



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:37 AM UTC

PDB ID : 8X0A / pdb_00008x0a
Title : Crystal Structure of MftR from mycobacterium tuberculosis
Authors : Peng, F.; Ke, Z.H.; Li, Y.
Deposited on : 2023-11-04
Resolution : 2.70 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

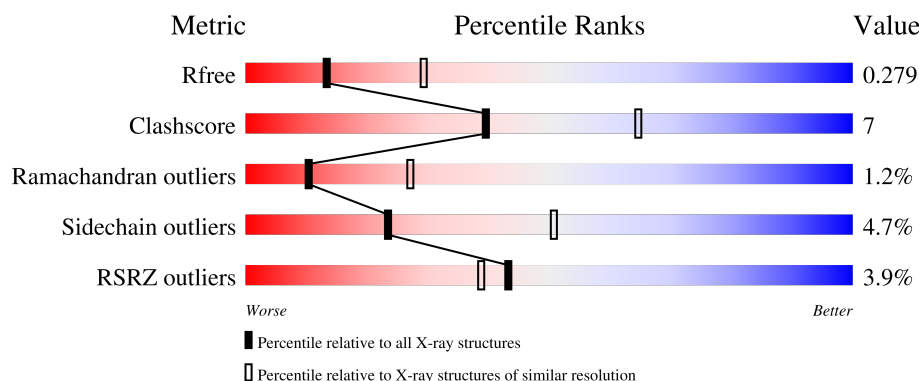
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	186	
1	B	186	
1	C	186	
1	D	186	
1	E	186	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7285 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 Putative mycofactocin biosynthesis transcriptional regulator MftR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	184	Total	C	N	O	S	0	0	0
			1451	920	261	265	5			
1	E	184	Total	C	N	O	S	0	0	0
			1451	920	261	265	5			
1	B	184	Total	C	N	O	S	0	0	0
			1451	920	261	265	5			
1	C	182	Total	C	N	O	S	0	0	0
			1432	908	256	263	5			
1	D	184	Total	C	N	O	S	0	0	0
			1451	920	261	265	5			

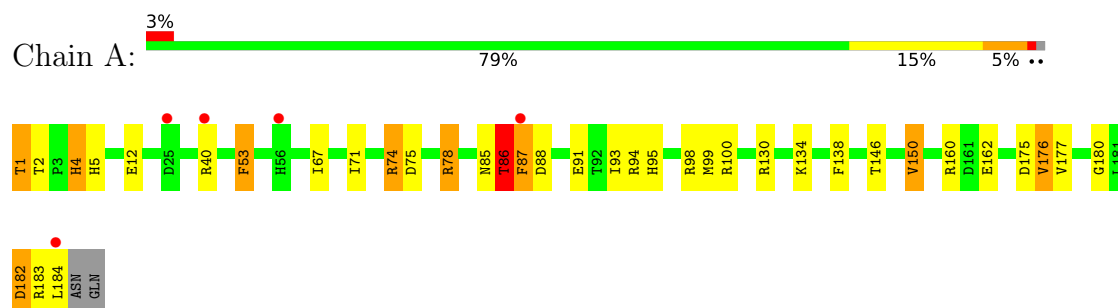
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	O	0	0
			10	10		
2	E	8	Total	O	0	0
			8	8		
2	B	12	Total	O	0	0
			12	12		
2	C	14	Total	O	0	0
			14	14		
2	D	5	Total	O	0	0
			5	5		

