



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:53 AM GMT

PDB ID : 2YR2
Title : Crystal Structure of the Hypothetical regulator from Sulfolobus tokodaii
Authors : Kumarevel, T.S.; Tanaka, T.; Nishio, M.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2007-04-02
Resolution : 2.01 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
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

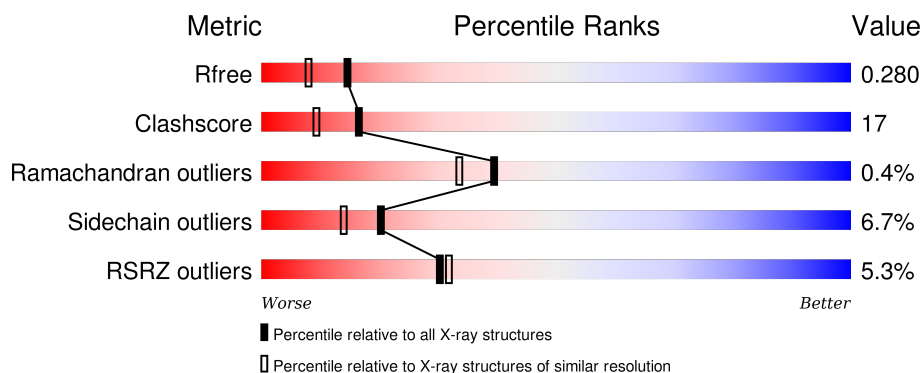
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

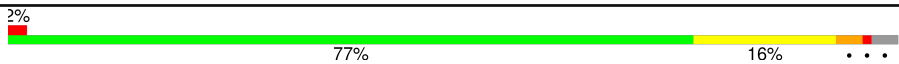
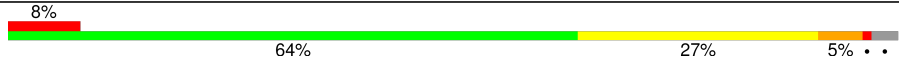
The reported resolution of this entry is 2.01 Å.

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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	146	
1	B	146	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2551 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 146aa long hypothetical transcriptional regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	142	Total	C	N	O	S	0	0	0
			1147	722	201	220	4			
1	B	141	Total	C	N	O	S	0	0	0
			1142	719	200	219	4			

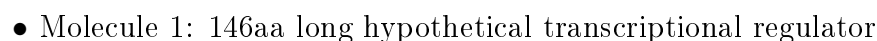
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	156	Total	O	0	0
			156	156		
3	B	105	Total	O	0	0
			105	105		

- Molecule 1: 146aa long hypothetical transcriptional regulator



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	106.26Å 41.93Å 70.58Å 90.00° 109.22° 90.00°	Depositor
Resolution (Å)	20.00 – 2.01 38.69 – 2.01	Depositor EDS
% Data completeness (in resolution range)	98.0 (20.00-2.01) 98.0 (38.69-2.01)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 2.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.234 , 0.279 0.234 , 0.280	Depositor DCC
R_{free} test set	1000 reflections (5.13%)	DCC
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 19540 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2551	wwPDB-VP
Average B, all atoms (Å ²)	34.0	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 33.53 % 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 7.8836e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN

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	0.48	1/1155 (0.1%)	0.61	1/1547 (0.1%)
1	B	0.75	4/1150 (0.3%)	0.70	2/1540 (0.1%)
All	All	0.63	5/2305 (0.2%)	0.66	3/3087 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	95	GLU	CD-OE2	-13.42	1.10	1.25
1	B	95	GLU	CD-OE1	-12.21	1.12	1.25
1	A	10	GLN	C-O	-8.55	1.07	1.23
1	B	95	GLU	C-O	-8.03	1.08	1.23
1	B	95	GLU	CB-CG	-6.42	1.40	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	133	LYS	CD-CE-NZ	7.37	128.66	111.70
1	A	6	GLU	N-CA-C	-7.02	92.05	111.00
1	B	87	GLU	N-CA-C	-5.73	95.53	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1147	0	1202	25	2
1	B	1142	0	1200	53	2
2	B	1	0	0	0	0
3	A	156	0	0	7	0
3	B	105	0	0	8	0
All	All	2551	0	2402	78	2

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 17.

