



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 20, 2016 – 01:54 PM EDT

PDB ID : 5B48
Title : 2-Oxoacid:Ferredoxin Oxidoreductase 1 from Sulfolobus tokodai
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Deposited on : 2016-04-01
Resolution : 2.50 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/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
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027939

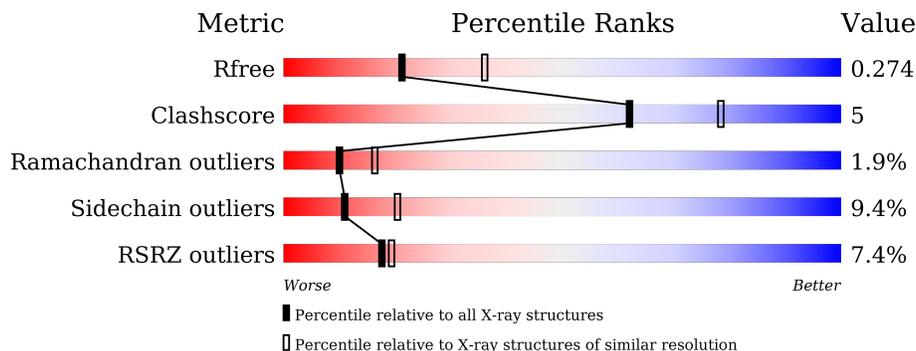
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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. 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	627	 5% 76% 13% • 8%
1	C	627	 7% 72% 15% • 11%
2	B	305	 10% 75% 15% • 6%
2	D	305	 5% 63% 11% •• 24%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13072 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 2-oxoacid--ferredoxin oxidoreductase alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	576	4546	2913	756	864	13	0	0	0
1	C	560	4421	2832	737	841	11	0	0	0

- Molecule 2 is a protein called 2-oxoacid--ferredoxin oxidoreductase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	286	2209	1415	377	412	5	0	0	0
2	D	231	1796	1155	309	331	1	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

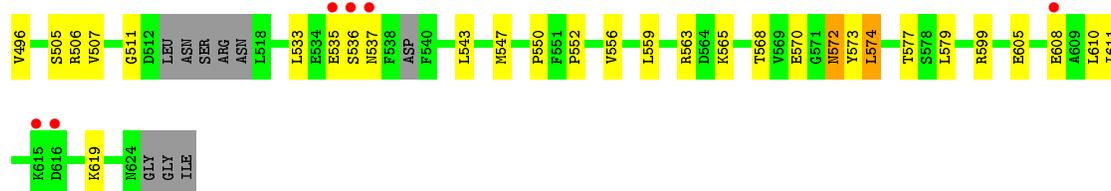
Chain	Residue	Modelled	Actual	Comment	Reference
B	5	THR	LYS	engineered mutation	UNP Q96Y68
D	5	THR	LYS	engineered mutation	UNP Q96Y68

- Molecule 3 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe₄S₄).

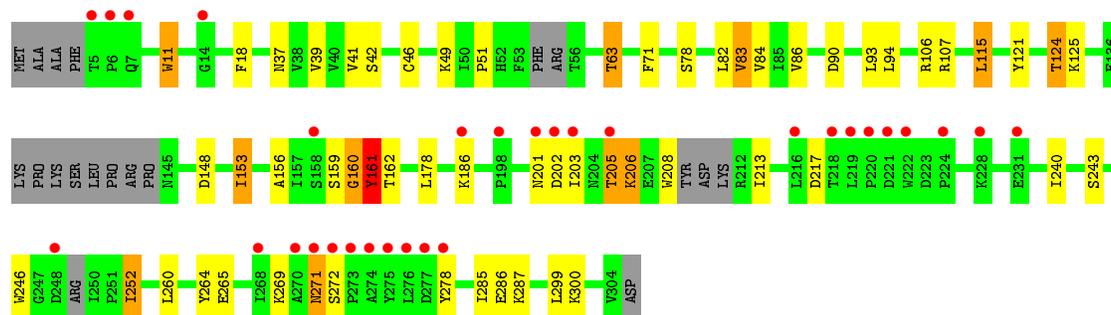
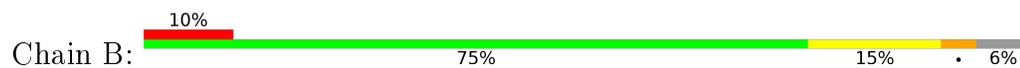
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Mg 1 1	0	0

- Molecule 6 is water.