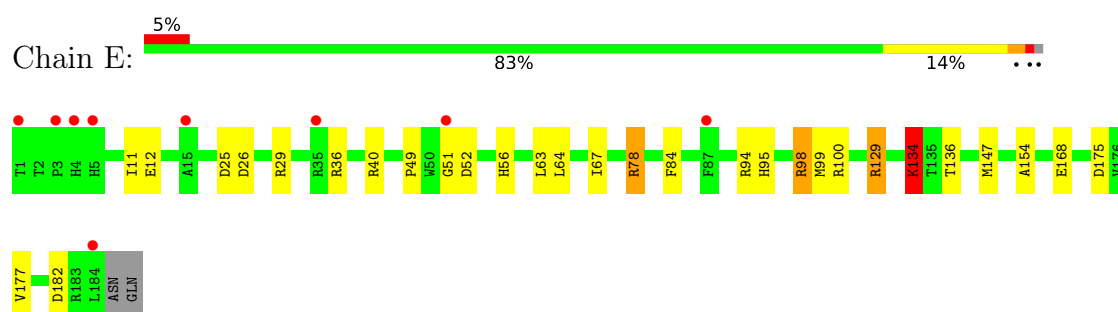
3 Residue-property plots [i](#)

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.

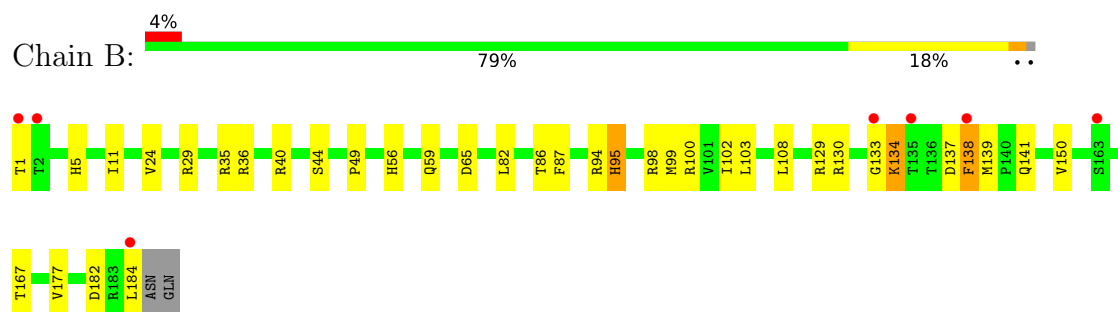
- Molecule 1: Putative mycofactocin biosynthesis transcriptional regulator MftR



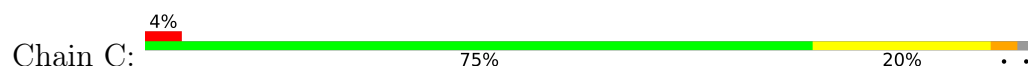
- Molecule 1: Putative mycofactocin biosynthesis transcriptional regulator MftR