All (78) 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:B:84:ARG:CB	1:B:89:ARG:HH12	1.54	1.19
1:B:89:ARG:HH11	1:B:89:ARG:HA	1.22	1.04
1:B:84:ARG:HB3	1:B:89:ARG:HH12	1.22	1.02
1:B:84:ARG:HB2	1:B:89:ARG:HH12	1.29	0.96
1:B:84:ARG:HB3	1:B:89:ARG:NH1	1.81	0.96
1:B:84:ARG:CB	1:B:89:ARG:NH1	2.36	0.88
1:B:85:ASP:O	1:B:86:ARG:HB2	1.73	0.86
1:B:89:ARG:HD3	3:B:1065:HOH:O	1.77	0.84
1:B:89:ARG:HH11	1:B:89:ARG:CA	1.91	0.84
1:B:6:GLU:HG3	1:B:8:ARG:H	1.43	0.84
1:A:139:LYS:O	1:A:143:GLU:HG3	1.80	0.81
1:A:27:ASN:ND2	1:A:37:TYR:H	1.86	0.74
1:B:53:MET:CG	1:B:94:ILE:HD11	2.19	0.73
1:B:89:ARG:CD	3:B:1065:HOH:O	2.35	0.73
1:B:11:ILE:O	1:B:15:ILE:HG12	1.89	0.71
1:B:39:ASP:OD2	1:B:73:LYS:HE3	1.91	0.71
1:B:27:ASN:ND2	1:B:37:TYR:H	1.90	0.69
1:B:53:MET:HG2	1:B:94:ILE:HD11	1.76	0.68
1:B:79:LEU:HD22	1:B:100:GLY:HA2	1.77	0.67
1:B:89:ARG:HB3	1:B:89:ARG:NH1	2.09	0.67
1:A:53:MET:HE1	1:A:82:ARG:HH21	1.62	0.65
1:B:89:ARG:HA	1:B:89:ARG:NH1	2.05	0.65
1:A:6:GLU:O	1:A:9:ILE:HG22	1.96	0.65
1:A:82:ARG:HB2	3:A:175:HOH:O	1.97	0.64
1:B:79:LEU:HD23	1:B:99:LYS:HG3	1.79	0.63
1:A:79:LEU:HD23	1:A:99:LYS:HG2	1.82	0.61
1:B:89:ARG:HH11	1:B:89:ARG:CB	2.13	0.60
1:B:53:MET:HG3	1:B:94:ILE:HD11	1.82	0.60
1:B:89:ARG:NE	3:B:1065:HOH:O	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:ARG:O	1:B:144:ILE:HD13	2.04	0.58
1:B:138:LEU:O	1:B:142:GLU:HB2	2.04	0.58
1:B:84:ARG:HB2	1:B:89:ARG:NH1	2.10	0.58
1:B:141:ILE:O	1:B:141:ILE:HD13	2.04	0.58
1:A:53:MET:CE	1:A:82:ARG:HH21	2.17	0.57
1:B:79:LEU:CD2	1:B:99:LYS:HG3	2.35	0.57
1:B:89:ARG:NH1	1:B:89:ARG:CB	2.67	0.57
1:B:121:ASP:HA	3:B:1053:HOH:O	2.06	0.55
1:B:145:SER:O	1:B:146:GLN:HG2	2.06	0.55
1:B:6:GLU:HG3	1:B:7:ASN:N	2.20	0.54
1:A:77:MET:HE3	3:A:267:HOH:O	2.08	0.53
1:B:6:GLU:CG	1:B:7:ASN:N	2.71	0.53
1:A:83:VAL:HG12	3:A:281:HOH:O	2.07	0.53
1:A:136:LYS:HE3	3:A:178:HOH:O	2.09	0.52
1:A:32:GLU:HG2	3:A:185:HOH:O	2.09	0.52
1:A:124:GLU:O	1:A:128:ILE:HD13	2.10	0.51
1:A:64:GLN:NE2	3:A:152:HOH:O	2.44	0.51
1:B:128:ILE:HG13	1:B:129:LEU:N	2.25	0.51
1:A:39:ASP:OD2	1:A:73:LYS:HE2	2.10	0.50
1:B:96:ILE:HD13	1:B:101:LEU:HG	1.95	0.48
1:B:6:GLU:HA	3:B:1054:HOH:O	2.13	0.48
1:B:16:ALA:O	1:B:20:ARG:HG3	2.14	0.48
1:A:10:GLN:HB2	1:A:10:GLN:HE21	1.51	0.47
1:A:85:ASP:O	1:A:89:ARG:HD3	2.16	0.46
1:A:44:ARG:HA	1:A:104:PHE:CZ	2.51	0.46
1:B:92:ILE:HD12	1:B:92:ILE:N	2.30	0.46
1:A:79:LEU:CD2	1:A:99:LYS:HG2	2.46	0.45
1:A:27:ASN:HD21	1:A:37:TYR:H	1.63	0.45
1:B:141:ILE:O	1:B:145:SER:OG	2.35	0.44
1:B:26:LEU:HD13	1:B:37:TYR:CE1	2.52	0.44
1:B:98:GLU:HG3	1:B:99:LYS:N	2.32	0.44
1:B:7:ASN:O	1:B:11:ILE:HG13	2.18	0.44
1:B:109:GLU:HG3	3:B:1090:HOH:O	2.17	0.43
1:A:139:LYS:HG3	3:A:195:HOH:O	2.18	0.43
1:B:143:GLU:HG3	1:B:144:ILE:HD12	2.00	0.43
1:B:28:ARG:HH11	1:B:28:ARG:HG2	1.84	0.43
1:A:10:GLN:HG3	1:A:10:GLN:O	2.18	0.43
1:B:56:LEU:HD12	1:B:94:ILE:HD13	2.00	0.43
1:A:9:ILE:HD13	1:A:9:ILE:O	2.19	0.43
1:B:85:ASP:O	1:B:86:ARG:CB	2.52	0.42
1:A:53:MET:HE1	1:A:68:THR:HA	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ARG:NH1	1:A:92:ILE:HD11	2.35	0.42
1:B:78:GLY:HA2	3:B:1006:HOH:O	2.20	0.41
1:B:6:GLU:CD	1:B:7:ASN:H	2.19	0.41
1:B:27:ASN:HD21	1:B:37:TYR:H	1.65	0.41
1:B:59:ARG:HD3	3:B:1052:HOH:O	2.19	0.41
1:B:48:ASP:CG	1:B:48:ASP:O	2.59	0.41
1:B:28:ARG:NH1	1:B:28:ARG:HG2	2.37	0.40
1:A:111:TYR:HD1	1:A:111:TYR:O	2.04	0.40

