



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 02:03 pm BST

PDB ID : 4X0X
Title : The structure of AhpE from Mycobacterium tuberculosis revisited
Authors : Pallo, A.; Messens, J.
Deposited on : 2014-11-24
Resolution : 1.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

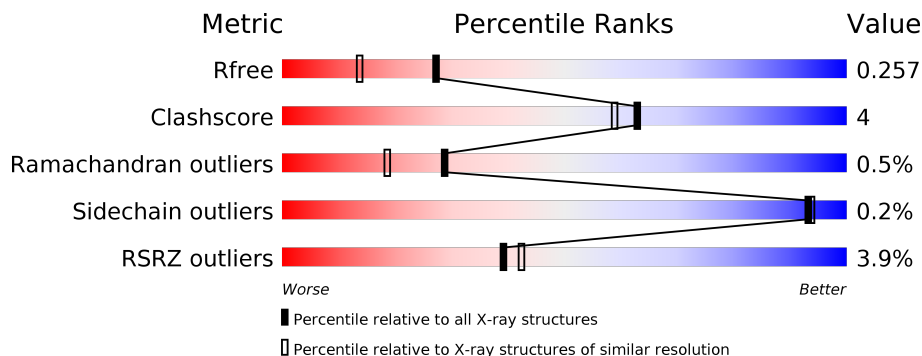
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 on 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	154	 5% 87% 11% ..
1	B	154	 2% 90% 9% .
1	C	154	 3% 93% 7%
1	D	154	 6% 88% 10% .

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 5105 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 peroxiredoxin MT2298.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	153	1198	765	208	222	3	0	2	0
1	B	154	1198	764	209	222	3	0	1	0
1	C	154	1198	764	209	222	3	0	1	0
1	D	154	1200	766	213	218	3	0	1	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	expression tag	UNP P9WIE2
B	0	HIS	-	expression tag	UNP P9WIE2
C	0	HIS	-	expression tag	UNP P9WIE2
D	0	HIS	-	expression tag	UNP P9WIE2

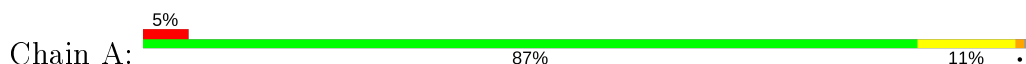
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	76	Total	O	0	0
			76	76		
2	B	86	Total	O	0	0
			86	86		
2	C	82	Total	O	0	0
			82	82		
2	D	67	Total	O	0	0
			67	67		

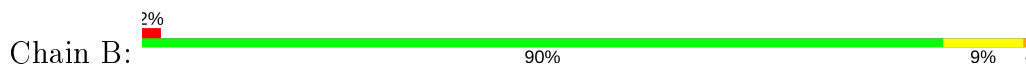
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

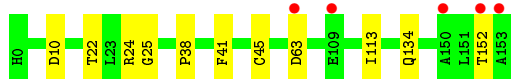
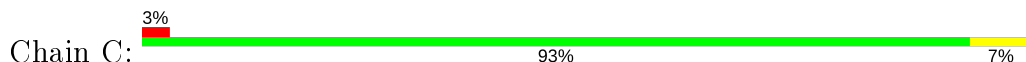
- Molecule 1: Putative peroxiredoxin MT2298



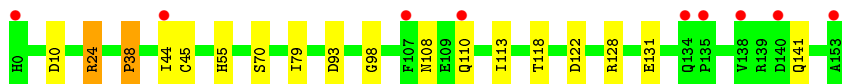
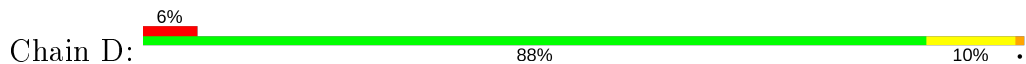
- Molecule 1: Putative peroxiredoxin MT2298