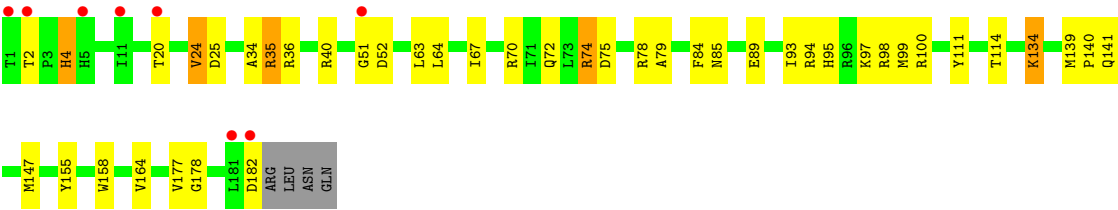


- Molecule 1: Putative mycofactocin biosynthesis transcriptional regulator MftR

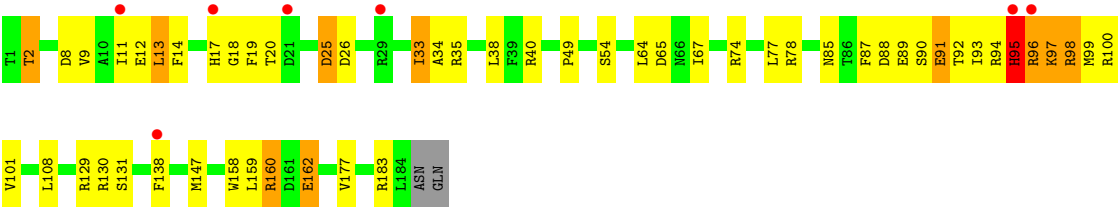


- Molecule 1: Putative mycofactocin biosynthesis transcriptional regulator MftR





● Molecule 1: Putative mycofactocin biosynthesis transcriptional regulator MftR



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.75Å 81.43Å 120.46Å 90.00° 107.36° 90.00°	Depositor
Resolution (Å)	47.38 – 2.70 47.38 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.4 (47.38-2.70) 99.4 (47.38-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0350	Depositor
R, R_{free}	0.229 , 0.278 0.230 , 0.279	Depositor DCC
R_{free} test set	1380 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	51.9	Xtriage
Anisotropy	0.223	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7285	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 4.95% 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	0.83	3/1482 (0.2%)	1.56	16/2010 (0.8%)
1	B	0.81	3/1482 (0.2%)	1.30	2/2010 (0.1%)
1	C	0.76	0/1463	1.43	10/1985 (0.5%)
1	D	0.79	1/1482 (0.1%)	1.35	6/2010 (0.3%)
1	E	0.75	0/1482	1.33	7/2010 (0.3%)
All	All	0.79	7/7391 (0.1%)	1.40	41/10025 (0.4%)

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	10
1	B	0	5
1	C	0	6
1	D	0	6
1	E	0	8
All	All	0	35

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	56	HIS	CE1-NE2	6.34	1.38	1.32
1	D	96	ARG	N-CA	6.21	1.52	1.46
1	A	5	HIS	CE1-NE2	5.95	1.38	1.32
1	B	95	HIS	CE1-NE2	5.80	1.38	1.32
1	A	87	PHE	C-N	5.27	1.40	1.33
1	A	4	HIS	CE1-NE2	5.03	1.37	1.32
1	B	5	HIS	CE1-NE2	5.03	1.37	1.32

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	ASP	N-CA-CB	-23.49	76.85	110.44
1	A	87	PHE	CA-C-O	-14.85	105.62	121.36
1	C	140	PRO	CB-CA-C	14.66	135.76	111.56
1	C	141	GLN	N-CA-C	10.76	123.08	111.36
1	C	141	GLN	N-CA-CB	-10.13	95.16	110.16
1	D	96	ARG	N-CA-C	9.67	123.04	109.31
1	A	87	PHE	CA-CB-CG	9.62	123.42	113.80
1	C	51	GLY	N-CA-C	8.69	133.77	113.18
1	C	36	ARG	CB-CG-CD	8.48	130.82	111.30
1	A	87	PHE	N-CA-C	8.41	120.34	108.74
1	C	52	ASP	N-CA-CB	8.38	124.66	110.49
1	D	131	SER	CA-C-N	-7.33	114.14	122.42
1	D	131	SER	C-N-CA	-7.33	114.14	122.42
1	E	52	ASP	N-CA-CB	7.15	123.17	110.79
1	B	182	ASP	CA-CB-CG	6.90	119.50	112.60
1	B	59	GLN	CB-CA-C	6.38	121.70	110.85
1	E	51	GLY	N-CA-C	6.29	128.08	113.18
1	A	138	PHE	CA-CB-CG	6.28	120.08	113.80
1	A	87	PHE	N-CA-CB	-6.19	100.87	111.21
1	D	96	ARG	CB-CA-C	-6.13	100.01	114.16
1	C	52	ASP	N-CA-C	-6.11	97.78	110.80
1	A	182	ASP	CA-CB-CG	6.06	118.66	112.60
1	E	134	LYS	CA-C-N	5.98	128.57	120.38
1	E	134	LYS	C-N-CA	5.98	128.57	120.38
1	E	52	ASP	N-CA-C	-5.78	98.31	109.24
1	C	36	ARG	NE-CZ-NH2	5.70	124.33	119.20
1	A	53	PHE	N-CA-CB	5.58	119.91	110.49
1	E	12	GLU	CB-CG-CD	5.40	121.78	112.60
1	C	134	LYS	CB-CA-C	5.40	118.29	110.79
1	D	2	THR	CA-C-O	-5.36	112.82	120.16
1	E	182	ASP	CA-CB-CG	5.27	117.87	112.60
1	A	1	THR	CA-CB-OG1	-5.26	101.71	109.60
1	A	134	LYS	CA-C-N	5.25	127.58	120.38
1	A	134	LYS	C-N-CA	5.25	127.58	120.38
1	A	86	THR	CA-C-O	-5.24	114.98	121.11
1	A	180	GLY	CA-C-O	-5.24	118.82	122.22
1	A	176	VAL	N-CA-CB	5.23	116.67	110.55
1	C	89	GLU	CB-CG-CD	5.20	121.44	112.60
1	A	12	GLU	CB-CG-CD	5.13	121.32	112.60
1	D	131	SER	N-CA-C	5.12	119.99	113.18
1	A	2	THR	N-CA-CB	5.05	115.18	110.39