All (2) 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:87:GLU:OE2	1:B:85:ASP:O[4_454]	2.17	0.03
1:A:143:GLU:OE1	1:B:140:ARG:NH1[4_444]	2.19	0.01

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	140/146 (96%)	140 (100%)	0	0	100	100
1	B	139/146 (95%)	135 (97%)	3 (2%)	1 (1%)	26	19
All	All	279/292 (96%)	275 (99%)	3 (1%)	1 (0%)	39	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	145	SER

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	127/132 (96%)	121 (95%)	6 (5%)	32	27
1	B	127/132 (96%)	116 (91%)	11 (9%)	13	7
All	All	254/264 (96%)	237 (93%)	17 (7%)	20	14

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLU
1	A	9	ILE
1	A	10	GLN
1	A	27	ASN
1	A	99	LYS
1	A	111	TYR
1	B	10	GLN
1	B	48	ASP
1	B	86	ARG
1	B	89	ARG
1	B	91	LYS
1	B	95	GLU
1	B	111	TYR
1	B	128	ILE
1	B	133	LYS
1	B	141	ILE
1	B	145	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	27	ASN
1	A	34	ASN
1	A	64	GLN
1	A	105	ASN

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Mol	Chain	Res	Type
1	B	10	GLN
1	B	27	ASN
1	B	34	ASN
1	B	105	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	142/146 (97%)	-0.05	3 (2%) 67 67	17, 26, 40, 72	0
1	B	141/146 (96%)	0.34	12 (8%) 13 14	23, 37, 50, 84	0
All	All	283/292 (96%)	0.15	15 (5%) 30 32	17, 31, 50, 84	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	145	SER	4.4
1	B	146	GLN	4.1
1	B	145	SER	3.8
1	B	101	LEU	3.2
1	B	98	GLU	3.1
1	A	146	GLN	2.9
1	B	144	ILE	2.7
1	A	5	ASN	2.2
1	B	73	LYS	2.2
1	B	89	ARG	2.1
1	B	32	GLU	2.1
1	B	116	ASN	2.1
1	B	86	ARG	2.0
1	B	120	GLY	2.0
1	B	125	ASP	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](#)

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
2	ZN	B	1001	1/1	1.00	0.07	-0.86	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.