



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 10:04 AM EDT

PDB ID : 7YF5 / pdb_00007yf5
Title : Crystal Structure of the UDPGA Binding Domain of the Human Phase II Metabolizing Enzyme UDP-Glucuronosyltransferase 2B10
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Deposited on : 2022-07-07
Resolution : 1.53 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

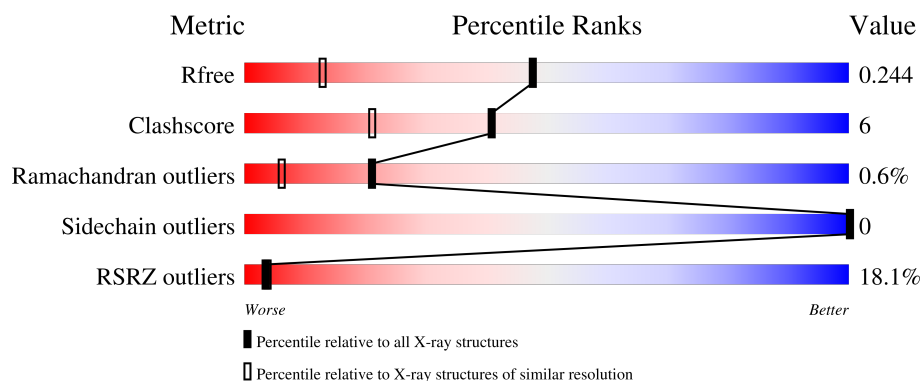
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.53 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}	180053	5890 (1.54-1.50)
Clashscore	190562	6116 (1.54-1.50)
Ramachandran outliers	187476	6002 (1.54-1.50)
Sidechain outliers	187428	5999 (1.54-1.50)
RSRZ outliers	180081	5891 (1.54-1.50)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	174	21% 86% 13% .
1	B	174	14% 84% 9% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2778 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.

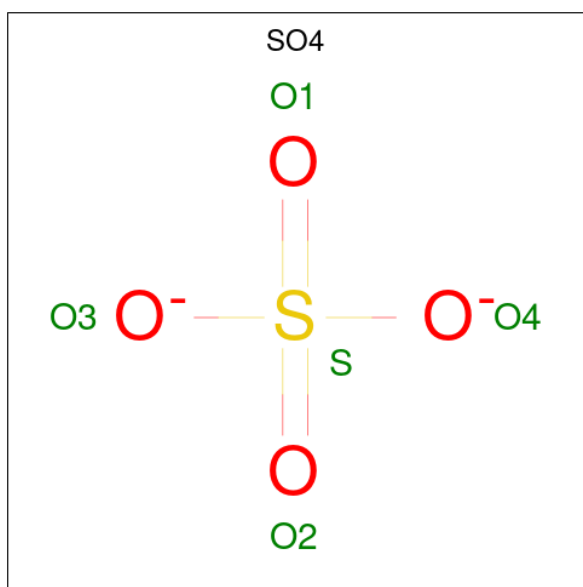
- Molecule 1 is a protein called UDP-glucuronosyltransferase 2B10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	163	Total	C	N	O	S	0	0	0
			1266	808	219	232	7			
1	A	174	Total	C	N	O	S	0	0	0
			1342	854	234	246	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	276	GLY	-	expression tag	UNP P36537
B	277	SER	-	expression tag	UNP P36537
B	278	ALA	-	expression tag	UNP P36537
B	279	MET	-	expression tag	UNP P36537
B	280	GLY	-	expression tag	UNP P36537
B	281	SER	-	expression tag	UNP P36537
A	276	GLY	-	expression tag	UNP P36537
A	277	SER	-	expression tag	UNP P36537
A	278	ALA	-	expression tag	UNP P36537
A	279	MET	-	expression tag	UNP P36537
A	280	GLY	-	expression tag	UNP P36537
A	281	SER	-	expression tag	UNP P36537

