



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 08:50 am BST

PDB ID : 5LWV
Title : Human OGT in complex with UDP and fused substrate peptide (HCF1)
Authors : Raimi, O.; Rafie, K.; Kapuria, V.; Herr, W.; van Aalten, D.
Deposited on : 2016-09-19
Resolution : 1.90 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 following versions of software and data (see [references ⓘ](#)) 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

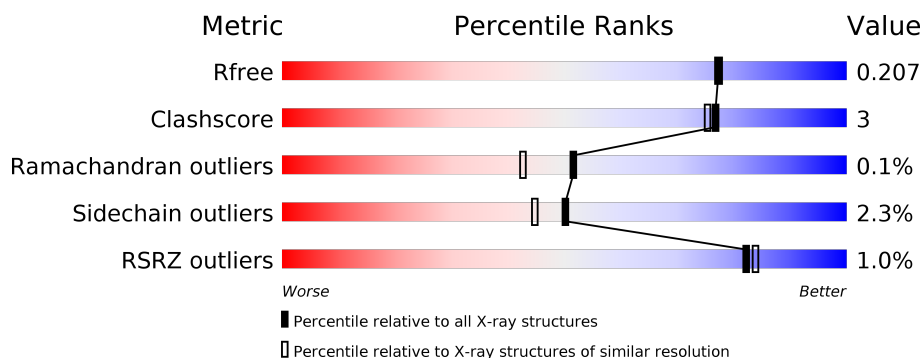
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 $\leq 5\%$. 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	749	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0;">%</div> <div style="position: absolute; top: 10px; left: 50%; transform: translate(-50%, -50%);">86%</div> <div style="position: absolute; top: 10px; right: 0;">8% • 5%</div> </div> </div>

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
3	GOL	A	1202	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6302 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.

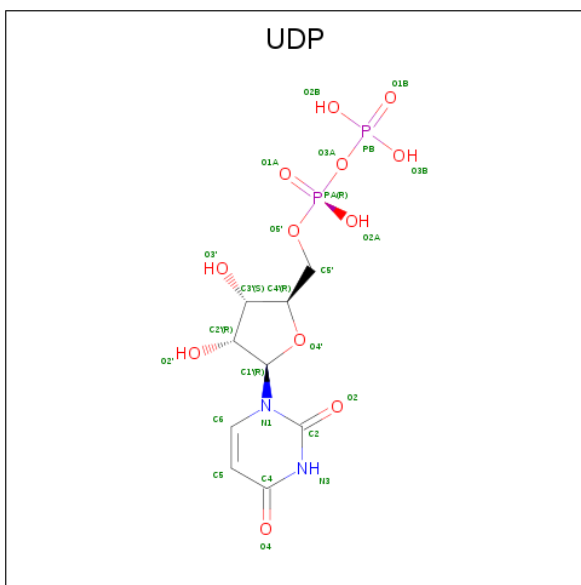
- Molecule 1 is a protein called Host cell factor 1,UDP-N-acetylglucosamine--peptide N-acetyl glucosaminyltransferase 110 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	711	Total	C	N	O	S	0	1	0
			5596	3543	979	1037	37			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	288	HIS	-	expression tag	UNP P51610
A	289	HIS	-	expression tag	UNP P51610
A	290	HIS	-	expression tag	UNP P51610
A	291	HIS	-	expression tag	UNP P51610
A	292	HIS	-	expression tag	UNP P51610
A	293	HIS	-	expression tag	UNP P51610
A	312	GLY	-	linker	UNP P51610
A	313	GLY	-	linker	UNP P51610
A	314	GLY	-	linker	UNP P51610

- Molecule 2 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		

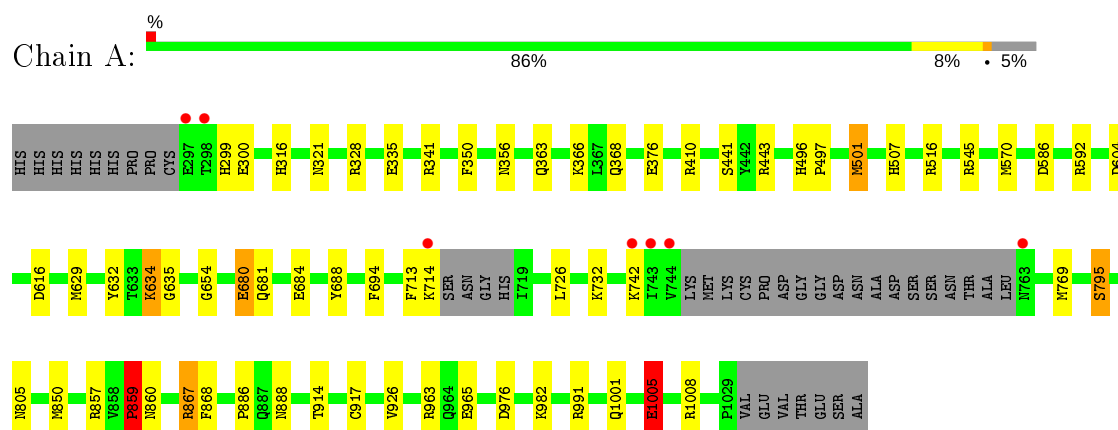
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	665	Total	O	0	0
			665	665		

