Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 01:11 PM GMT

PDB ID	:	3T7J
Title	:	Crystal structure of Rtt107p (residues 820-1070)
Authors	:	Li, X.; Li, F.; Wu, J.; Shi, Y.
Deposited on	:	2011-07-30
Resolution	:	2.04 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

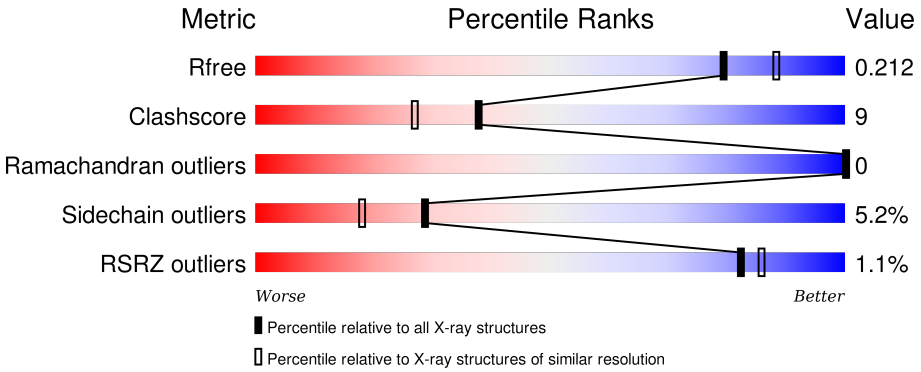
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.04 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	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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 <=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	256	% 74% 15% • 7%
1	B	256	% 74% 17% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3988 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 Regulator of Ty1 transposition protein 107.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	9	0	0
			1926	1251	320	349	6			
1	B	237	Total	C	N	O	S	10	0	0
			1926	1251	320	349	6			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	815	GLY	-	EXPRESSION TAG	UNP P38850
A	816	SER	-	EXPRESSION TAG	UNP P38850
A	817	PRO	-	EXPRESSION TAG	UNP P38850
A	818	HIS	-	EXPRESSION TAG	UNP P38850
A	819	MET	-	EXPRESSION TAG	UNP P38850
B	815	GLY	-	EXPRESSION TAG	UNP P38850
B	816	SER	-	EXPRESSION TAG	UNP P38850
B	817	PRO	-	EXPRESSION TAG	UNP P38850
B	818	HIS	-	EXPRESSION TAG	UNP P38850
B	819	MET	-	EXPRESSION TAG	UNP P38850

- Molecule 2 is water.

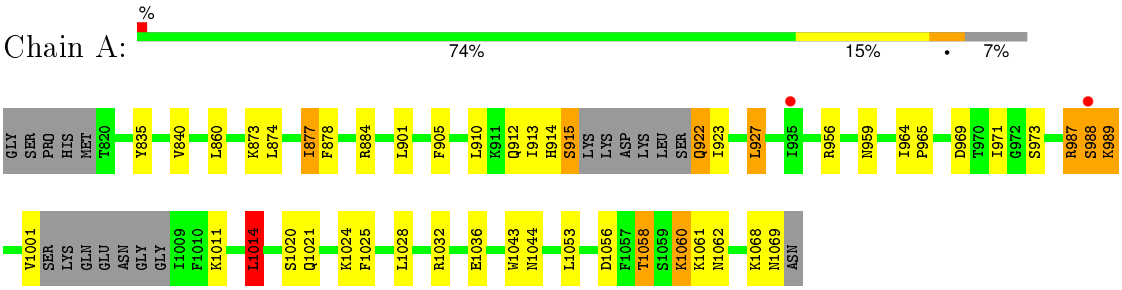
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	72	Total	O	0	0
			72	72		
2	B	64	Total	O	0	0
			64	64		

3 Residue-property plots (i)

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.

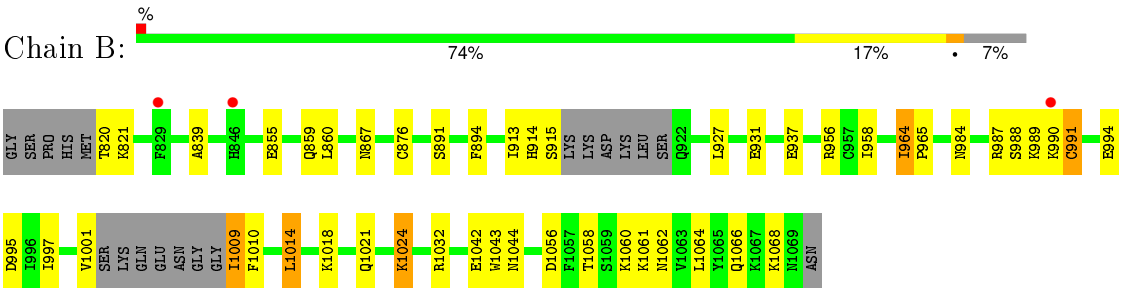
- Molecule 1: Regulator of Ty1 transposition protein 107

