Full wwPDB X-ray Structure Validation Report ⓘ

Jun 24, 2024 – 09:28 PM EDT

PDB ID	:	6ITZ
Title	:	Peroxiredoxin from Thermococcus kodakaraensis
Authors	:	Nakamura, T.; Himiyama, T.
Deposited on	:	2018-11-27
Resolution	:	2.96 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:

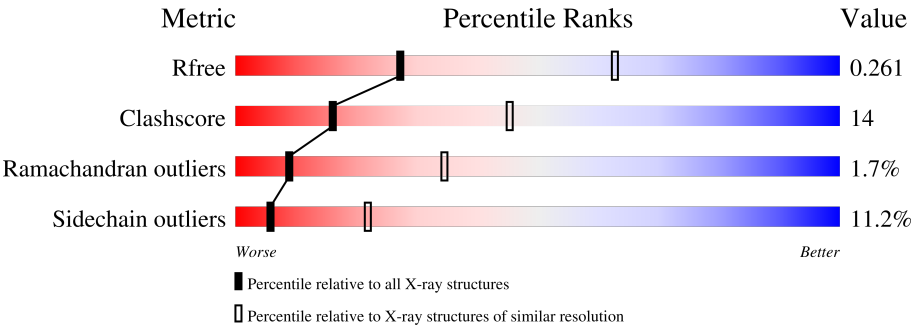
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.96 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	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	216	60% 34% 5%
1	B	216	65% 29% 5%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 3430 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 Peroxiredoxin.

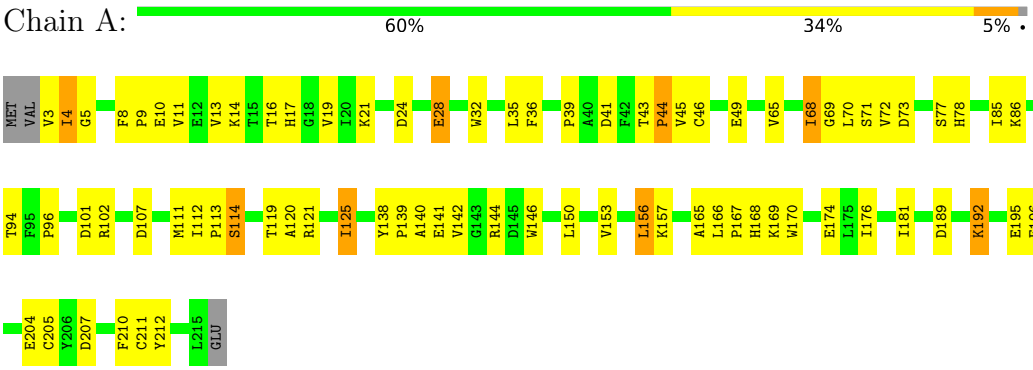
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1715	1117	279	313	6			
1	B	213	Total	C	N	O	S	0	0	0
			1715	1117	279	313	6			

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. 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. 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.

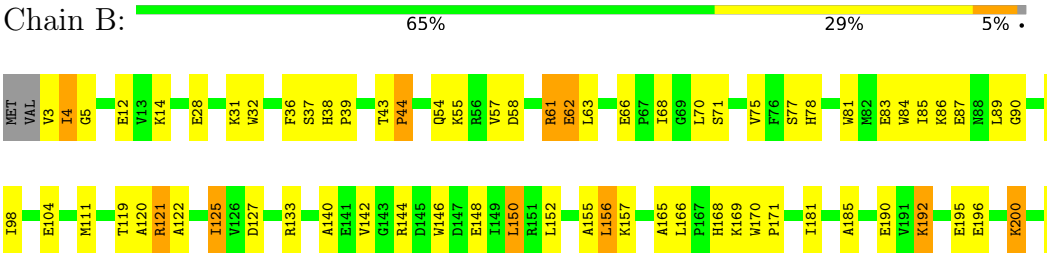
- Molecule 1: Peroxiredoxin

Chain A:



- Molecule 1: Peroxiredoxin

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, α, β, γ	142.18Å 142.18Å 49.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.55 – 2.96 34.15 – 2.96	Depositor EDS
% Data completeness (in resolution range)	99.5 (35.55-2.96) 99.4 (34.15-2.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.63 (at 2.95Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.218 , 0.274 0.206 , 0.261	Depositor DCC
R_{free} test set	577 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å^2)	55.3	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , -5.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.40$, $< L^2 > = 0.22$	Xtriage
Estimated twinning fraction	0.236 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	3430	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.70	0/1761	0.75	0/2389
1	B	0.68	0/1761	0.79	2/2389 (0.1%)
All	All	0.69	0/3522	0.77	2/4778 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	61	ARG	NE-CZ-NH2	-7.74	116.43	120.30
1	B	61	ARG	NE-CZ-NH1	5.27	122.93	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	1715	0	1705	58	0
1	B	1715	0	1703	53	0
All	All	3430	0	3408	99	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 14.