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		

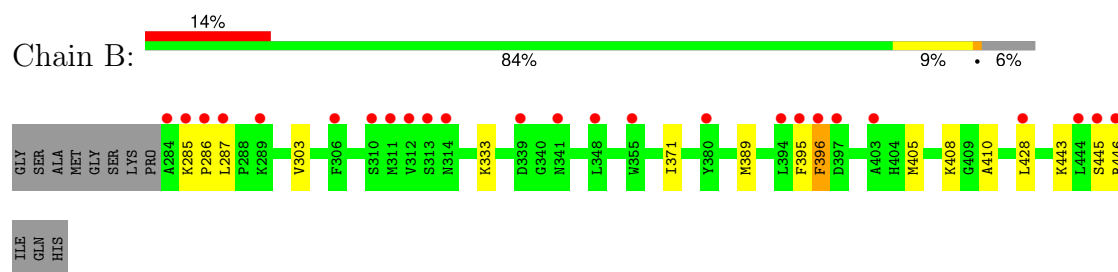
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	86	Total	O	0	0
			86	86		
3	A	79	Total	O	0	0
			79	79		

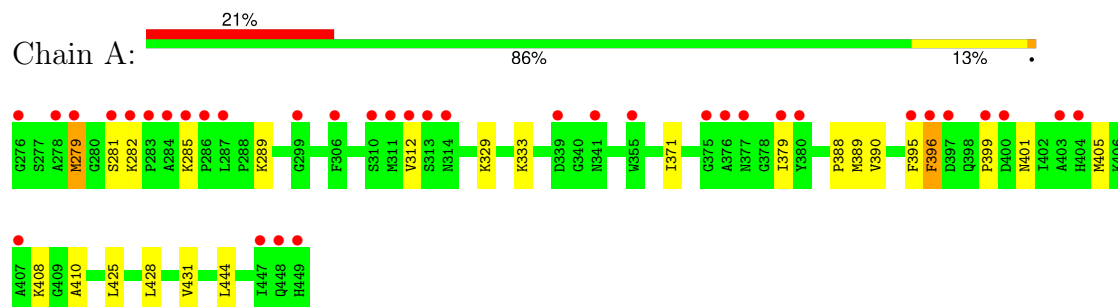
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: UDP-glucuronosyltransferase 2B10



• Molecule 1: UDP-glucuronosyltransferase 2B10



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.91Å 58.39Å 68.87Å 90.00° 98.14° 90.00°	Depositor
Resolution (Å)	34.09 – 1.53 34.09 – 1.53	Depositor EDS
% Data completeness (in resolution range)	93.9 (34.09-1.53) 93.9 (34.09-1.53)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 1.53Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.215 , 0.245 0.216 , 0.244	Depositor DCC
R_{free} test set	2412 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.156	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 30.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2778	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.38	1/1372 (0.1%)	0.51	0/1856
1	B	0.33	0/1294	0.53	0/1752
All	All	0.35	1/2666 (0.0%)	0.52	0/3608

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	279	MET	SD-CE	-6.32	1.63	1.79

There are no bond angle outliers.

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	1342	0	1354	20	0
1	B	1266	0	1278	14	0
2	B	5	0	0	0	0
3	A	79	0	0	1	0
3	B	86	0	0	0	0
All	All	2778	0	2632	34	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 (34) 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:285:LYS:HZ2	1:B:287:LEU:H	1.15	0.89
1:B:395:PHE:HB3	1:B:396:PHE:CD2	2.27	0.69
1:A:405:MET:HE3	1:A:410:ALA:HB3	1.80	0.64
1:A:289:LYS:N	1:A:289:LYS:HD3	2.15	0.60
1:A:395:PHE:HB3	1:A:396:PHE:CE1	2.37	0.60
1:B:371:ILE:HD11	1:B:428:LEU:HD21	1.84	0.59
1:B:408:LYS:HA	1:B:408:LYS:HE2	1.88	0.55
1:A:379:ILE:HD12	1:A:401:ASN:HB3	1.88	0.55
1:A:388:PRO:HB2	1:A:431:VAL:HG13	1.91	0.52
1:A:329:LYS:HD3	1:A:425:LEU:HD22	1.90	0.52
1:B:405:MET:HE3	1:B:410:ALA:HB3	1.92	0.52
1:A:389:MET:HE1	1:A:405:MET:HE2	1.90	0.52
1:B:445:SER:O	1:B:446:ARG:HG2	2.09	0.52
1:B:285:LYS:HD2	1:B:286:PRO:HD2	1.93	0.51
1:B:395:PHE:HB3	1:B:396:PHE:CE2	2.46	0.51
1:A:371:ILE:HD11	1:A:428:LEU:HD21	1.92	0.51
1:A:395:PHE:HB3	1:A:396:PHE:CD1	2.46	0.51
1:B:446:ARG:NE	1:B:446:ARG:HA	2.27	0.50
1:A:312:VAL:HG12	1:A:312:VAL:O	2.13	0.48
1:B:389:MET:HE1	1:B:405:MET:HE2	1.97	0.47
1:A:396:PHE:CD1	1:A:396:PHE:N	2.81	0.46
1:A:285:LYS:HA	1:A:285:LYS:HD2	1.79	0.46
1:A:333:LYS:NZ	3:A:504:HOH:O	2.49	0.46
1:A:281:SER:C	1:A:282:LYS:HD3	2.41	0.45
1:A:371:ILE:HD12	1:A:371:ILE:N	2.32	0.44
1:A:408:LYS:HD2	1:A:444:LEU:HD21	2.01	0.43
1:A:371:ILE:HG13	1:A:390:VAL:HB	2.01	0.42
1:A:396:PHE:O	1:A:399:PRO:HD2	2.19	0.42
1:B:303:VAL:HG12	1:B:333:LYS:HB2	2.02	0.41
1:B:443:LYS:HE3	1:B:443:LYS:HB3	1.76	0.41
1:A:279:MET:H	1:A:279:MET:HG2	1.68	0.41
1:B:396:PHE:CD2	1:B:396:PHE:N	2.87	0.41
1:B:389:MET:CE	1:B:405:MET:HE2	2.51	0.40
1:A:289:LYS:N	1:A:289:LYS:CD	2.82	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	172/174 (99%)	168 (98%)	3 (2%)	1 (1%)	21	6
1	B	161/174 (92%)	157 (98%)	3 (2%)	1 (1%)	21	6
All	All	333/348 (96%)	325 (98%)	6 (2%)	2 (1%)	21	6

