

Full wwPDB X-ray Structure Validation Report (i)

May 13, 2020 - 02:37 am BST

PDB ID : 5LEV

Title: Crystal structure of human UDP-N-acetylglucosamine-dolichyl-phosphate N-

acetylglucosaminephosphotransferase (DPAGT1) (V264G mutant)

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Deposited on : 2016-06-30

Resolution : 3.20 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 (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 : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.11

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

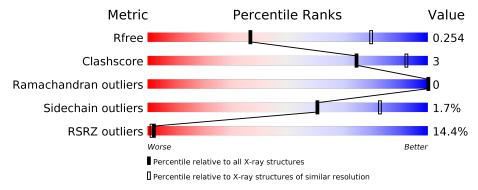
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

The reported resolution of this entry is 3.20 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar \ resolution} \\ (\#{\rm Entries, \ resolution \ \ range(\AA)}) \end{array}$
R_{free}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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 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 <=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		
			13%		
1	A	409	82%	9%	8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	UNL	A	501	_	_	-	X



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2882 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-N-acetylglucosamine--dolichyl-phosphate N-acetylglucosaminephosphotransferase.

Mol	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
1	A	375	Total 2846	C 1896	N 449	O 483	S 18	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q9H3H5
A	264	GLY	VAL	engineered mutation	UNP Q9H3H5

• Molecule 2 is UNKNOWN LIGAND (three-letter code: UNL) (formula:).

\mathbf{M}	ol	Chain	Residues	Atoms		ZeroOcc	AltConf		
2	2	A	1	Total 22	C 15	O 6	P 1	0	0

• Molecule 3 is water.

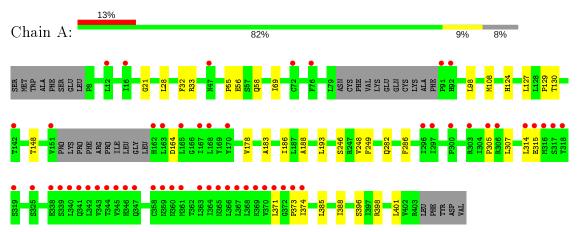
\mathbf{Mol}	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
3	A	14	Total O 14 14	0	0



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 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-N-acetylglucosamine--dolichyl-phosphate N-acetylglucosaminephosphotransfe rase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants	103.25Å 103.25Å 239.15Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.71 - 3.20	Depositor
Resolution (A)	51.63 - 3.20	EDS
% Data completeness	98.2 (26.71-3.20)	Depositor
(in resolution range)	98.2 (51.63-3.20)	EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.92 (at 3.19Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
P. P.	0.234 , 0.240	Depositor
R, R_{free}	0.251 , 0.254	DCC
R_{free} test set	666 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	122.5	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.30 , 129.8	EDS
L-test for twinning ²	$ < L >=0.49, < L^2>=0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o , F_c correlation	0.85	EDS
Total number of atoms	2882	wwPDB-VP
Average B, all atoms (Å ²)	142.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.00% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

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: UNL

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).

Mal	Chain	Bond	lengths	Bond angles	
Mol Chain	RMSZ	# Z >5	RMSZ	# Z > 5	
1	Α	0.44	0/2914	0.59	0/3975

There are no bond length outliers.

There are no bond angle outliers.

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	2846	0	2850	18	0
2	A	22	0	0	0	0
3	A	14	0	0	0	0
All	All	2882	0	2850	18	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 3.

All (18) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1			Clash overlap (Å)	
1:A:186:ILE:HD11	1:A:286:PHE:HA	1.64	0.80	

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Atom-1	Atom-2	$egin{array}{ll} ext{Interatomic} \ ext{distance } (ext{Å}) \end{array}$	Clash overlap (Å)
1:A:21:GLY:HA3	1:A:69:ILE:HD11	1.69	0.73
1:A:58:GLN:HG3	1:A:248:VAL:HB	1.80	0.64
1:A:55:PRO:HG2	1:A:248:VAL:HG12	1.80	0.61
1:A:396:SER:HA	1:A:401:LEU:HB2	1.92	0.50
1:A:56:GLU:HA	1:A:249:PHE:HB2	1.93	0.50
1:A:371:LEU:HB2	1:A:374:ILE:HD11	1.92	0.50
1:A:148:THR:HA	1:A:164:ASP:HA	1.94	0.49
1:A:188:ALA:HA	1:A:193:LEU:HD23	1.95	0.48
1:A:315:GLU:HG3	1:A:373:PRO:HB3	1.96	0.47
1:A:108:MET:HE1	1:A:129:PRO:HB3	1.97	0.47
1:A:124:HIS:HA	1:A:127:LEU:HD12	1.97	0.47
1:A:183:ALA:HB1	1:A:282:GLN:HB3	1.96	0.46
1:A:385:LEU:HA	1:A:388:ILE:HD12	1.97	0.46
1:A:129:PRO:HB2	1:A:178:VAL:HG22	1.98	0.45
1:A:305:PRO:HB2	1:A:314:LEU:HB3	2.01	0.43
1:A:33:ARG:HG3	1:A:55:PRO:HG3	2.02	0.41
1:A:28:LEU:HD22	1:A:32:PHE:HE1	1.86	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	369/409 (90%)	360 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	298/361 (82%)	293 (98%)	5 (2%)	60 83	