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: Host cell factor 1,UDP-N-acetylglucosamine--peptide N-acetylglucosaminyltransferase 110 kDa subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	168.02Å 168.02Å 162.26Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.50 – 1.90 48.50 – 1.93	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.50-1.90) 100.0 (48.50-1.93)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.169 , 0.199 0.180 , 0.207	Depositor DCC
R_{free} test set	5042 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6302	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 3.33% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PO4, UDP

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	1.03	12/5727 (0.2%)	1.05	31/7771 (0.4%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	965	GLU	CD-OE1	-6.99	1.18	1.25
1	A	795	SER	CB-OG	-6.71	1.33	1.42
1	A	376	GLU	CD-OE2	6.47	1.32	1.25
1	A	1005	GLU	CD-OE2	6.04	1.32	1.25
1	A	335	GLU	CD-OE2	5.76	1.31	1.25
1	A	680	GLU	CD-OE1	-5.66	1.19	1.25
1	A	516	ARG	CZ-NH1	5.47	1.40	1.33
1	A	867	ARG	CZ-NH2	5.33	1.40	1.33
1	A	632	TYR	CE1-CZ	5.17	1.45	1.38
1	A	684	GLU	CD-OE1	5.07	1.31	1.25
1	A	688	TYR	CE1-CZ	5.06	1.45	1.38
1	A	441	SER	CB-OG	5.04	1.48	1.42

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	443	ARG	NE-CZ-NH2	16.82	128.71	120.30
1	A	443	ARG	NE-CZ-NH1	-13.62	113.49	120.30
1	A	545	ARG	NE-CZ-NH1	11.19	125.90	120.30
1	A	850	MET	CG-SD-CE	-9.29	85.34	100.20
1	A	991	ARG	NE-CZ-NH2	-8.35	116.13	120.30
1	A	516	ARG	NE-CZ-NH2	-8.11	116.24	120.30
1	A	1005	GLU	OE1-CD-OE2	8.01	132.92	123.30
1	A	545	ARG	NE-CZ-NH2	-7.92	116.34	120.30
1	A	867	ARG	NE-CZ-NH2	-7.71	116.44	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	516	ARG	NE-CZ-NH1	7.18	123.89	120.30
1	A	410	ARG	NE-CZ-NH2	-7.10	116.75	120.30
1	A	1008	ARG	NE-CZ-NH2	6.69	123.65	120.30
1	A	616	ASP	CB-CG-OD1	6.60	124.24	118.30
1	A	410	ARG	NE-CZ-NH1	6.49	123.55	120.30
1	A	616	ASP	CB-CG-OD2	-6.38	112.56	118.30
1	A	443	ARG	CD-NE-CZ	6.20	132.28	123.60
1	A	991	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	A	592	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	A	570	MET	CG-SD-CE	-5.61	91.22	100.20
1	A	604	ASP	CB-CG-OD2	-5.61	113.25	118.30
1	A	991	ARG	CG-CD-NE	-5.56	100.13	111.80
1	A	976	ASP	CB-CG-OD1	5.49	123.24	118.30
1	A	443	ARG	CB-CG-CD	5.47	125.82	111.60
1	A	857	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	A	963	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	A	501	MET	CG-SD-CE	-5.35	91.63	100.20
1	A	586	ASP	CB-CG-OD1	5.33	123.09	118.30
1	A	604	ASP	CB-CG-OD1	5.30	123.07	118.30
1	A	341	ARG	NE-CZ-NH1	-5.10	117.75	120.30
1	A	769	MET	CG-SD-CE	-5.07	92.09	100.20
1	A	867	ARG	CG-CD-NE	-5.01	101.29	111.80

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	5596	0	5551	34	0
2	A	25	0	11	0	0
3	A	6	0	7	5	0
4	A	10	0	0	0	0
5	A	665	0	0	6	0
All	All	6302	0	5569	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 3.