There are no chirality outliers.

All (35) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	100	ARG	Sidechain
1	A	130	ARG	Sidechain
1	A	160	ARG	Sidechain
1	A	40	ARG	Sidechain
1	A	74	ARG	Sidechain
1	A	78	ARG	Sidechain
1	A	85	ASN	Mainchain
1	A	86	THR	Mainchain
1	A	94	ARG	Sidechain
1	A	98	ARG	Sidechain
1	B	100	ARG	Sidechain
1	B	130	ARG	Sidechain
1	B	35	ARG	Sidechain
1	B	40	ARG	Sidechain
1	B	94	ARG	Sidechain
1	C	100	ARG	Sidechain
1	C	35	ARG	Sidechain
1	C	40	ARG	Sidechain
1	C	70	ARG	Sidechain
1	C	94	ARG	Sidechain
1	C	98	ARG	Sidechain
1	D	100	ARG	Sidechain
1	D	130	ARG	Sidechain
1	D	160	ARG	Sidechain
1	D	35	ARG	Sidechain
1	D	40	ARG	Sidechain
1	D	78	ARG	Sidechain
1	E	100	ARG	Sidechain
1	E	129	ARG	Sidechain
1	E	29	ARG	Sidechain
1	E	36	ARG	Sidechain
1	E	40	ARG	Sidechain
1	E	78	ARG	Sidechain
1	E	94	ARG	Sidechain
1	E	98	ARG	Sidechain

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	1451	0	1438	20	0
1	B	1451	0	1438	18	0
1	C	1432	0	1414	14	1
1	D	1451	0	1438	40	0
1	E	1451	0	1438	12	0
2	A	10	0	0	0	0
2	B	12	0	0	0	0
2	C	14	0	0	2	0
2	D	5	0	0	0	0
2	E	8	0	0	0	0
All	All	7285	0	7166	96	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 7.