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	LEU
1	A	130	THR
1	A	246	SER
1	A	307	LEU
1	A	398	ARG

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

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

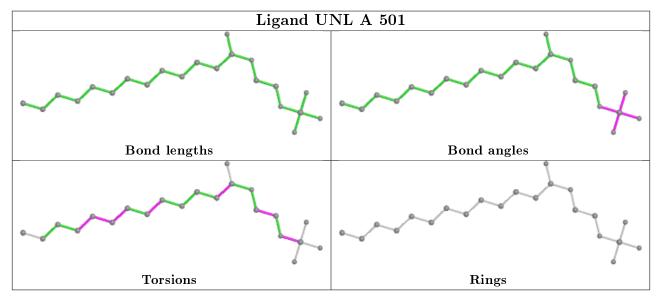
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.



The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

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, 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	$\langle { m RSRZ} \rangle$	$\#\mathrm{RSRZ}{>}2$		$OWAB(A^2)$	Q < 0.9
1	A	375/409 (91%)	0.68	54 (14%)	2 1	81, 135, 221, 248	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	315	GLU	7.3
1	A	316	MET	7.1
1	A	365	ASN	6.2
1	A	368	LEU	6.1
1	A	366	LEU	5.7
1	A	91	PRO	5.5
1	A	367	LEU	5.2
1	A	358	CYS	5.0
1	A	370	VAL	4.9
1	A	314	LEU	4.7
1	A	305	PRO	4.6
1	A	317	SER	4.6
1	A	342	LEU	4.5
1	A	345	VAL	4.5
1	A	359	ASN	4.4
1	A	318	TYR	4.2
1	A	167	ILE	4.1
1	A	369	LYS	4.1
1	A	163	LEU	4.0
1	A	339	SER	4.0
1	A	325	SER	3.8
1	A	373	PRO	3.6
1	A	344	THR	3.6
1	A	92	HIS	3.5
1	A	319	SER	3.5
1	A	16	ILE	3.5
1	A	374	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	162	HIS	3.3
1	A	12	LEU	3.3
1	A	360	ASN	3.1
1	A	372	GLY	3.0
1	A	343	VAL	2.9
1	A	371	LEU	2.9
1	A	300	PRO	2.9
1	A	363	LEU	2.8
1	A	168	LEU	2.7
1	A	47	ASN	2.7
1	A	306	ARG	2.7
1	A	142	THR	2.6
1	A	303	ARG	2.5
1	A	76	PHE	2.5
1	A	341	GLN	2.5
1	A	347	GLN	2.5
1	A	170	TYR	2.4
1	A	361	MET	2.4
1	A	297	ILE	2.4
1	A	72	CYS	2.3
1	A	340	LEU	2.2
1	A	165	LEU	2.2
1	A	364	ILE	2.2
1	A	346	HIS	2.1
1	A	151	VAL	2.0
1	A	296	ILE	2.0
1	A	338	GLU	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)

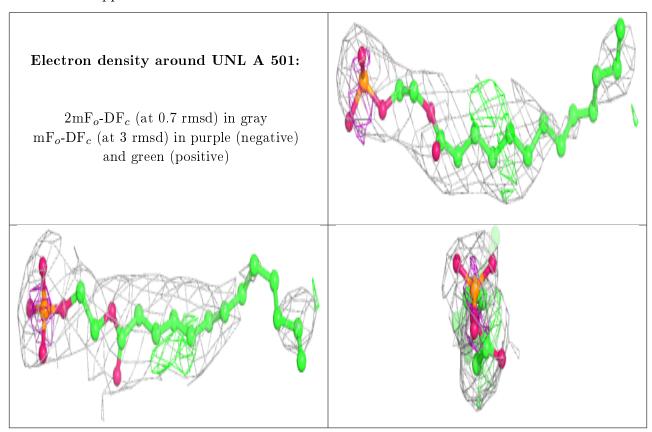
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,



median, 95^{th} 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	${f B\text{-factors}}({f \AA}^2)$	Q < 0.9
2	UNL	A	501	22/-	0.69	0.43	123,137,167,168	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers (i)

There are no such residues in this entry.

