



wwPDB X-ray Structure Validation Summary Report i

Feb 26, 2014 – 09:58 PM GMT

PDB ID : 3R2W
Title : Crystal Structure of UDP-glucose Pyrophosphorylase of Homo Sapiens
Authors : Zheng, X.; Yu, Q.
Deposited on : 2011-03-14
Resolution : 3.60 Å(reported)

This is a wwPDB 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/ValidationPDFNotes.html>

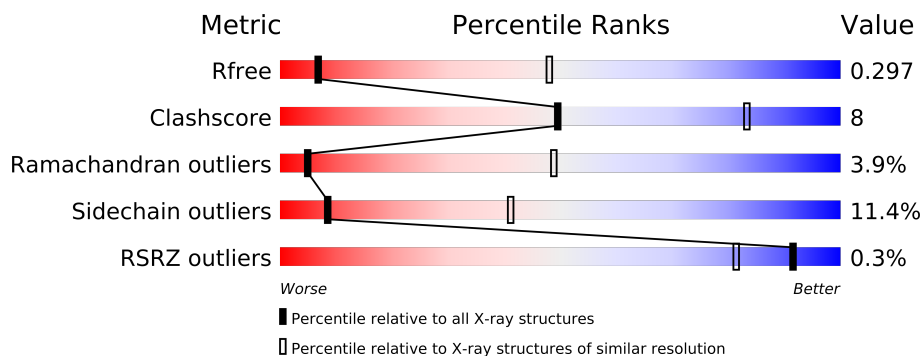
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.60 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	1020 (3.86-3.34)
Clashscore	79885	1155 (3.80-3.40)
Ramachandran outliers	78287	1109 (3.80-3.40)
Sidechain outliers	78261	1108 (3.80-3.40)
RSRZ outliers	66119	1000 (3.84-3.36)

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	528	
1	B	528	
1	C	528	
1	D	528	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 13938 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 UTP--glucose-1-phosphateuridylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	465	Total	C	N	O	S	0	0	0
			3539	2253	604	672	10			
1	B	467	Total	C	N	O	S	0	0	0
			3534	2251	605	667	11			
1	C	450	Total	C	N	O	S	0	0	0
			3299	2088	557	646	8			
1	D	468	Total	C	N	O	S	0	0	0
			3566	2275	608	672	11			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-30	MET	-	EXPRESSION TAG	UNP Q16851
A	-29	GLY	-	EXPRESSION TAG	UNP Q16851
A	-28	SER	-	EXPRESSION TAG	UNP Q16851
A	-27	SER	-	EXPRESSION TAG	UNP Q16851
A	-26	HIS	-	EXPRESSION TAG	UNP Q16851
A	-25	HIS	-	EXPRESSION TAG	UNP Q16851
A	-24	HIS	-	EXPRESSION TAG	UNP Q16851
A	-23	HIS	-	EXPRESSION TAG	UNP Q16851
A	-22	HIS	-	EXPRESSION TAG	UNP Q16851
A	-21	HIS	-	EXPRESSION TAG	UNP Q16851
A	-20	SER	-	EXPRESSION TAG	UNP Q16851
A	-19	SER	-	EXPRESSION TAG	UNP Q16851
A	-18	GLY	-	EXPRESSION TAG	UNP Q16851
A	-17	LEU	-	EXPRESSION TAG	UNP Q16851
A	-16	VAL	-	EXPRESSION TAG	UNP Q16851
A	-15	PRO	-	EXPRESSION TAG	UNP Q16851
A	-14	ARG	-	EXPRESSION TAG	UNP Q16851
A	-13	GLY	-	EXPRESSION TAG	UNP Q16851
A	-12	SER	-	EXPRESSION TAG	UNP Q16851
A	-11	HIS	-	EXPRESSION TAG	UNP Q16851
B	-30	MET	-	EXPRESSION TAG	UNP Q16851