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:A:680:GLU:OE2	5:A:1302:HOH:O	1.79	1.00
1:A:860:ASN:N	5:A:1301:HOH:O	1.72	0.89
1:A:501:MET:CE	1:A:917:CYS:SG	2.61	0.88
1:A:681:GLN:NE2	5:A:1304:HOH:O	2.09	0.84
1:A:501:MET:HE1	1:A:917:CYS:SG	2.18	0.84
1:A:321:ASN:HD21	1:A:356:ASN:HD22	1.24	0.82
1:A:368:GLN:H	3:A:1202:GOL:H31	1.45	0.81
1:A:501:MET:HE3	1:A:917:CYS:SG	2.22	0.79
1:A:859:PRO:O	5:A:1303:HOH:O	2.00	0.78
1:A:501:MET:CE	1:A:917:CYS:HB2	2.27	0.64
1:A:328:ARG:HH11	1:A:363:GLN:HE22	1.47	0.62
1:A:867:ARG:HD2	5:A:1305:HOH:O	2.02	0.60
1:A:300:GLU:HB2	1:A:634:LYS:HE3	1.86	0.56
1:A:501:MET:CE	1:A:917:CYS:CB	2.83	0.56
1:A:634:LYS:HA	1:A:634:LYS:HE2	1.88	0.55
1:A:886:PRO:HB2	1:A:888:ASN:HD21	1.73	0.54
1:A:299:HIS:HA	1:A:634:LYS:HB2	1.90	0.52
1:A:328:ARG:HH11	1:A:363:GLN:NE2	2.09	0.52
1:A:886:PRO:HB2	1:A:888:ASN:ND2	2.26	0.51
1:A:316:HIS:HD2	5:A:1881:HOH:O	1.95	0.48
1:A:366:LYS:C	3:A:1202:GOL:H12	2.34	0.47
1:A:366:LYS:HA	3:A:1202:GOL:H12	1.95	0.47
1:A:634:LYS:HD3	1:A:635:GLY:H	1.81	0.46
1:A:1001:GLN:O	1:A:1005:GLU:HG2	2.15	0.46
1:A:507:HIS:CG	1:A:681:GLN:HG2	2.52	0.45
1:A:713:PHE:O	1:A:714:LYS:HB2	2.16	0.45
1:A:714:LYS:C	1:A:714:LYS:HD3	2.37	0.45
1:A:366:LYS:HA	3:A:1202:GOL:C1	2.48	0.43
1:A:634:LYS:HA	1:A:634:LYS:CE	2.49	0.43
1:A:366:LYS:CA	3:A:1202:GOL:H12	2.49	0.42
1:A:629:MET:O	1:A:654:GLY:HA3	2.19	0.42
1:A:507:HIS:ND1	1:A:681:GLN:HG2	2.35	0.42
1:A:299:HIS:HD2	1:A:300:GLU:O	2.03	0.41
1:A:496:HIS:CG	1:A:497:PRO:HD2	2.56	0.41

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	706/749 (94%)	691 (98%)	14 (2%)	1 (0%)	51	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	859	PRO

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	607/639 (95%)	593 (98%)	14 (2%)	50	45

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	350	PHE
1	A	634	LYS
1	A	694	PHE
1	A	726	LEU
1	A	732	LYS
1	A	742	LYS
1	A	795	SER
1	A	805	ASN
1	A	859	PRO
1	A	868	PHE

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Mol	Chain	Res	Type
1	A	914	THR
1	A	926	VAL
1	A	982	LYS
1	A	1005	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	299	HIS
1	A	316	HIS
1	A	321	ASN
1	A	363	GLN
1	A	402	GLN
1	A	620	GLN
1	A	888	ASN
1	A	990	GLN

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](#)

4 ligands are 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
4	PO4	A	1203	-	4,4,4	1.72	1 (25%)	6,6,6	1.98	1 (16%)
4	PO4	A	1204	-	4,4,4	1.28	0	6,6,6	1.42	2 (33%)
2	UDP	A	1201	-	20,26,26	1.04	2 (10%)	25,40,40	1.36	4 (16%)
3	GOL	A	1202	-	5,5,5	1.53	1 (20%)	5,5,5	2.73	2 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	UDP	A	1201	-	-	4/14/32/32	0/2/2/2
3	GOL	A	1202	-	-	2/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1203	PO4	P-O2	-3.12	1.45	1.54
3	A	1202	GOL	O1-C1	-3.04	1.29	1.42
2	A	1201	UDP	O4'-C1'	2.68	1.44	1.41
2	A	1201	UDP	PB-O3B	-2.68	1.44	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1202	GOL	O1-C1-C2	4.79	133.15	110.20
3	A	1202	GOL	O2-C2-C3	3.41	124.16	109.12
4	A	1203	PO4	O3-P-O2	3.27	118.46	107.97
4	A	1204	PO4	O2-P-O1	-2.67	101.14	110.89
2	A	1201	UDP	O4'-C1'-C2'	-2.35	103.48	106.93
2	A	1201	UDP	O2A-PA-O1A	2.31	123.64	112.24
2	A	1201	UDP	O2B-PB-O1B	2.03	118.64	110.68
4	A	1204	PO4	O3-P-O2	2.02	114.46	107.97
2	A	1201	UDP	C6-N1-C2	-2.02	118.00	121.20