All (96) 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:D:8:ASP:OD2	1:D:94:ARG:NH1	1.96	0.97
1:A:87:PHE:HB3	1:A:91:GLU:CG	2.02	0.89
1:A:87:PHE:HB3	1:A:91:GLU:HG2	1.63	0.80
1:A:183:ARG:NH1	1:D:183:ARG:NH2	2.33	0.76
1:E:147:MET:HA	1:E:147:MET:HE2	1.69	0.75
1:A:87:PHE:HB3	1:A:91:GLU:HG3	1.71	0.73
1:B:133:GLY:O	1:B:134:LYS:HB2	1.91	0.70
1:C:24:VAL:HG13	1:C:35:ARG:HH12	1.57	0.69
1:D:96:ARG:HG2	1:D:97:LYS:H	1.57	0.68
1:A:86:THR:C	1:A:87:PHE:CG	2.73	0.67
1:C:85:ASN:OD1	1:C:155:TYR:OH	2.13	0.65
1:D:8:ASP:CG	1:D:94:ARG:NH1	2.56	0.64
1:B:1:THR:O	1:B:1:THR:HG22	1.98	0.63
1:D:8:ASP:OD1	1:D:94:ARG:HD3	2.00	0.62
1:D:93:ILE:C	1:D:95:HIS:H	2.07	0.61
1:C:111:TYR:O	1:C:114:THR:OG1	2.17	0.61
1:D:95:HIS:HB3	1:D:158:TRP:HH2	1.64	0.61
1:D:96:ARG:HG3	1:D:158:TRP:CE2	2.35	0.61
1:C:2:THR:HG22	1:C:4:HIS:H	1.66	0.60
1:D:97:LYS:O	1:D:99:MET:HE2	2.03	0.59
1:C:34:ALA:HB1	2:C:210:HOH:O	2.03	0.59
1:D:96:ARG:HD2	1:D:162:GLU:HA	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:77:LEU:HD13	1:D:147:MET:HE1	1.83	0.58
1:D:96:ARG:CG	1:D:97:LYS:H	2.17	0.57
1:D:87:PHE:HB3	1:D:91:GLU:HB2	1.87	0.56
1:D:90:SER:HA	1:D:93:ILE:HD12	1.85	0.56
1:B:134:LYS:HB2	1:B:137:ASP:OD2	2.05	0.56
1:A:183:ARG:NH1	1:D:183:ARG:CZ	2.68	0.56
1:A:146:THR:HG23	1:A:177:VAL:HG21	1.89	0.55
1:D:88:ASP:OD1	1:D:89:GLU:N	2.41	0.54
1:D:85:ASN:HA	1:D:87:PHE:CZ	2.44	0.53
1:C:158:TRP:HA	1:C:164:VAL:CG2	2.39	0.52
1:D:95:HIS:HB3	1:D:158:TRP:CH2	2.44	0.52
1:D:97:LYS:C	1:D:98:ARG:O	2.52	0.52
1:B:65:ASP:HA	1:B:129:ARG:HH22	1.75	0.51
1:D:77:LEU:HD13	1:D:147:MET:CE	2.40	0.51
1:D:93:ILE:C	1:D:95:HIS:N	2.66	0.51
1:C:63:LEU:HD13	1:C:84:PHE:CE2	2.46	0.50
1:E:63:LEU:HD13	1:E:84:PHE:CE2	2.46	0.50
1:B:11:ILE:HD12	1:B:98:ARG:HA	1.92	0.50
1:B:82:LEU:O	1:B:86:THR:HG23	2.11	0.50
1:A:1:THR:HA	1:A:4:HIS:HB3	1.94	0.50
1:E:11:ILE:HD11	1:E:49:PRO:HB3	1.93	0.50
1:A:78:ARG:NH1	1:A:175:ASP:OD1	2.43	0.50
1:D:9:VAL:O	1:D:13:LEU:HG	2.11	0.49
1:A:146:THR:CG2	1:A:177:VAL:HG21	2.42	0.49
1:A:71:ILE:HD12	1:C:79:ALA:HB1	1.93	0.48
1:A:183:ARG:HH11	1:D:183:ARG:NH2	2.09	0.48