All (99) 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:24:ASP:O	1:A:28:GLU:HB2	1.69	0.91
1:A:121:ARG:NH1	1:A:140:ALA:O	2.09	0.86
1:A:199:ALA:O	1:A:200:LYS:HB2	1.82	0.80
1:A:125:ILE:HD11	1:A:156:LEU:HD23	1.65	0.79
1:A:44:PRO:HB2	1:B:181:ILE:HG21	1.63	0.78
1:B:58:ASP:O	1:B:62:GLU:HB2	1.84	0.76
1:A:181:ILE:HD13	1:B:44:PRO:HB2	1.67	0.75
1:A:168:HIS:ND1	1:A:169:LYS:HB2	2.05	0.72
1:B:54:GLN:HE22	1:B:94:THR:CG2	2.04	0.70
1:A:167:PRO:HD2	1:A:170:TRP:HB2	1.74	0.68
1:B:54:GLN:HE22	1:B:94:THR:HG22	1.59	0.67
1:A:174:GLU:OE1	1:B:55:LYS:HE2	1.94	0.66
1:A:49:GLU:OE2	1:A:144:ARG:N	2.23	0.66
1:B:32:TRP:HZ3	1:B:125:ILE:HD11	1.60	0.66
1:A:39:PRO:HD2	1:A:46:CYS:SG	2.39	0.61
1:A:204:GLU:HB2	1:A:212:TYR:CE1	2.36	0.60
1:A:85:ILE:HG23	1:A:89:LEU:HD12	1.84	0.60
1:A:73:ASP:HB2	1:A:78:HIS:CE1	2.36	0.59
1:B:70:LEU:C	1:B:70:LEU:HD23	2.23	0.59
1:A:142:VAL:HG21	1:B:156:LEU:HD13	1.83	0.59
1:A:41:ASP:OD1	1:A:78:HIS:ND1	2.35	0.59
1:B:32:TRP:CZ3	1:B:125:ILE:HD11	2.36	0.59
1:B:148:GLU:HA	1:B:148:GLU:OE1	2.01	0.58
1:A:10:GLU:HB3	1:A:21:LYS:NZ	2.18	0.58
1:B:196:GLU:O	1:B:200:LYS:HG2	2.03	0.58
1:A:192:LYS:O	1:A:196:GLU:HG3	2.04	0.58
1:B:121:ARG:NH1	1:B:140:ALA:O	2.37	0.57
1:B:125:ILE:HD13	1:B:133:ARG:HB2	1.86	0.57
1:A:70:LEU:HD22	1:A:111:MET:HE1	1.89	0.54
1:A:125:ILE:HD12	1:A:153:VAL:HG22	1.90	0.53
1:A:199:ALA:O	1:A:200:LYS:CB	2.56	0.53
1:B:121:ARG:HG2	1:B:144:ARG:CZ	2.38	0.53
1:A:36:PHE:O	1:A:69:GLY:HA2	2.09	0.53
1:A:205:CYS:SG	1:A:211:CYS:SG	3.07	0.53
1:A:121:ARG:HB2	1:A:138:TYR:O	2.08	0.52
1:B:61:ARG:HD2	1:B:66:GLU:OE2	2.09	0.52
1:B:127:ASP:OD2	1:B:133:ARG:HD3	2.09	0.52
1:B:165:ALA:O	1:B:166:LEU:HD13	2.08	0.52
1:A:65:VAL:HG21	1:A:153:VAL:HG11	1.92	0.51
1:A:125:ILE:CD1	1:A:156:LEU:HD23	2.39	0.51
1:B:43:THR:HB	1:B:44:PRO:HD2	1.92	0.51
1:A:4:ILE:HG22	1:A:5:GLY:N	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:GLU:O	1:A:92:GLU:HG3	2.09	0.51
1:B:36:PHE:HA	1:B:122:ALA:O	2.12	0.49
1:B:38:HIS:CE1	1:B:71:SER:HB3	2.48	0.49
1:B:54:GLN:NE2	1:B:94:THR:HG22	2.27	0.49
1:A:114:SER:HB3	1:B:4:ILE:O	2.12	0.49
1:B:75:VAL:O	1:B:78:HIS:HB2	2.13	0.49
1:A:165:ALA:O	1:A:166:LEU:HD13	2.13	0.48
1:A:10:GLU:HB3	1:A:21:LYS:HZ1	1.76	0.48
1:A:142:VAL:HG22	1:B:155:ALA:HB1	1.94	0.48
1:B:4:ILE:CG2	1:B:5:GLY:N	2.76	0.48
1:A:35:LEU:HA	1:A:68:ILE:O	2.14	0.48
1:B:37:SER:HB3	1:B:111:MET:SD	2.54	0.48
1:B:125:ILE:HD12	1:B:156:LEU:HD23	1.95	0.47
1:B:31:LYS:HD3	1:B:66:GLU:HG2	1.96	0.47
1:A:68:ILE:HD12	1:A:96:PRO:HB2	1.97	0.47
1:B:168:HIS:ND1	1:B:169:LYS:HB2	2.29	0.47
1:B:4:ILE:HG23	1:B:5:GLY:N	2.29	0.47
1:B:85:ILE:HG23	1:B:89:LEU:HD12	1.95	0.47
1:A:167:PRO:CD	1:A:170:TRP:HB2	2.42	0.47
1:A:146:TRP:HA	1:A:146:TRP:CE3	2.50	0.46
1:B:121:ARG:HG2	1:B:144:ARG:NE	2.30	0.46
1:A:142:VAL:HG11	1:B:152:LEU:HD12	1.96	0.46
1:B:86:LYS:O	1:B:90:GLY:HA2	2.15	0.46
1:A:107:ASP:OD1	1:A:112:ILE:HG13	2.15	0.46
1:A:32:TRP:CE2	1:A:157:LYS:HE2	2.51	0.45
1:A:113:PRO:HG3	1:A:120:ALA:HB2	1.97	0.45
1:B:39:PRO:HD3	1:B:120:ALA:O	2.17	0.45
1:B:63:LEU:O	1:B:157:LYS:HE3	2.16	0.45
1:A:168:HIS:CE1	1:A:169:LYS:HB2	2.51	0.45
1:A:43:THR:HB	1:A:44:PRO:CD	2.46	0.45
1:B:70:LEU:HB2	1:B:98:ILE:HB	1.99	0.45
1:A:44:PRO:CB	1:B:181:ILE:HG21	2.40	0.45
1:B:192:LYS:HE3	1:B:192:LYS:HA	1.98	0.44
1:A:45:VAL:HB	1:A:121:ARG:HH12	1.81	0.44
1:A:139:PRO:HB2	1:A:141:GLU:OE1	2.18	0.44
1:A:167:PRO:O	1:A:168:HIS:C	2.57	0.43
1:B:121:ARG:HD2	1:B:142:VAL:O	2.18	0.43
1:A:72:VAL:HG22	1:A:72:VAL:O	2.17	0.43
1:B:185:ALA:HB1	1:B:190:GLU:HB3	1.99	0.43
1:B:170:TRP:CD1	1:B:171:PRO:HA	2.53	0.43
1:A:138:TYR:CE2	1:B:152:LEU:HD13	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:ILE:CG2	1:A:5:GLY:N	2.81	0.43
1:A:71:SER:OG	1:A:78:HIS:HE1	2.02	0.42
1:A:65:VAL:HG21	1:A:153:VAL:CG1	2.49	0.42
1:A:181:ILE:HG21	1:B:44:PRO:HB2	2.00	0.42
1:B:68:ILE:HG13	1:B:96:PRO:HG2	2.01	0.42
1:B:121:ARG:CD	1:B:142:VAL:O	2.67	0.42
1:B:146:TRP:O	1:B:150:LEU:HB2	2.18	0.42
1:A:142:VAL:CG2	1:B:156:LEU:HD13	2.49	0.42
1:B:81:TRP:O	1:B:84:TRP:HB3	2.21	0.41
1:A:196:GLU:O	1:A:199:ALA:O	2.39	0.41
1:B:146:TRP:HA	1:B:146:TRP:CE3	2.56	0.41
1:B:83:GLU:O	1:B:87:GLU:HG2	2.20	0.41
1:A:8:PHE:HA	1:A:9:PRO:HD2	1.93	0.41
1:A:49:GLU:HG2	1:B:168:HIS:HB2	2.03	0.41
1:A:16:THR:OG1	1:A:17:HIS:HD2	2.04	0.40
1:A:146:TRP:HA	1:A:146:TRP:HE3	1.86	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	211/216 (98%)	189 (90%)	17 (8%)	5 (2%)	6 26
1	B	211/216 (98%)	195 (92%)	14 (7%)	2 (1%)	17 51
All	All	422/432 (98%)	384 (91%)	31 (7%)	7 (2%)	9 34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	200	LYS
1	A	4	ILE