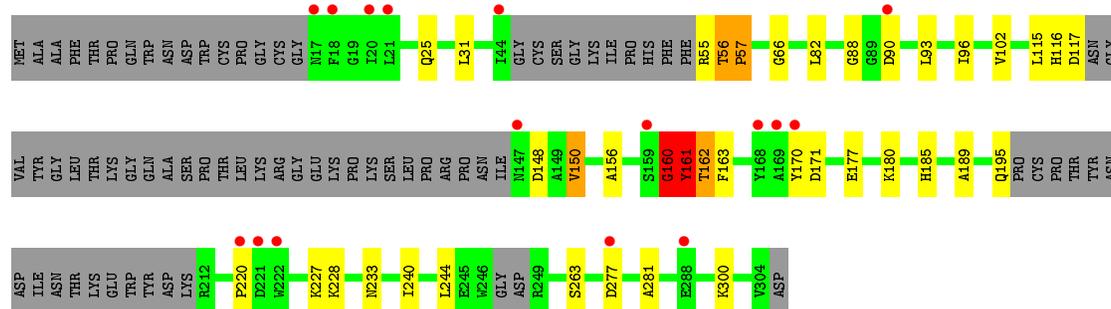
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	25	Total O 25 25	0	0
6	B	5	Total O 5 5	0	0
6	C	23	Total O 23 23	0	0
6	D	8	Total O 8 8	0	0



- Molecule 2: 2-oxoacid--ferredoxin oxidoreductase beta subunit



- Molecule 2: 2-oxoacid--ferredoxin oxidoreductase beta subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.15Å 145.88Å 170.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.47 – 2.50 36.47 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.9 (36.47-2.50) 99.0 (36.47-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.211 , 0.278 0.212 , 0.274	Depositor DCC
R_{free} test set	3274 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	61.0	Xtrriage
Anisotropy	0.026	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13072	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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

Bond lengths and bond angles in the following residue types are not validated in this section: MG, SF4, TDN

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	0.68	0/4629	0.86	6/6253 (0.1%)
1	C	0.65	0/4498	0.84	5/6069 (0.1%)
2	B	0.69	2/2254 (0.1%)	0.90	4/3060 (0.1%)
2	D	0.63	0/1825	0.85	1/2471 (0.0%)
All	All	0.67	2/13206 (0.0%)	0.86	16/17853 (0.1%)

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
1	C	0	1
2	B	0	2
2	D	0	1
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	217	ASP	CB-CG	6.70	1.65	1.51
2	B	265	GLU	CD-OE1	-5.92	1.19	1.25

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	455	ARG	NE-CZ-NH1	8.60	124.60	120.30
1	A	310	ARG	NE-CZ-NH2	-6.75	116.93	120.30
1	C	422	LEU	CA-CB-CG	6.73	130.77	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	11	TRP	CA-CB-CG	5.90	124.92	113.70
1	A	58	ARG	NE-CZ-NH1	5.88	123.24	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	255	ILE	Peptide
2	B	159	SER	Peptide
2	B	51	PRO	Peptide
1	C	511	GLY	Peptide
2	D	160	GLY	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	4546	0	4572	66	0
1	C	4421	0	4428	39	0
2	B	2209	0	2227	30	0
2	D	1796	0	1855	13	0
3	B	8	0	0	0	0
4	B	30	0	0	0	0
5	B	1	0	0	0	0
6	A	25	0	0	1	0
6	B	5	0	0	0	0
6	C	23	0	0	1	0
6	D	8	0	0	0	0
All	All	13072	0	13082	141	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 5.