1:C:72:GLN:HB2	2:C:214:HOH:O	2.12	0.48
1:D:49:PRO:HG2	1:D:108:LEU:HD21	1.96	0.48
1:B:102:ILE:HA	1:B:108:LEU:HD23	1.96	0.48
1:D:65:ASP:HA	1:D:129:ARG:NH2	2.28	0.48
1:A:95:HIS:CE1	1:A:99:MET:HE3	2.49	0.47
1:B:49:PRO:HG2	1:B:108:LEU:HD21	1.97	0.47
1:B:95:HIS:CE1	1:B:99:MET:HE3	2.50	0.47
1:D:33:ILE:HD12	1:D:34:ALA:O	2.14	0.47
1:B:29:ARG:HH11	1:B:29:ARG:HG2	1.78	0.47
1:B:11:ILE:HD11	1:B:49:PRO:HB3	1.97	0.47
1:B:86:THR:HG22	1:B:167:THR:HG21	1.97	0.46
1:C:64:LEU:O	1:C:67:ILE:HG22	2.16	0.46
1:A:183:ARG:HH12	1:D:183:ARG:CZ	2.29	0.46
1:B:134:LYS:HD2	1:B:134:LYS:HA	1.66	0.46
1:E:11:ILE:HD12	1:E:98:ARG:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:ARG:HH22	1:C:78:ARG:NH1	2.14	0.45
1:E:78:ARG:NH1	1:E:175:ASP:OD1	2.49	0.45
1:D:14:PHE:CD1	1:D:19:PHE:HA	2.51	0.45
1:E:154:ALA:HA	1:B:138:PHE:CZ	2.52	0.45
1:C:95:HIS:CE1	1:C:99:MET:HE3	2.52	0.45
1:E:95:HIS:CE1	1:E:99:MET:HE3	2.51	0.45
1:D:96:ARG:HH11	1:D:162:GLU:HB2	1.82	0.44
1:E:64:LEU:O	1:E:129:ARG:NH2	2.51	0.44
1:D:33:ILE:CD1	1:D:38:LEU:HB2	2.47	0.44
1:A:150:VAL:HG21	1:A:177:VAL:HG11	2.00	0.44
1:D:25:ASP:O	1:D:26:ASP:C	2.61	0.44
1:C:74:ARG:HG3	1:C:75:ASP:N	2.31	0.43
1:D:160:ARG:HH11	1:D:160:ARG:HG3	1.83	0.43
1:E:56:HIS:NE2	1:E:84:PHE:HB3	2.34	0.43
1:D:99:MET:HE1	1:D:159:LEU:O	2.18	0.43
1:B:1:THR:O	1:B:1:THR:CG2	2.64	0.43
1:C:93:ILE:HG22	1:C:97:LYS:HE3	2.01	0.43
1:D:95:HIS:CB	1:D:158:TRP:HH2	2.31	0.43
1:D:9:VAL:HA	1:D:12:GLU:OE1	2.19	0.43
1:A:74:ARG:NH2	1:A:75:ASP:OD1	2.52	0.42
1:D:64:LEU:O	1:D:67:ILE:HG22	2.20	0.42
1:D:88:ASP:HB3	1:D:91:GLU:HG3	2.01	0.42
1:E:134:LYS:HG2	1:E:136:THR:HG22	2.01	0.42
1:A:93:ILE:HD11	1:B:29:ARG:O	2.20	0.41
1:D:18:GLY:O	1:D:19:PHE:C	2.63	0.41
1:A:162:GLU:O	1:A:162:GLU:HG2	2.21	0.41
1:E:64:LEU:O	1:E:67:ILE:HG22	2.21	0.41
1:E:25:ASP:O	1:E:26:ASP:C	2.64	0.41
1:B:138:PHE:HD1	1:B:138:PHE:HA	1.77	0.41
1:D:92:THR:O	1:D:96:ARG:N	2.41	0.41
1:A:183:ARG:C	1:A:184:LEU:HG	2.46	0.41
1:B:87:PHE:CE2	1:B:95:HIS:CE1	3.09	0.40
1:D:89:GLU:O	1:D:93:ILE:HG13	2.20	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:C:178:GLY:O	1:C:182:ASP:OD1[2_656]	1.84	0.36