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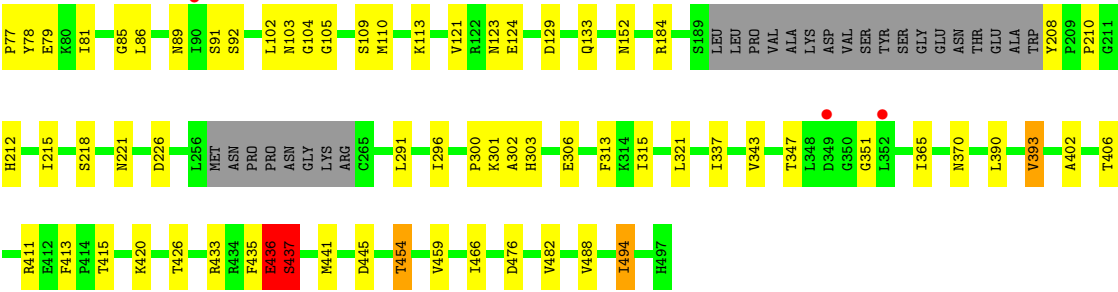
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-29	GLY	-	EXPRESSION TAG	UNP Q16851
B	-28	SER	-	EXPRESSION TAG	UNP Q16851
B	-27	SER	-	EXPRESSION TAG	UNP Q16851
B	-26	HIS	-	EXPRESSION TAG	UNP Q16851
B	-25	HIS	-	EXPRESSION TAG	UNP Q16851
B	-24	HIS	-	EXPRESSION TAG	UNP Q16851
B	-23	HIS	-	EXPRESSION TAG	UNP Q16851
B	-22	HIS	-	EXPRESSION TAG	UNP Q16851
B	-21	HIS	-	EXPRESSION TAG	UNP Q16851
B	-20	SER	-	EXPRESSION TAG	UNP Q16851
B	-19	SER	-	EXPRESSION TAG	UNP Q16851
B	-18	GLY	-	EXPRESSION TAG	UNP Q16851
B	-17	LEU	-	EXPRESSION TAG	UNP Q16851
B	-16	VAL	-	EXPRESSION TAG	UNP Q16851
B	-15	PRO	-	EXPRESSION TAG	UNP Q16851
B	-14	ARG	-	EXPRESSION TAG	UNP Q16851
B	-13	GLY	-	EXPRESSION TAG	UNP Q16851
B	-12	SER	-	EXPRESSION TAG	UNP Q16851
B	-11	HIS	-	EXPRESSION TAG	UNP Q16851
C	-30	MET	-	EXPRESSION TAG	UNP Q16851
C	-29	GLY	-	EXPRESSION TAG	UNP Q16851
C	-28	SER	-	EXPRESSION TAG	UNP Q16851
C	-27	SER	-	EXPRESSION TAG	UNP Q16851
C	-26	HIS	-	EXPRESSION TAG	UNP Q16851
C	-25	HIS	-	EXPRESSION TAG	UNP Q16851
C	-24	HIS	-	EXPRESSION TAG	UNP Q16851
C	-23	HIS	-	EXPRESSION TAG	UNP Q16851
C	-22	HIS	-	EXPRESSION TAG	UNP Q16851
C	-21	HIS	-	EXPRESSION TAG	UNP Q16851
C	-20	SER	-	EXPRESSION TAG	UNP Q16851
C	-19	SER	-	EXPRESSION TAG	UNP Q16851
C	-18	GLY	-	EXPRESSION TAG	UNP Q16851
C	-17	LEU	-	EXPRESSION TAG	UNP Q16851
C	-16	VAL	-	EXPRESSION TAG	UNP Q16851
C	-15	PRO	-	EXPRESSION TAG	UNP Q16851
C	-14	ARG	-	EXPRESSION TAG	UNP Q16851
C	-13	GLY	-	EXPRESSION TAG	UNP Q16851
C	-12	SER	-	EXPRESSION TAG	UNP Q16851
C	-11	HIS	-	EXPRESSION TAG	UNP Q16851
D	-30	MET	-	EXPRESSION TAG	UNP Q16851
D	-29	GLY	-	EXPRESSION TAG	UNP Q16851
D	-28	SER	-	EXPRESSION TAG	UNP Q16851

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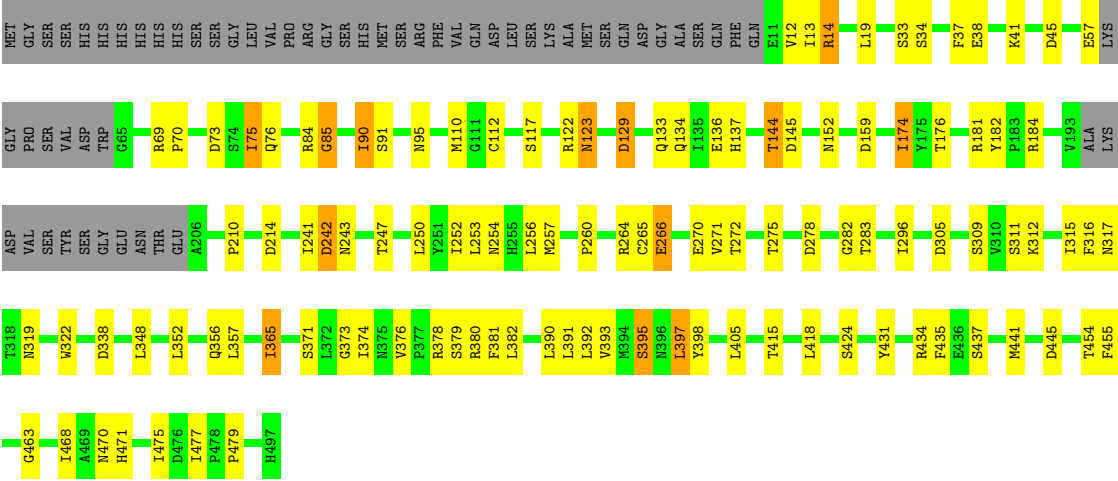
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Chain	Residue	Modelled	Actual	Comment	Reference
D	-27	SER	-	EXPRESSION TAG	UNP Q16851
D	-26	HIS	-	EXPRESSION TAG	UNP Q16851
D	-25	HIS	-	EXPRESSION TAG	UNP Q16851
D	-24	HIS	-	EXPRESSION TAG	UNP Q16851
D	-23	HIS	-	EXPRESSION TAG	UNP Q16851
D	-22	HIS	-	EXPRESSION TAG	UNP Q16851
D	-21	HIS	-	EXPRESSION TAG	UNP Q16851
D	-20	SER	-	EXPRESSION TAG	UNP Q16851
D	-19	SER	-	EXPRESSION TAG	UNP Q16851
D	-18	GLY	-	EXPRESSION TAG	UNP Q16851
D	-17	LEU	-	EXPRESSION TAG	UNP Q16851
D	-16	VAL	-	EXPRESSION TAG	UNP Q16851
D	-15	PRO	-	EXPRESSION TAG	UNP Q16851
D	-14	ARG	-	EXPRESSION TAG	UNP Q16851
D	-13	GLY	-	EXPRESSION TAG	UNP Q16851
D	-12	SER	-	EXPRESSION TAG	UNP Q16851
D	-11	HIS	-	EXPRESSION TAG	UNP Q16851