- Molecule 1: Putative peroxiredoxin MT2298



- Molecule 1: Putative peroxiredoxin MT2298



4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	148.20Å 148.20Å 33.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.90 37.05 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.00-1.90) 99.9 (37.05-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 1.89Å)	Xtrriage
Refinement program	REFMAC 3.15	Depositor
R, R_{free}	0.219 , 0.249 0.229 , 0.257	Depositor DCC
R_{free} test set	2986 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtrriage
Anisotropy	0.326	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.065 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5105	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.92	0/1234	0.96	5/1681 (0.3%)
1	B	0.93	0/1231	0.90	2/1675 (0.1%)
1	C	0.95	0/1231	0.88	2/1675 (0.1%)
1	D	0.92	0/1234	0.96	5/1679 (0.3%)
All	All	0.93	0/4930	0.93	14/6710 (0.2%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	24[A]	ARG	NE-CZ-NH1	6.81	123.70	120.30
1	D	24[B]	ARG	NE-CZ-NH1	6.81	123.70	120.30
1	D	24[A]	ARG	NE-CZ-NH2	-6.75	116.92	120.30
1	D	24[B]	ARG	NE-CZ-NH2	-6.75	116.92	120.30
1	A	24	ARG	NE-CZ-NH2	-6.63	116.98	120.30
1	C	10	ASP	CB-CG-OD1	6.30	123.97	118.30
1	A	10[A]	ASP	CB-CG-OD1	6.20	123.88	118.30
1	A	10[B]	ASP	CB-CG-OD1	6.20	123.88	118.30
1	D	10	ASP	CB-CG-OD1	6.16	123.84	118.30
1	A	24	ARG	NE-CZ-NH1	6.08	123.34	120.30
1	A	123	ARG	NE-CZ-NH2	-5.63	117.48	120.30
1	B	10	ASP	CB-CG-OD1	5.45	123.21	118.30
1	C	10	ASP	CB-CG-OD2	-5.15	113.67	118.30
1	B	139	ARG	NE-CZ-NH2	-5.11	117.75	120.30

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	1198	0	1168	9	1
1	B	1198	0	1167	11	0
1	C	1198	0	1167	8	0
1	D	1200	0	1170	14	0
2	A	76	0	0	2	0
2	B	86	0	0	2	0
2	C	82	0	0	2	0
2	D	67	0	0	2	0
All	All	5105	0	4672	41	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 4.

All (41) 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:24[A]:ARG:NH1	2:D:201:HOH:O	2.14	0.79
1:C:63:ASP:HB3	1:C:152:THR:HG21	1.66	0.76
1:C:25:GLY:O	2:C:201:HOH:O	2.07	0.73
1:A:0:HIS:CB	2:A:231:HOH:O	2.41	0.67
1:B:38:PRO:HB3	1:B:113:ILE:HD12	1.79	0.64
1:C:38:PRO:HB3	1:C:113:ILE:HD12	1.81	0.63
1:D:55:HIS:CD2	1:D:141:GLN:HE22	2.18	0.61
1:C:134:GLN:NE2	2:C:202:HOH:O	2.14	0.60
1:D:44:ILE:HG23	1:D:45:CYS:N	2.15	0.59
1:D:38:PRO:HD2	1:D:45:CYS:SG	2.47	0.55
1:A:3:ASN:CG	1:A:4:VAL:H	2.10	0.54
1:B:82:THR:HG21	2:B:266:HOH:O	2.07	0.54
1:D:108:ASN:OD1	1:D:110:GLN:HG3	2.06	0.54
1:D:38:PRO:HB3	1:D:113:ILE:HD12	1.90	0.53
1:D:55:HIS:HB2	1:D:141:GLN:NE2	2.24	0.53
1:C:22:THR:HG22	1:C:24:ARG:HG2	1.91	0.53
1:A:38:PRO:HB3	1:A:113:ILE:HD12	1.91	0.52
1:B:22:THR:HG22	1:B:24:ARG:HG2	1.91	0.52
1:D:44:ILE:CG2	1:D:45:CYS:N	2.73	0.51
1:B:51:GLN:HG2	1:B:141:GLN:OE1	2.10	0.51
1:B:27:ARG:NE	2:B:261:HOH:O	2.33	0.50
1:A:38:PRO:HD2	1:A:45:CYS:SG	2.55	0.47
1:A:1:MET:HE2	1:A:119:PHE:CE2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:VAL:HG12	1:A:132:MET:CE	2.47	0.45
1:B:133:LYS:HE2	1:B:137:GLU:HB3	1.98	0.45
1:C:41:PHE:HB3	1:D:79:ILE:HG13	1.98	0.45
1:D:118:THR:HB	1:D:131:GLU:HG2	1.99	0.45
1:D:55:HIS:CG	1:D:141:GLN:HE22	2.34	0.44
1:B:22:THR:CG2	1:B:24:ARG:HG2	2.47	0.43
1:D:141:GLN:CD	2:D:257:HOH:O	2.56	0.43
1:B:133:LYS:CE	1:B:137:GLU:HB3	2.49	0.43
1:B:51:GLN:HG2	1:B:141:GLN:CD	2.38	0.43
1:C:22:THR:CG2	1:C:24:ARG:HG2	2.48	0.43
1:B:41:PHE:HE2	1:B:76:THR:HG23	1.84	0.43
1:A:37:PHE:CZ	1:A:70:SER:HB3	2.53	0.43
1:C:38:PRO:HD2	1:C:45:CYS:SG	2.59	0.42
1:D:93:ASP:OD2	1:D:98:GLY:HA2	2.19	0.42
1:B:78:LYS:O	1:B:82:THR:HG23	2.19	0.42
1:A:122:ASP:HB3	1:A:128:ARG:HG3	2.02	0.41
1:A:18:GLN:NE2	2:A:262:HOH:O	2.48	0.40
1:D:122:ASP:HB3	1:D:128:ARG:HG3	2.04	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:24:ARG:NH2	1:A:138:VAL:O[1_554]	1.96	0.24