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Mol	Chain	Res	Type
1	B	4	ILE
1	A	201	GLY
1	B	44	PRO
1	A	44	PRO
1	A	176	ILE

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	183/186 (98%)	160 (87%)	23 (13%)	4 17
1	B	183/186 (98%)	165 (90%)	18 (10%)	8 27
All	All	366/372 (98%)	325 (89%)	41 (11%)	6 22

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	VAL
1	A	11	VAL
1	A	13	VAL
1	A	14	LYS
1	A	19	VAL
1	A	28	GLU
1	A	68	ILE
1	A	77	SER
1	A	86	LYS
1	A	92	GLU
1	A	94	THR
1	A	101	ASP
1	A	102	ARG
1	A	114	SER
1	A	119	THR
1	A	125	ILE
1	A	150	LEU
1	A	156	LEU

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Mol	Chain	Res	Type
1	A	189	ASP
1	A	192	LYS
1	A	195	GLU
1	A	207	ASP
1	A	210	PHE
1	B	3	VAL
1	B	12	GLU
1	B	14	LYS
1	B	28	GLU
1	B	57	VAL
1	B	62	GLU
1	B	77	SER
1	B	104	GLU
1	B	119	THR
1	B	121	ARG
1	B	125	ILE
1	B	150	LEU
1	B	156	LEU
1	B	192	LYS
1	B	195	GLU
1	B	200	LYS
1	B	210	PHE
1	B	215	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	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 monosaccharides 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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.