● Molecule 1: UTP--glucose-1-phosphateuridylyltransferase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	140.44Å 140.44Å 311.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.60 20.00 – 3.60	Depositor EDS
% Data completeness (in resolution range)	94.0 (20.00-3.60) 94.0 (20.00-3.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 3.62Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.247 , 0.304 0.250 , 0.297	Depositor DCC
R_{free} test set	1983 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	113.6	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.22 , 54.8	EDS
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 39283 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13938	wwPDB-VP
Average B, all atoms (Å ²)	144.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality i

5.1 Standard geometry i

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	0.61	2/3602 (0.1%)	0.72	1/4894 (0.0%)
1	B	0.58	0/3596	0.79	1/4889 (0.0%)
1	C	0.51	0/3353	0.66	0/4578
1	D	0.67	0/3630	0.83	3/4936 (0.1%)
All	All	0.60	2/14181 (0.0%)	0.75	5/19297 (0.0%)

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	311	SER	CB-OG	10.91	1.56	1.42
1	A	379	SER	CB-OG	7.32	1.51	1.42

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	382	LEU	CA-CB-CG	5.99	129.08	115.30
1	D	33	SER	N-CA-C	-5.80	95.35	111.00
1	A	444	LEU	CA-CB-CG	5.35	127.60	115.30
1	B	86	LEU	CA-CB-CG	5.28	127.45	115.30
1	D	33	SER	C-N-CA	5.09	134.43	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	91	SER	Peptide

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	3539	0	116	20	0
1	B	3534	0	89	42	0
1	C	3299	0	132	18	0
1	D	3566	0	192	33	0
All	All	13938	0	529	110	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 8.

The worst 5 of 110 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:129:ASP:O	1:D:133:GLN:CG	2.34	0.73
1:B:265:CYS:CA	1:B:266:GLU:CB	2.67	0.72
1:B:75:ILE:HA	1:B:373:GLY:O	1.90	0.71
1:D:90:ILE:N	1:D:91:SER:CA	2.54	0.70
1:B:441:MET:O	1:B:443:GLU:N	2.25	0.70

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	459/528 (87%)	376 (82%)	64 (14%)	19 (4%)	4	47
1	B	461/528 (87%)	377 (82%)	58 (13%)	26 (6%)	3	37
1	C	442/528 (84%)	363 (82%)	65 (15%)	14 (3%)	6	55
1	D	462/528 (88%)	384 (83%)	65 (14%)	13 (3%)	8	59
All	All	1824/2112 (86%)	1500 (82%)	252 (14%)	72 (4%)	5	49

5 of 72 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	SER
1	A	257	MET
1	A	263	LYS
1	B	265	CYS
1	B	266	GLU

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	381/470 (81%)	345 (91%)	36 (9%)	13	53
1	B	380/470 (81%)	331 (87%)	49 (13%)	6	34
1	C	341/470 (73%)	310 (91%)	31 (9%)	14	55
1	D	385/470 (82%)	331 (86%)	54 (14%)	5	30
All	All	1487/1880 (79%)	1317 (89%)	170 (11%)	8	41

5 of 170 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	397	LEU
1	C	226	ASP
1	D	393	VAL
1	B	424	SER
1	C	81	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	212	HIS

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	465/528 (88%)	-0.10	2 (0%) 90 76	80, 145, 200, 335	0
1	B	467/528 (88%)	-0.20	0 100 100	72, 130, 184, 231	0
1	C	450/528 (85%)	0.02	4 (0%) 81 57	85, 190, 241, 272	0
1	D	468/528 (88%)	-0.24	0 100 100	58, 114, 164, 227	0
All	All	1850/2112 (87%)	-0.13	6 (0%) 91 81	58, 135, 218, 335	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	55	LEU	2.8
1	A	208	TYR	2.7
1	C	349	ASP	2.1
1	A	209	PRO	2.1
1	C	90	ILE	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.