The worst 5 of 141 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:184:ASN:CA	1:A:187:ILE:HD11	1.25	1.55
1:A:184:ASN:HA	1:A:187:ILE:CD1	1.49	1.40
1:A:184:ASN:C	1:A:187:ILE:HD11	1.66	1.16
1:A:183:VAL:O	1:A:187:ILE:CD1	1.93	1.15
1:A:184:ASN:CA	1:A:187:ILE:CD1	2.16	1.14

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	560/627 (89%)	522 (93%)	26 (5%)	12 (2%)	9	14
1	C	536/627 (86%)	506 (94%)	22 (4%)	8 (2%)	13	22
2	B	276/305 (90%)	248 (90%)	23 (8%)	5 (2%)	11	18
2	D	221/305 (72%)	201 (91%)	14 (6%)	6 (3%)	6	9
All	All	1593/1864 (86%)	1477 (93%)	85 (5%)	31 (2%)	10	16

5 of 31 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	15	ILE
1	A	16	ASP
1	A	116	PHE
1	A	120	GLN
1	A	419	TYR

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	488/535 (91%)	451 (92%)	37 (8%)	16	30
1	C	473/535 (88%)	420 (89%)	53 (11%)	7	14
2	B	239/256 (93%)	217 (91%)	22 (9%)	11	21
2	D	194/256 (76%)	175 (90%)	19 (10%)	10	19
All	All	1394/1582 (88%)	1263 (91%)	131 (9%)	11	20

5 of 131 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	39	TYR
1	C	232	LEU
2	D	162	THR
1	C	52	LEU
1	C	144	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	ASN
1	A	591	ASN
1	C	11	GLN
1	C	144	ASN

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](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	SF4	B	401	2	0,12,12	0.00	-	0,24,24	0.00	-
4	TDN	B	402	5	25,31,31	2.33	8 (32%)	28,46,46	2.53	9 (32%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SF4	B	401	2	-	0/0/48/48	0/6/5/5
4	TDN	B	402	5	-	1/18/23/23	0/2/2/2

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	402	TDN	C5-S1	-8.26	1.58	1.74
4	B	402	TDN	C7'-C5'	-3.67	1.43	1.51
4	B	402	TDN	CM2-C2'	-3.12	1.39	1.49
4	B	402	TDN	PB-O3B	-2.47	1.46	1.54
4	B	402	TDN	PB-O2B	-2.01	1.47	1.54

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	402	TDN	C6-C5-S1	-3.93	114.74	120.24
4	B	402	TDN	CM4-C4-C5	-3.27	121.98	128.91
4	B	402	TDN	N1'-C2'-N3'	-2.32	121.14	125.50
4	B	402	TDN	O2B-PB-O1B	2.54	118.91	110.63
4	B	402	TDN	CM2-C2'-N3'	2.59	121.62	117.20

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	402	TDN	C4-C5-C6-C7

There are no ring outliers.

No monomer is involved in short contacts.

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	576/627 (91%)	0.31	33 (5%) 27 31	37, 62, 107, 143	0
1	C	560/627 (89%)	0.32	42 (7%) 17 19	34, 62, 106, 131	0
2	B	286/305 (93%)	0.53	31 (10%) 8 8	38, 66, 101, 146	0
2	D	231/305 (75%)	0.38	16 (6%) 20 22	42, 65, 100, 137	0
All	All	1653/1864 (88%)	0.36	122 (7%) 17 19	34, 63, 105, 146	0

The worst 5 of 122 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	220	PRO	7.5
2	B	6	PRO	6.2
1	A	194	PHE	6.2
2	D	170	TYR	5.8
1	C	277	ILE	5.3

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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors

of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	TDN	B	402	30/30	0.95	0.14	0.01	47,58,66,68	0
3	SF4	B	401	8/8	0.99	0.03	-1.69	55,58,65,66	0
5	MG	B	403	1/1	0.84	0.08	-3.56	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.