Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 05:07 PM GMT

PDB ID	:	4HA7
Title	:	Structural insights into the reduction mechanism of Saccharomyces cerevisia Riboflavin Biosynthesis Reductase Rib7
Authors	:	Lv, Z.; Sun, J.; Liu, Y.
Deposited on	:	2012-09-25
Resolution	:	2.10 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:

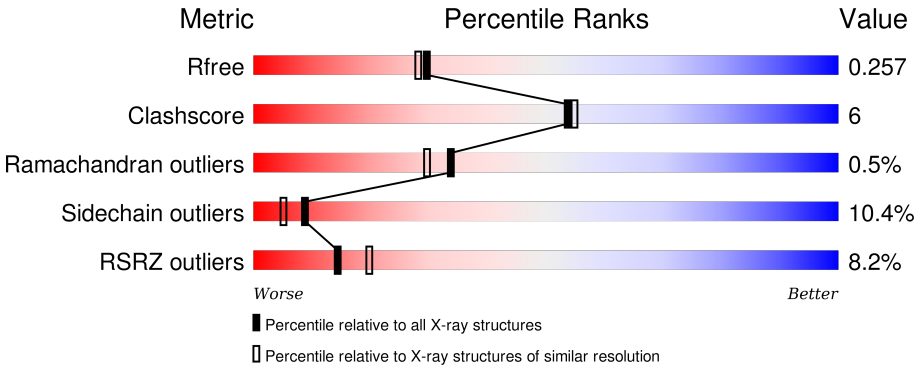
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	249	3% 66% 18% • 11%
1	B	249	11% 64% 16% • 18%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3514 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 2,5-diamino-6-ribosylamino-4(3H)-pyrimidinone 5'-phosphate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1741	1111	299	324	7			
1	B	205	Total	C	N	O	S	0	0	0
			1632	1043	279	304	6			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP P33312
A	-3	PRO	-	EXPRESSION TAG	UNP P33312
A	-2	LEU	-	EXPRESSION TAG	UNP P33312
A	-1	GLY	-	EXPRESSION TAG	UNP P33312
A	0	SER	-	EXPRESSION TAG	UNP P33312
B	-4	GLY	-	EXPRESSION TAG	UNP P33312
B	-3	PRO	-	EXPRESSION TAG	UNP P33312
B	-2	LEU	-	EXPRESSION TAG	UNP P33312
B	-1	GLY	-	EXPRESSION TAG	UNP P33312
B	0	SER	-	EXPRESSION TAG	UNP P33312

- Molecule 2 is water.

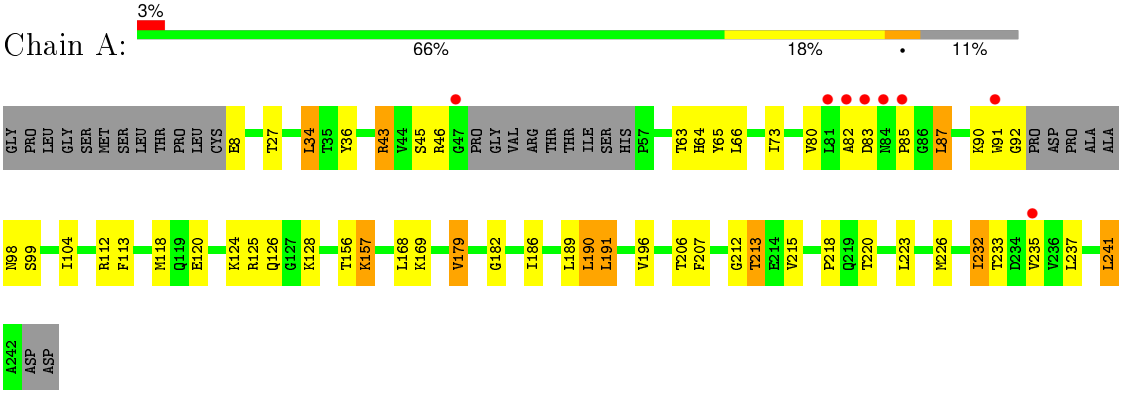
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	95	Total	O	0	0
			95	95		
2	B	46	Total	O	0	0
			46	46		

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.