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	153/154 (99%)	148 (97%)	4 (3%)	1 (1%)	22	12
1	B	153/154 (99%)	148 (97%)	4 (3%)	1 (1%)	22	12

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	153/154 (99%)	148 (97%)	5 (3%)	0	100	100
1	D	153/154 (99%)	150 (98%)	2 (1%)	1 (1%)	22	12
All	All	612/616 (99%)	594 (97%)	15 (2%)	3 (0%)	29	18

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	38	PRO
1	B	38	PRO
1	A	38	PRO

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	125/124 (101%)	125 (100%)	0	100	100
1	B	124/124 (100%)	124 (100%)	0	100	100
1	C	124/124 (100%)	124 (100%)	0	100	100
1	D	123/124 (99%)	122 (99%)	1 (1%)	81	82
All	All	496/496 (100%)	495 (100%)	1 (0%)	93	94

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	70	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	GLN
1	D	0	HIS
1	D	141	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 carbohydrates 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

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	153/154 (99%)	0.47	7 (4%) 32 35	14, 21, 40, 59	0
1	B	154/154 (100%)	0.30	3 (1%) 66 69	13, 20, 32, 70	0
1	C	154/154 (100%)	0.44	5 (3%) 47 50	14, 20, 34, 73	0
1	D	154/154 (100%)	0.59	9 (5%) 23 25	15, 22, 42, 63	0
All	All	615/616 (99%)	0.45	24 (3%) 39 42	13, 21, 37, 73	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	153	ALA	10.0
1	C	153	ALA	5.6
1	D	153	ALA	4.8
1	A	135	PRO	4.5
1	C	152	THR	4.2
1	D	135	PRO	3.9
1	C	150	ALA	3.2
1	D	0	HIS	3.2
1	D	44	ILE	2.9
1	D	138	VAL	2.8
1	D	134	GLN	2.8
1	D	107	PHE	2.7
1	A	134	GLN	2.6
1	C	63	ASP	2.6
1	B	151	LEU	2.5
1	A	45	CYS	2.5
1	D	140	ASP	2.5
1	A	110	GLN	2.4
1	D	110	GLN	2.4
1	B	152	THR	2.3
1	A	151	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	109	GLU	2.0
1	A	24	ARG	2.0
1	A	1	MET	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 carbohydrates 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.