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	182/186 (98%)	176 (97%)	4 (2%)	2 (1%)	11	29
1	B	182/186 (98%)	172 (94%)	6 (3%)	4 (2%)	5	14
1	C	180/186 (97%)	174 (97%)	6 (3%)	0	100	100
1	D	182/186 (98%)	170 (93%)	8 (4%)	4 (2%)	5	14
1	E	182/186 (98%)	176 (97%)	5 (3%)	1 (0%)	24	48
All	All	908/930 (98%)	868 (96%)	29 (3%)	11 (1%)	10	27

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	53	PHE
1	E	134	LYS
1	B	134	LYS
1	B	139	MET
1	D	98	ARG
1	B	141	GLN
1	A	182	ASP
1	D	2	THR
1	D	95	HIS
1	D	97	LYS
1	B	44	SER

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	148/150 (99%)	145 (98%)	3 (2%)	48	76
1	B	148/150 (99%)	141 (95%)	7 (5%)	23	51
1	C	146/150 (97%)	137 (94%)	9 (6%)	16	39
1	D	148/150 (99%)	134 (90%)	14 (10%)	8	20
1	E	148/150 (99%)	146 (99%)	2 (1%)	59	82
All	All	738/750 (98%)	703 (95%)	35 (5%)	23	51

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	ILE
1	A	150	VAL
1	A	176	VAL
1	E	168	GLU
1	E	177	VAL
1	B	24	VAL
1	B	36	ARG
1	B	103	LEU
1	B	138	PHE
1	B	150	VAL
1	B	177	VAL
1	B	184	LEU
1	C	4	HIS
1	C	20	THR
1	C	24	VAL
1	C	25	ASP
1	C	74	ARG
1	C	134	LYS
1	C	139	MET
1	C	147	MET
1	C	177	VAL
1	D	11	ILE
1	D	13	LEU
1	D	17	HIS
1	D	20	THR
1	D	25	ASP
1	D	33	ILE
1	D	54	SER
1	D	74	ARG
1	D	91	GLU
1	D	95	HIS

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Mol	Chain	Res	Type
1	D	101	VAL
1	D	138	PHE
1	D	162	GLU
1	D	177	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	56	HIS
1	A	61	GLN
1	A	109	GLN
1	A	141	GLN
1	E	4	HIS
1	B	5	HIS
1	C	66	ASN
1	C	141	GLN
1	D	61	GLN

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	184/186 (98%)	0.03	5 (2%) 56 53	25, 37, 63, 93	0
1	B	184/186 (98%)	0.08	7 (3%) 44 40	27, 41, 71, 94	0
1	C	182/186 (97%)	0.28	8 (4%) 39 35	29, 46, 69, 106	0
1	D	184/186 (98%)	0.41	7 (3%) 44 40	23, 46, 88, 120	0
1	E	184/186 (98%)	0.29	9 (4%) 35 31	23, 48, 81, 117	0
All	All	918/930 (98%)	0.22	36 (3%) 43 39	23, 43, 80, 120	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	51	GLY	4.2
1	B	184	LEU	4.2
1	A	87	PHE	4.1
1	D	95	HIS	3.5
1	C	1	THR	3.5
1	A	184	LEU	3.5
1	D	96	ARG	3.2
1	C	181	LEU	3.2
1	E	1	THR	3.2
1	E	87	PHE	3.0
1	E	4	HIS	2.9
1	A	25	ASP	2.9
1	C	5	HIS	2.9
1	A	40	ARG	2.8
1	E	184	LEU	2.7
1	D	11	ILE	2.6
1	C	182	ASP	2.6
1	D	29	ARG	2.5
1	E	51	GLY	2.5
1	E	15	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	163	SER	2.5
1	B	138	PHE	2.5
1	B	133	GLY	2.3
1	D	17	HIS	2.3
1	D	21	ASP	2.2
1	E	5	HIS	2.2
1	B	135	THR	2.1
1	E	35	ARG	2.1
1	C	2	THR	2.1
1	E	3	PRO	2.1
1	B	2	THR	2.1
1	D	138	PHE	2.1
1	B	1	THR	2.0
1	C	20	THR	2.0
1	A	56	HIS	2.0
1	C	11	ILE	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.