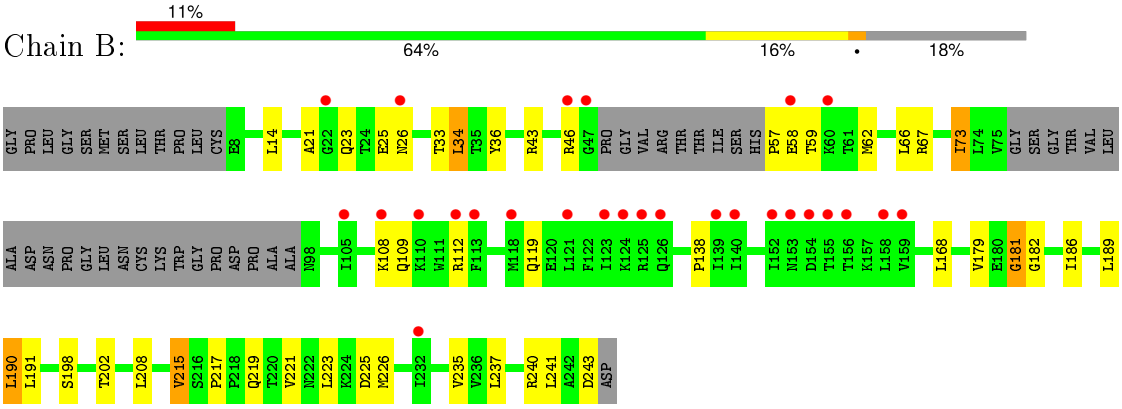
- Molecule 1: 2,5-diamino-6-ribosylamino-4(3H)-pyrimidinone 5'-phosphate reductase

Chain A:



- Molecule 1: 2,5-diamino-6-ribosylamino-4(3H)-pyrimidinone 5'-phosphate reductase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	45.50Å 68.62Å 148.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.75 – 2.10 36.75 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.6 (36.75-2.10) 97.6 (36.75-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.33 (at 2.10Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.200 , 0.266 0.193 , 0.257	Depositor DCC
R_{free} test set	1370 reflections (4.99%)	DCC
Wilson B-factor (Å^2)	41.7	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 60.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 27561 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3514	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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	0.42	0/1777	0.64	0/2413
1	B	0.36	0/1665	0.60	0/2260
All	All	0.39	0/3442	0.62	0/4673

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	181	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	1741	0	1765	28	0
1	B	1632	0	1658	19	0
2	A	95	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	46	0	0	0	0
All	All	3514	0	3423	44	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 6.

All (44) 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:225:ASP:OD1	1:B:240:ARG:NH1	2.22	0.72
1:A:8:GLU:OE1	2:A:332:HOH:O	2.05	0.72
1:A:45:SER:HB2	1:A:213:THR:HG23	1.77	0.67
1:A:235:VAL:HG21	1:B:235:VAL:HG21	1.79	0.65
1:B:109:GLN:HE22	1:B:138:PRO:HB3	1.61	0.64
1:A:206:THR:HA	1:B:223:LEU:HD13	1.81	0.63
1:A:99:SER:O	1:A:128:LYS:NZ	2.28	0.62
1:A:186:ILE:HG22	1:A:190:LEU:HD22	1.83	0.59
1:A:83:ASP:O	1:A:85:PRO:HD3	2.03	0.58
1:A:85:PRO:HB2	1:A:87:LEU:HD13	1.85	0.58
1:A:182:GLY:O	1:A:186:ILE:HG12	2.06	0.56
1:A:63:THR:OG1	2:A:383:HOH:O	2.17	0.55
1:A:36:TYR:CE1	1:A:215:VAL:HG11	2.43	0.54
1:A:191:LEU:HD21	1:A:218:PRO:HD2	1.90	0.53
1:A:80:VAL:HG21	1:A:104:ILE:HG23	1.91	0.53
1:A:196:VAL:HB	1:A:241:LEU:HD22	1.91	0.52
1:A:98:ASN:ND2	2:A:356:HOH:O	2.37	0.52
1:B:181:GLY:HA2	1:B:186:ILE:HD11	1.92	0.51
1:B:109:GLN:NE2	1:B:138:PRO:HB3	2.24	0.51
1:B:67:ARG:HE	1:B:73:ILE:HD12	1.76	0.50
1:A:232:ILE:HG13	1:A:233:THR:HG23	1.94	0.50
1:B:190:LEU:HD23	1:B:217:PRO:HD3	1.92	0.50
1:A:64:HIS:HB3	1:A:90:LYS:O	2.11	0.49
1:A:27:THR:O	1:A:169:LYS:NZ	2.30	0.49
1:B:21:ALA:O	1:B:25:GLU:HG3	2.13	0.48
1:B:36:TYR:CE1	1:B:215:VAL:HG11	2.50	0.46
1:B:33:THR:HG22	1:B:198:SER:HB2	1.97	0.46
1:B:14:LEU:HD11	1:B:62:MET:HG3	1.98	0.46
1:A:65:TYR:HB2	1:A:92:GLY:HA3	1.98	0.44
1:B:182:GLY:O	1:B:186:ILE:HG12	2.17	0.44
1:A:156:THR:O	1:A:157:LYS:HD3	2.17	0.44
1:A:73:ILE:HD13	1:A:87:LEU:HG	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:LEU:HD23	1:A:179:VAL:HG22	2.00	0.44
1:A:207:PHE:HB2	1:B:221:VAL:HB	1.99	0.44
1:A:43:ARG:HG2	1:A:212:GLY:C	2.39	0.43
1:B:57:PRO:HA	1:B:58:GLU:HA	1.63	0.42
1:B:34:LEU:HD23	1:B:179:VAL:HG22	2.01	0.42
1:A:90:LYS:HG2	1:A:91:TRP:N	2.35	0.41
1:B:208:LEU:HD23	1:B:208:LEU:HA	1.88	0.41
1:A:90:LYS:HB3	1:A:90:LYS:HE2	1.82	0.41
1:A:113:PHE:HE2	1:A:118:MET:HG2	1.86	0.41
1:A:124:LYS:O	1:A:126:GLN:HG3	2.21	0.41
1:B:58:GLU:OE1	1:B:58:GLU:N	2.55	0.40
1:B:59:THR:HG21	1:B:202:THR:HG21	2.02	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	215/249 (86%)	206 (96%)	8 (4%)	1 (0%)	34	30
1	B	199/249 (80%)	184 (92%)	14 (7%)	1 (0%)	34	30
All	All	414/498 (83%)	390 (94%)	22 (5%)	2 (0%)	34	30

