



Full wwPDB X-ray Structure Validation Report

Nov 26, 2014 – 04:05 AM EST

PDB ID : 3WT2
Title : Crystal structure of the b'-a' domain of thermophilic fungal protein disulfide isomerase (oxidized form)
Authors : Inagaki, K.; Satoh, T.; Itoh, S.G.; Okumura, H.; Kato, K.
Deposited on : 2014-04-02
Resolution : 3.30 Å(reported)

This is a full wwPDB validation 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/ValidationPDFNotes.html>

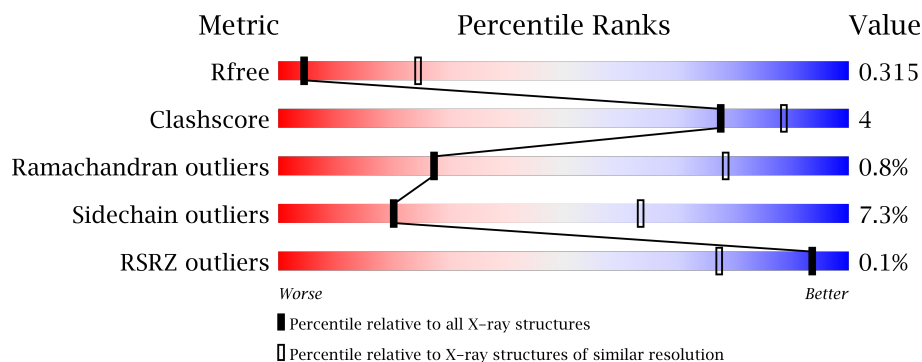
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable24195
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.1.3
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable24195

1 Overall quality at a glance

The reported resolution of this entry is 3.30 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}	66092	1341 (3.40-3.20)
Clashscore	79885	1696 (3.40-3.20)
Ramachandran outliers	78287	1664 (3.40-3.20)
Sidechain outliers	78261	1662 (3.40-3.20)
RSRZ outliers	66119	1342 (3.40-3.20)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	247	
1	B	247	
1	C	247	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5637 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Protein disulfide-isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1879	1214	299	363	3			
1	B	242	Total	C	N	O	S	0	0	0
			1879	1214	299	363	3			
1	C	242	Total	C	N	O	S	0	0	0
			1879	1214	299	363	3			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	203	GLY	-	EXPRESSION TAG	UNP P55059
A	204	PRO	-	EXPRESSION TAG	UNP P55059
A	205	LEU	-	EXPRESSION TAG	UNP P55059
A	206	GLY	-	EXPRESSION TAG	UNP P55059
A	207	SER	-	EXPRESSION TAG	UNP P55059
B	203	GLY	-	EXPRESSION TAG	UNP P55059
B	204	PRO	-	EXPRESSION TAG	UNP P55059
B	205	LEU	-	EXPRESSION TAG	UNP P55059
B	206	GLY	-	EXPRESSION TAG	UNP P55059
B	207	SER	-	EXPRESSION TAG	UNP P55059
C	203	GLY	-	EXPRESSION TAG	UNP P55059
C	204	PRO	-	EXPRESSION TAG	UNP P55059
C	205	LEU	-	EXPRESSION TAG	UNP P55059
C	206	GLY	-	EXPRESSION TAG	UNP P55059
C	207	SER	-	EXPRESSION TAG	UNP P55059

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Protein disulfide-isomerase

Chain A: 



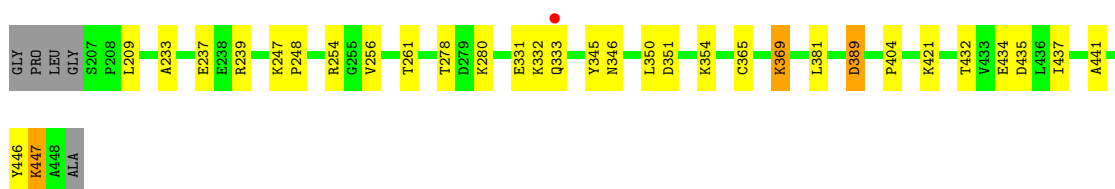
- Molecule 1: Protein disulfide-isomerase

Chain B: 



- Molecule 1: Protein disulfide-isomerase

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	96.31Å 96.31Å 69.49Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.30 19.97 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-3.30) 99.7 (19.97-3.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.55 (at 3.29Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.221 , 0.313 0.226 , 0.315	Depositor DCC
R_{free} test set	516 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	77.7	Xtriage
Anisotropy	0.113	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 28.8	EDS
Estimated twinning fraction	0.000 for -h,-k,l 0.038 for h,-h-k,-l 0.000 for -k,-h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 10736 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5637	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 66.15 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 6.3330e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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	0.64	0/1922	0.76	0/2604
1	B	0.60	0/1922	0.75	0/2604
1	C	0.60	0/1922	0.74	0/2604
All	All	0.61	0/5766	0.75	0/7812

