILE	2.4
1	A	400	LYS	2.4
1	B	372	TYR	2.3
1	A	397	GLU	2.3
1	A	364	LYS	2.3
1	B	394	LEU	2.2
1	A	264	ALA	2.2
1	B	363	GLU	2.2
1	A	373	ASP	2.1
1	B	5	TYR	2.1
1	B	133	HIS	2.1
1	A	406	GLU	2.0
1	A	372	TYR	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.