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	23	GLN
1	A	82	ALA

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	198/221 (90%)	177 (89%)	21 (11%)	8 5
1	B	187/221 (85%)	168 (90%)	19 (10%)	9 5
All	All	385/442 (87%)	345 (90%)	40 (10%)	9 5

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LEU
1	A	43	ARG
1	A	46	ARG
1	A	66	LEU
1	A	87	LEU
1	A	112	ARG
1	A	120	GLU
1	A	125	ARG
1	A	157	LYS
1	A	168	LEU
1	A	179	VAL
1	A	189	LEU
1	A	190	LEU
1	A	191	LEU
1	A	213	THR
1	A	220	THR
1	A	223	LEU
1	A	226	MET
1	A	232	ILE
1	A	237	LEU
1	A	241	LEU
1	B	26	ASN
1	B	34	LEU
1	B	43	ARG
1	B	46	ARG
1	B	66	LEU
1	B	73	ILE

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Mol	Chain	Res	Type
1	B	108	LYS
1	B	112	ARG
1	B	119	GLN
1	B	168	LEU
1	B	189	LEU
1	B	190	LEU
1	B	191	LEU
1	B	215	VAL
1	B	219	GLN
1	B	226	MET
1	B	237	LEU
1	B	241	LEU
1	B	243	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	221/249 (88%)	0.35	8 (3%) 46 55	30, 47, 86, 143	0
1	B	205/249 (82%)	0.60	27 (13%) 4 6	34, 57, 112, 143	0
All	All	426/498 (85%)	0.47	35 (8%) 14 20	30, 51, 106, 143	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	91	TRP	7.4
1	B	155	THR	6.5
1	B	47	GLY	6.2
1	B	125	ARG	4.9
1	B	152	ILE	4.3
1	B	126	GLN	4.1
1	B	123	ILE	3.9
1	A	81	LEU	3.6
1	A	82	ALA	3.6
1	B	153	ASN	3.3
1	A	83	ASP	3.3
1	B	60	LYS	3.2
1	A	85	PRO	3.2
1	B	108	LYS	3.2
1	B	124	LYS	3.2
1	B	156	THR	3.0
1	B	232	ILE	2.7
1	B	121	LEU	2.7
1	B	158	LEU	2.7
1	B	113	PHE	2.6
1	B	139	ILE	2.6
1	B	159	VAL	2.5
1	A	84	ASN	2.5
1	B	112	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	46	ARG	2.4
1	B	154	ASP	2.4
1	B	118	MET	2.4
1	B	140	ILE	2.3
1	B	22	GLY	2.2
1	A	47	GLY	2.2
1	B	26	ASN	2.1
1	B	110	LYS	2.1
1	B	105	ILE	2.0
1	B	58	GLU	2.0
1	A	235	VAL	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.