



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 1EZK / pdb_00001ezk
Title : Crystal structure of recombinant tryparedoxin I
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Deposited on : 2000-05-11
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

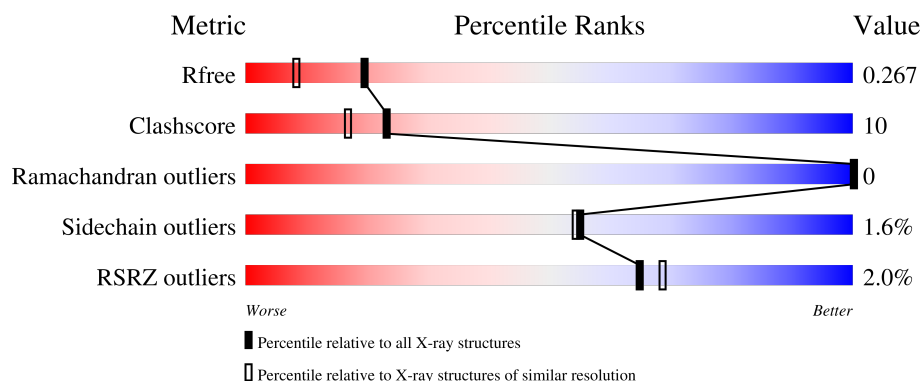
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (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 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	153	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1349 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 TRYPAREDOXIN I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	0	0
			1194	775	193	222	4			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	147	LEU	-	expression tag	UNP O96438
A	148	GLU	-	expression tag	UNP O96438
A	149	HIS	-	expression tag	UNP O96438
A	150	HIS	-	expression tag	UNP O96438
A	151	HIS	-	expression tag	UNP O96438
A	152	HIS	-	expression tag	UNP O96438
A	153	HIS	-	expression tag	UNP O96438

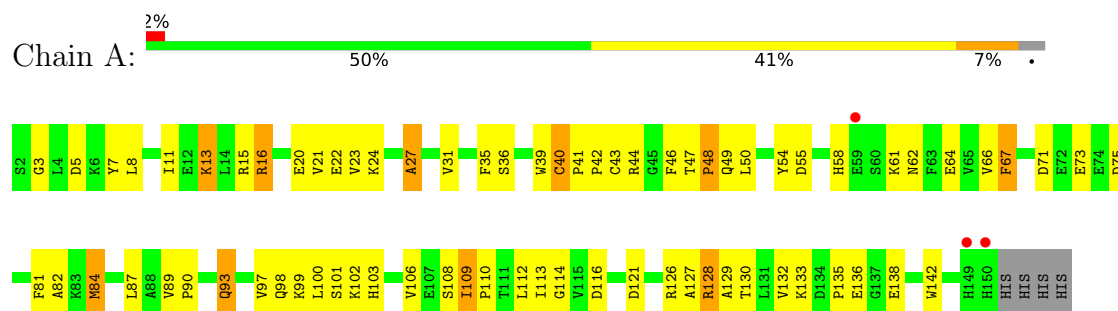
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	155	Total	O	0	0
			155	155		

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: TRYPAREDOXIN I



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	38.10Å 39.73Å 99.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.90 49.59 – 1.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-1.90) 96.4 (49.59-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.42 (at 1.91Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.198 , 0.283 0.183 , 0.267	Depositor DCC
R_{free} test set	585 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.038 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	1349	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.30% 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	1.80	19/1232 (1.5%)	2.51	83/1671 (5.0%)

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	1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	84	MET	SD-CE	-20.78	1.27	1.79
1	A	50	LEU	N-CA	7.16	1.55	1.46
1	A	36	SER	CA-CB	7.08	1.65	1.53
1	A	66	VAL	CA-C	6.41	1.61	1.52
1	A	114	GLY	CA-C	6.28	1.57	1.52
1	A	109	ILE	N-CA	6.14	1.54	1.46
1	A	100	LEU	N-CA	6.02	1.53	1.46
1	A	106	VAL	C-O	-5.85	1.18	1.23
1	A	44	ARG	CA-CB	5.54	1.61	1.53
1	A	35	PHE	C-O	-5.46	1.17	1.24
1	A	84	MET	CA-C	5.44	1.60	1.53
1	A	101	SER	C-N	5.40	1.41	1.33
1	A	42	PRO	N-CD	5.37	1.55	1.47
1	A	98	GLN	C-O	-5.35	1.17	1.24
1	A	128	ARG	N-CA	5.31	1.52	1.46
1	A	54	TYR	C-O	5.29	1.30	1.24
1	A	22	GLU	CA-CB	5.25	1.61	1.53
1	A	55	ASP	C-N	-5.18	1.27	1.33
1	A	90	PRO	N-CD	5.16	1.54	1.47