Chain A:



- Molecule 1: Regulator of Ty1 transposition protein 107

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	34.36Å 62.18Å 65.80Å 86.64° 75.38° 73.98°	Depositor
Resolution (Å)	32.05 – 2.04 32.05 – 2.04	Depositor EDS
% Data completeness (in resolution range)	94.6 (32.05-2.04) 92.4 (32.05-2.04)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.04 (at 2.05Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.203 , 0.254 0.204 , 0.212	Depositor DCC
R_{free} test set	1538 reflections (5.06%)	DCC
Wilson B-factor (Å^2)	29.8	Xtriage
Anisotropy	0.332	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 35.2	EDS
Estimated twinning fraction	0.397 for h,h-k,h-l 0.009 for -h,-h+k,-l 0.007 for -h,-k,-h+l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 31155 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3988	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.97% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.40	0/1958	0.56	1/2637 (0.0%)
1	B	0.38	0/1958	0.56	1/2637 (0.0%)
All	All	0.39	0/3916	0.56	2/5274 (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	1014	LEU	CA-CB-CG	5.98	129.05	115.30
1	A	1014	LEU	CA-CB-CG	5.64	128.26	115.30

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	1926	0	2004	35	0
1	B	1926	0	2004	37	0
2	A	72	0	0	1	0
2	B	64	0	0	1	0
All	All	3988	0	4008	72	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 9.