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	396	PHE
1	A	396	PHE

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	145/145 (100%)	145 (100%)	0	100	100
1	B	137/145 (94%)	137 (100%)	0	100	100
All	All	282/290 (97%)	282 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	B	433	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	501	-	4,4,4	0.31	0	6,6,6	0.33	0

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 [i](#)

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	174/174 (100%)	1.23	36 (20%) 2 3	18, 29, 49, 55	0
1	B	163/174 (93%)	1.00	25 (15%) 5 6	17, 26, 47, 66	0
All	All	337/348 (96%)	1.12	61 (18%) 3 4	17, 28, 49, 66	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	396	PHE	8.0
1	B	396	PHE	6.8
1	A	395	PHE	5.5
1	A	278	ALA	5.1
1	B	284	ALA	5.1
1	A	313	SER	4.9
1	A	276	GLY	4.9
1	B	355	TRP	4.4
1	A	312	VAL	4.3
1	B	395	PHE	4.2
1	A	380	TYR	4.2
1	B	339	ASP	4.0
1	A	448	GLN	4.0
1	B	446	ARG	3.8
1	B	286	PRO	3.7
1	A	339	ASP	3.7
1	A	286	PRO	3.7
1	B	311	MET	3.6
1	A	279	MET	3.6
1	B	287	LEU	3.5
1	A	281	SER	3.4
1	B	314	ASN	3.3
1	B	312	VAL	3.3
1	A	400	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	447	ILE	3.1
1	A	449	HIS	2.9
1	B	285	LYS	2.9
1	B	313	SER	2.9
1	A	403	ALA	2.8
1	B	348	LEU	2.8
1	A	310	SER	2.8
1	B	306	PHE	2.8
1	A	287	LEU	2.8
1	A	314	ASN	2.7
1	A	299	GLY	2.7
1	B	445	SER	2.7
1	B	310	SER	2.6
1	A	306	PHE	2.6
1	A	284	ALA	2.6
1	A	404	HIS	2.5
1	A	311	MET	2.5
1	A	407	ALA	2.4
1	A	379	ILE	2.4
1	B	444	LEU	2.4
1	A	399	PRO	2.4
1	B	394	LEU	2.4
1	A	377	ASN	2.3
1	B	341	ASN	2.3
1	A	285	LYS	2.3
1	A	375	GLY	2.3
1	B	403	ALA	2.2
1	A	355	TRP	2.2
1	A	283	PRO	2.1
1	B	289	LYS	2.1
1	A	282	LYS	2.1
1	A	397	ASP	2.1
1	A	376	ALA	2.1
1	A	341	ASN	2.1
1	B	428	LEU	2.0
1	B	397	ASP	2.0
1	B	380	TYR	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 [i](#)

There are no oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	B	501	5/5	0.91	0.12	29,30,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.