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1879	0	1869	12	0
1	B	1879	0	1869	19	0
1	C	1879	0	1869	16	0
All	All	5637	0	5607	42	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (42) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:388:LYS:HD2	1:C:369:LYS:HZ1	1.57	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:388:LYS:HD2	1:C:369:LYS:NZ	2.08	0.67
1:B:345:TYR:CE1	1:B:404:PRO:HG2	2.36	0.60
1:A:446:TYR:O	1:A:447:LYS:C	2.42	0.59
1:C:369:LYS:O	1:C:369:LYS:HG3	2.03	0.57
1:C:345:TYR:CE1	1:C:404:PRO:HG2	2.39	0.57
1:A:345:TYR:CE1	1:A:404:PRO:HG2	2.40	0.56
1:B:389:ASP:OD1	1:B:389:ASP:N	2.36	0.56
1:C:446:TYR:O	1:C:447:LYS:C	2.45	0.55
1:B:446:TYR:O	1:B:447:LYS:C	2.44	0.55
1:B:441:ALA:O	1:B:447:LYS:HA	2.09	0.51
1:A:441:ALA:O	1:A:447:LYS:HA	2.12	0.50
1:C:441:ALA:O	1:C:447:LYS:HA	2.12	0.50
1:B:233:ALA:HB3	1:B:239:ARG:HG3	1.94	0.50
1:C:381:LEU:CD2	1:C:437:ILE:HD11	2.42	0.50
1:A:389:ASP:OD1	1:A:389:ASP:N	2.37	0.49
1:A:233:ALA:HB3	1:A:239:ARG:HG3	1.94	0.49
1:C:233:ALA:HB3	1:C:239:ARG:HG3	1.95	0.49
1:A:381:LEU:CD2	1:A:437:ILE:HD11	2.43	0.48
1:B:381:LEU:CD2	1:B:437:ILE:HD11	2.43	0.47
1:B:388:LYS:HD2	1:C:369:LYS:CE	2.43	0.47
1:C:354:LYS:HE3	1:C:389:ASP:O	2.16	0.45
1:C:432:THR:HG23	1:C:435:ASP:H	1.81	0.45
1:B:354:LYS:HE3	1:B:389:ASP:O	2.16	0.45
1:B:346:ASN:OD1	1:B:350:LEU:HD12	2.17	0.45
1:A:247:LYS:HB3	1:A:248:PRO:HD3	1.98	0.45
1:B:388:LYS:HG3	1:C:369:LYS:HZ3	1.82	0.45
1:A:354:LYS:HE3	1:A:389:ASP:O	2.16	0.44
1:A:432:THR:HG23	1:A:435:ASP:H	1.82	0.44
1:C:247:LYS:HB3	1:C:248:PRO:HD3	1.99	0.44
1:B:432:THR:HG23	1:B:435:ASP:H	1.83	0.44
1:B:247:LYS:HB3	1:B:248:PRO:HD3	2.00	0.44
1:A:346:ASN:OD1	1:A:350:LEU:HD12	2.17	0.43
1:C:346:ASN:OD1	1:C:350:LEU:HD12	2.18	0.43
1:B:410:PHE:HA	1:B:411:PRO:C	2.40	0.42
1:A:446:TYR:O	1:A:448:ALA:N	2.53	0.42
1:C:381:LEU:HD23	1:C:437:ILE:HD11	2.02	0.41
1:A:350:LEU:HA	1:A:421:LYS:HG2	2.02	0.41
1:B:350:LEU:HA	1:B:421:LYS:HG2	2.02	0.41
1:B:319:LYS:HE3	1:C:237:GLU:OE1	2.20	0.41
1:B:304:THR:O	1:B:305:PHE:C	2.60	0.41
1:B:381:LEU:HD23	1:B:437:ILE:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	240/247 (97%)	230 (96%)	8 (3%)	2 (1%)	27	78
1	B	240/247 (97%)	230 (96%)	8 (3%)	2 (1%)	27	78
1	C	240/247 (97%)	230 (96%)	8 (3%)	2 (1%)	27	78
All	All	720/741 (97%)	690 (96%)	24 (3%)	6 (1%)	27	78

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	447	LYS
1	B	447	LYS
1	C	447	LYS
1	A	351	ASP
1	B	351	ASP
1	C	351	ASP

5.3.2 Protein sidechains ⓘ

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	197/199 (99%)	183 (93%)	14 (7%)	21	65
1	B	197/199 (99%)	182 (92%)	15 (8%)	19	61
1	C	197/199 (99%)	183 (93%)	14 (7%)	21	65
All	All	591/597 (99%)	548 (93%)	43 (7%)	20	63

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	209	LEU
1	A	254	ARG
1	A	256	VAL
1	A	261	THR
1	A	278	THR
1	A	280	LYS
1	A	331	GLU
1	A	332	LYS
1	A	333	GLN
1	A	365	CYS
1	A	369	LYS
1	A	389	ASP
1	A	421	LYS
1	A	434	GLU
1	B	209	LEU
1	B	254	ARG
1	B	256	VAL
1	B	261	THR
1	B	278	THR
1	B	280	LYS
1	B	331	GLU
1	B	332	LYS
1	B	333	GLN
1	B	365	CYS
1	B	369	LYS
1	B	388	LYS
1	B	389	ASP
1	B	421	LYS
1	B	434	GLU
1	C	209	LEU
1	C	254	ARG
1	C	256	VAL
1	C	261	THR
1	C	278	THR
1	C	280	LYS
1	C	331	GLU
1	C	332	LYS
1	C	333	GLN
1	C	365	CYS
1	C	369	LYS
1	C	389	ASP
1	C	421	LYS
1	C	434	GLU

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	273	ASN
1	B	273	ASN
1	C	273	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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	242/247 (97%)	-0.21	0	100	100	44, 69, 110, 136	0
1	B	242/247 (97%)	-0.06	0	100	100	50, 85, 120, 151	0
1	C	242/247 (97%)	-0.02	1 (0%)	90	57	56, 89, 128, 159	0
All	All	726/741 (97%)	-0.10	1 (0%)	93	77	44, 82, 124, 159	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	333	GLN	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.