All (72) 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:1018:LYS:HB3	1:B:1021:GLN:HG3	1.54	0.87
1:B:860:LEU:HD11	1:B:913:ILE:HD11	1.67	0.74
1:B:855:GLU:O	1:B:859:GLN:HG2	1.87	0.74
1:B:931:GLU:HG3	2:B:52:HOH:O	1.87	0.74
1:A:914:HIS:O	1:A:915:SER:HB3	1.86	0.74
1:B:989:LYS:O	1:B:989:LYS:HD3	1.90	0.72
1:A:860:LEU:HD11	1:A:913:ILE:HD11	1.70	0.72
1:B:987:ARG:HD2	1:B:990:LYS:NZ	2.07	0.69
1:A:1011:LYS:HE3	1:A:1036:GLU:HG3	1.76	0.68
1:A:1056:ASP:OD1	1:A:1058:THR:HB	1.95	0.67
1:A:959:ASN:HB2	1:A:1014:LEU:HD13	1.80	0.62
1:B:1009:ILE:HD12	1:B:1010:PHE:H	1.66	0.60
1:B:1056:ASP:OD1	1:B:1058:THR:HB	2.02	0.60
1:A:1061:LYS:HG2	1:A:1062:ASN:ND2	2.19	0.57
1:A:1060:LYS:HE3	1:A:1060:LYS:O	2.05	0.56
1:A:860:LEU:HD11	1:A:913:ILE:CD1	2.35	0.56
1:B:964:ILE:HD11	1:B:1043:TRP:HB2	1.88	0.55
1:A:835:TYR:CD1	1:A:927:LEU:HD11	2.42	0.55
1:B:987:ARG:HD2	1:B:990:LYS:HZ2	1.72	0.54
1:A:840:VAL:HG23	1:A:877:ILE:HD12	1.90	0.53
1:A:964:ILE:HD12	1:A:965:PRO:HD2	1.92	0.52
1:B:984:ASN:HB2	1:B:997:ILE:HD13	1.90	0.52
1:A:912:GLN:OE1	1:A:923:ILE:HA	2.10	0.52
1:A:914:HIS:O	1:A:915:SER:CB	2.58	0.51
1:A:956:ARG:HE	1:A:1001:VAL:HG23	1.76	0.51
1:B:987:ARG:HD2	1:B:990:LYS:HZ3	1.73	0.51
1:B:994:GLU:OE2	1:B:1032:ARG:NH2	2.45	0.50
1:B:958:ILE:C	1:B:958:ILE:HD12	2.32	0.49
1:B:990:LYS:HZ2	1:B:990:LYS:HB2	1.77	0.49
1:B:1044:ASN:HD22	1:B:1062:ASN:HD21	1.59	0.49
1:A:878:PHE:HA	1:A:901:LEU:O	2.14	0.48
1:B:839:ALA:HB2	1:B:876:CYS:HB3	1.96	0.48
1:A:840:VAL:HG13	1:A:874:LEU:HD13	1.96	0.47
1:A:989:LYS:HD2	1:A:989:LYS:N	2.28	0.47
1:B:914:HIS:O	1:B:915:SER:HB2	2.15	0.47
1:B:1060:LYS:HB3	1:B:1060:LYS:HE3	1.62	0.47
1:A:1024:LYS:NZ	1:A:1028:LEU:HD11	2.30	0.46
1:A:971:ILE:HG23	1:A:1043:TRP:CZ3	2.50	0.46
1:B:820:THR:HG22	1:B:821:LYS:N	2.31	0.46
1:B:914:HIS:O	1:B:915:SER:CB	2.64	0.46
1:A:964:ILE:O	1:A:964:ILE:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:820:THR:HG21	1:B:855:GLU:OE1	2.17	0.45
1:B:988:SER:HA	1:B:989:LYS:HA	1.53	0.45
1:A:860:LEU:CD1	1:A:913:ILE:HD11	2.41	0.45
1:B:1066:GLN:OE1	1:B:1068:LYS:HE2	2.18	0.44
1:A:905:PHE:CE1	1:A:927:LEU:HD12	2.52	0.44
1:A:989:LYS:CD	1:A:989:LYS:N	2.81	0.44
1:B:989:LYS:HG2	1:B:1024:LYS:HE2	1.99	0.44
1:A:1021:GLN:O	1:A:1025:PHE:CD2	2.71	0.44
1:B:989:LYS:HG2	1:B:1024:LYS:CE	2.48	0.44
1:A:884:ARG:NH2	1:A:1044:ASN:OD1	2.47	0.43
1:B:964:ILE:HA	1:B:965:PRO:HD3	1.89	0.43
1:B:1009:ILE:CD1	1:B:1010:PHE:H	2.31	0.43
1:A:835:TYR:HD1	1:A:927:LEU:HD11	1.81	0.43
1:B:956:ARG:HG3	1:B:1001:VAL:HG23	2.01	0.42
1:B:860:LEU:HD11	1:B:913:ILE:CD1	2.45	0.42
1:A:910:LEU:HD23	1:A:910:LEU:HA	1.80	0.42
1:B:989:LYS:CG	1:B:1024:LYS:HE2	2.49	0.42
1:B:1042:GLU:HG3	1:B:1064:LEU:HD11	2.01	0.42
1:A:1021:GLN:O	1:A:1025:PHE:HD2	2.03	0.41
1:A:1053:LEU:HA	1:A:1053:LEU:HD23	1.86	0.41
1:B:1061:LYS:HG3	1:B:1062:ASN:ND2	2.36	0.41
1:A:873:LYS:HA	1:A:873:LYS:HD3	1.94	0.41
1:A:987:ARG:O	1:A:987:ARG:HG2	2.21	0.41
1:B:1009:ILE:HD12	1:B:1010:PHE:N	2.34	0.41
1:B:991:CYS:HA	1:B:995:ASP:OD2	2.19	0.41
1:A:1032:ARG:NH1	2:A:61:HOH:O	2.27	0.41
1:A:922:GLN:CA	1:A:922:GLN:HE21	2.32	0.41
1:B:984:ASN:HB2	1:B:997:ILE:CD1	2.51	0.41
1:A:988:SER:HA	1:A:989:LYS:HA	1.52	0.40
1:B:891:SER:HA	1:B:894:PHE:CE1	2.57	0.40
1:A:922:GLN:NE2	1:A:922:GLN:CA	2.84	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	231/256 (90%)	226 (98%)	5 (2%)	0	100	100
1	B	231/256 (90%)	226 (98%)	5 (2%)	0	100	100
All	All	462/512 (90%)	452 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	221/237 (93%)	206 (93%)	15 (7%)	20	11
1	B	221/237 (93%)	213 (96%)	8 (4%)	42	34
All	All	442/474 (93%)	419 (95%)	23 (5%)	29	18

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	877	ILE
1	A	915	SER
1	A	922	GLN
1	A	927	LEU
1	A	969	ASP
1	A	973	SER
1	A	987	ARG
1	A	988	SER
1	A	989	LYS
1	A	1014	LEU
1	A	1020	SER
1	A	1058	THR
1	A	1060	LYS
1	A	1068	LYS
1	A	1069	ASN

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Mol	Chain	Res	Type
1	B	867	ASN
1	B	927	LEU
1	B	937	GLU
1	B	964	ILE
1	B	991	CYS
1	B	1009	ILE
1	B	1014	LEU
1	B	1024	LYS

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	1062	ASN
1	B	1030	ASN
1	B	1062	ASN
1	B	1066	GLN

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

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, 95^{th} 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	237/256 (92%)	-0.23	2 (0%)	87	90	23, 34, 55, 67	40 (16%)
1	B	237/256 (92%)	-0.07	3 (1%)	79	83	23, 36, 56, 66	38 (16%)
All	All	474/512 (92%)	-0.15	5 (1%)	82	86	23, 36, 55, 67	78 (16%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	990	LYS	3.1
1	B	829	PHE	2.3
1	A	935	ILE	2.3
1	B	846	HIS	2.2
1	A	988	SER	2.1

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.