All (83) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	GLN	OE1-CD-NE2	19.22	141.82	122.60
1	A	113	ILE	CA-C-N	-13.02	112.82	122.33
1	A	113	ILE	C-N-CA	-13.02	112.82	122.33
1	A	66	VAL	O-C-N	12.60	135.26	123.03
1	A	22	GLU	CA-C-O	-10.23	109.08	121.36
1	A	46	PHE	CA-CB-CG	-9.48	104.32	113.80
1	A	15	ARG	NE-CZ-NH2	-9.29	110.84	119.20
1	A	16	ARG	CD-NE-CZ	9.28	137.39	124.40
1	A	67	PHE	O-C-N	9.26	134.32	122.93
1	A	41	PRO	CA-C-O	-9.18	107.26	120.56
1	A	41	PRO	CA-C-N	8.76	128.41	119.56
1	A	41	PRO	C-N-CA	8.76	128.41	119.56
1	A	128	ARG	CA-C-O	-8.11	111.96	120.55
1	A	113	ILE	CA-C-O	-7.62	112.29	120.36
1	A	49	GLN	CG-CD-NE2	7.60	127.80	116.40
1	A	132	VAL	CA-C-O	-7.21	112.86	120.57
1	A	15	ARG	CA-CB-CG	-7.05	99.99	114.10
1	A	41	PRO	N-CA-C	7.00	119.24	110.70
1	A	3	GLY	CA-C-O	-6.77	110.21	119.01
1	A	3	GLY	N-CA-C	-6.71	107.29	114.67
1	A	71	ASP	CA-C-O	-6.71	113.65	121.16
1	A	39	TRP	O-C-N	-6.59	113.05	122.41
1	A	31	VAL	O-C-N	6.58	130.17	123.20
1	A	40	CYS	O-C-N	6.54	126.87	121.31
1	A	126	ARG	NE-CZ-NH2	-6.53	113.32	119.20
1	A	5	ASP	O-C-N	-6.48	115.25	122.12
1	A	127	ALA	CA-C-N	-6.45	111.64	120.28
1	A	127	ALA	C-N-CA	-6.45	111.64	120.28
1	A	61	LYS	CA-C-O	-6.39	111.54	119.28
1	A	49	GLN	OE1-CD-NE2	-6.33	116.27	122.60
1	A	62	ASN	OD1-CG-ND2	6.28	128.88	122.60
1	A	108	SER	O-C-N	6.16	129.55	123.46
1	A	24	LYS	CA-C-O	-6.14	113.17	120.10
1	A	121	ASP	CA-CB-CG	6.13	118.73	112.60
1	A	15	ARG	NH1-CZ-NH2	6.04	127.16	119.30
1	A	31	VAL	CA-C-O	-5.94	114.21	120.39
1	A	55	ASP	CA-C-N	5.94	128.49	120.65
1	A	55	ASP	C-N-CA	5.94	128.49	120.65
1	A	75	ASP	CA-CB-CG	-5.92	106.68	112.60
1	A	5	ASP	CA-CB-CG	-5.91	106.69	112.60
1	A	93	GLN	CG-CD-NE2	-5.90	107.55	116.40
1	A	133	LYS	CA-C-O	-5.87	112.18	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	16	ARG	NE-CZ-NH2	-5.86	113.93	119.20
1	A	127	ALA	CA-C-O	-5.86	114.21	120.42
1	A	40	CYS	CA-C-O	-5.81	114.99	119.46
1	A	89	VAL	O-C-N	-5.80	114.69	121.07
1	A	82	ALA	O-C-N	-5.80	115.44	122.22
1	A	93	GLN	CB-CG-CD	-5.79	102.75	112.60
1	A	102	LYS	CA-C-O	5.79	126.56	120.42
1	A	39	TRP	CA-C-N	5.73	129.22	121.20
1	A	39	TRP	C-N-CA	5.73	129.22	121.20
1	A	128	ARG	CA-C-N	5.69	130.30	120.58
1	A	128	ARG	C-N-CA	5.69	130.30	120.58
1	A	48	PRO	CA-C-N	5.68	128.35	120.29
1	A	48	PRO	C-N-CA	5.68	128.35	120.29
1	A	99	LYS	CA-C-N	-5.65	112.27	120.29
1	A	99	LYS	C-N-CA	-5.65	112.27	120.29
1	A	101	SER	CA-C-O	5.62	126.50	120.55
1	A	27	ALA	CA-C-N	5.59	131.09	122.20
1	A	27	ALA	C-N-CA	5.59	131.09	122.20
1	A	44	ARG	NE-CZ-NH2	-5.58	114.18	119.20
1	A	129	ALA	N-CA-CB	5.55	118.47	110.20
1	A	103	HIS	CA-CB-CG	-5.42	108.39	113.80
1	A	97	VAL	CA-C-O	5.35	126.52	120.85
1	A	114	GLY	O-C-N	5.32	128.48	123.48
1	A	67	PHE	CA-C-O	-5.28	114.74	120.81
1	A	44	ARG	CG-CD-NE	-5.28	100.39	112.00
1	A	138	GLU	CA-CB-CG	-5.27	103.56	114.10
1	A	50	LEU	CA-C-N	5.24	127.16	120.56
1	A	50	LEU	C-N-CA	5.24	127.16	120.56
1	A	35	PHE	O-C-N	5.22	129.17	123.22
1	A	21	VAL	CA-C-O	-5.20	116.48	121.63
1	A	135	PRO	CA-C-O	5.19	127.43	119.55
1	A	93	GLN	CA-CB-CG	-5.18	103.73	114.10
1	A	50	LEU	N-CA-CB	-5.16	102.35	110.30
1	A	93	GLN	CG-CD-OE1	-5.16	110.48	120.80
1	A	54	TYR	CB-CA-C	-5.14	102.59	110.81
1	A	36	SER	CA-CB-OG	-5.13	100.84	111.10
1	A	93	GLN	CB-CA-C	-5.11	102.54	111.23
1	A	116	ASP	CA-C-O	-5.11	114.87	120.69
1	A	71	ASP	N-CA-CB	5.10	117.47	109.97
1	A	44	ARG	NE-CZ-NH1	5.08	126.58	121.50
1	A	16	ARG	NE-CZ-NH1	5.01	126.51	121.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	13	LYS	Mainchain

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	1194	0	1138	23	0
2	A	155	0	0	6	0
All	All	1349	0	1138	23	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 (23) 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:13:LYS:HE3	1:A:20:GLU:OE2	1.89	0.72
1:A:67:PHE:CD2	1:A:84:MET:HE1	2.32	0.64
1:A:128:ARG:HH11	1:A:128:ARG:HG2	1.62	0.63
1:A:81:PHE:HA	1:A:84:MET:HE2	1.81	0.62
1:A:8:LEU:HB3	1:A:11:ILE:HG13	1.83	0.60
1:A:58:HIS:HD2	2:A:282:HOH:O	1.84	0.59
1:A:128:ARG:HG2	1:A:128:ARG:NH1	2.18	0.57
1:A:16:ARG:NH2	1:A:64:GLU:OE2	2.39	0.55
1:A:130:THR:HG1	1:A:142:TRP:CG	2.26	0.54
1:A:7:TYR:O	1:A:8:LEU:HD23	2.10	0.51
1:A:81:PHE:HA	1:A:84:MET:CE	2.41	0.51
1:A:81:PHE:HD1	1:A:84:MET:HE3	1.77	0.50
1:A:73:GLU:HG3	2:A:180:HOH:O	2.13	0.48
1:A:136:GLU:HB2	2:A:184:HOH:O	2.13	0.47
1:A:109:ILE:HA	1:A:110:PRO:C	2.41	0.45
1:A:23:VAL:HG11	2:A:171:HOH:O	2.17	0.44
1:A:128:ARG:NH1	1:A:128:ARG:CG	2.80	0.44
1:A:13:LYS:HG2	2:A:304:HOH:O	2.18	0.43
1:A:47:THR:N	1:A:48:PRO:HD2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:ALA:HB2	2:A:291:HOH:O	2.18	0.43
1:A:16:ARG:HB2	1:A:87:LEU:CD2	2.49	0.42
1:A:67:PHE:CG	1:A:84:MET:HE1	2.55	0.42
1:A:40:CYS:CB	1:A:43:CYS:SG	3.07	0.42

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	147/153 (96%)	144 (98%)	3 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	127/131 (97%)	125 (98%)	2 (2%)	55	54

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	GLN
1	A	112	LEU

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

Mol	Chain	Res	Type
1	A	58	HIS
1	A	98	GLN
1	A	150	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 [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	149/153 (97%)	0.00	3 (2%) 65 69	13, 26, 44, 68	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	150	HIS	3.2
1	A	149	HIS	3.0
1	A	59	GLU	2.7

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.