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1201	UDP	C2'-C1'-N1-C6

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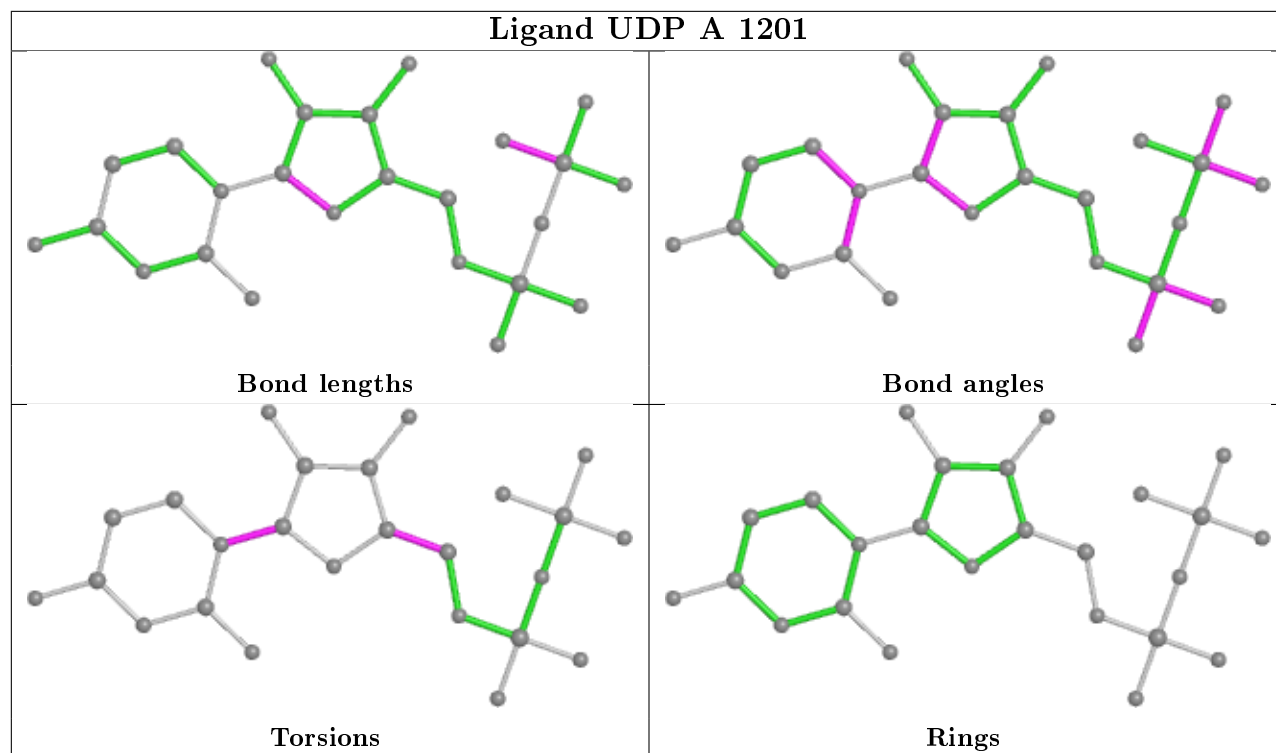
Mol	Chain	Res	Type	Atoms
2	A	1201	UDP	O4'-C1'-N1-C6
3	A	1202	GOL	C1-C2-C3-O3
3	A	1202	GOL	O2-C2-C3-O3
2	A	1201	UDP	O4'-C4'-C5'-O5'
2	A	1201	UDP	C3'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1202	GOL	5	0

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 than 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, 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	711/749 (94%)	-0.31	7 (0%) 82 84	19, 27, 47, 104	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	744	VAL	5.0
1	A	297	GLU	4.9
1	A	298	THR	3.3
1	A	743	ILE	3.1
1	A	714	LYS	2.6
1	A	742	LYS	2.1
1	A	763	ASN	2.1

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, 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.

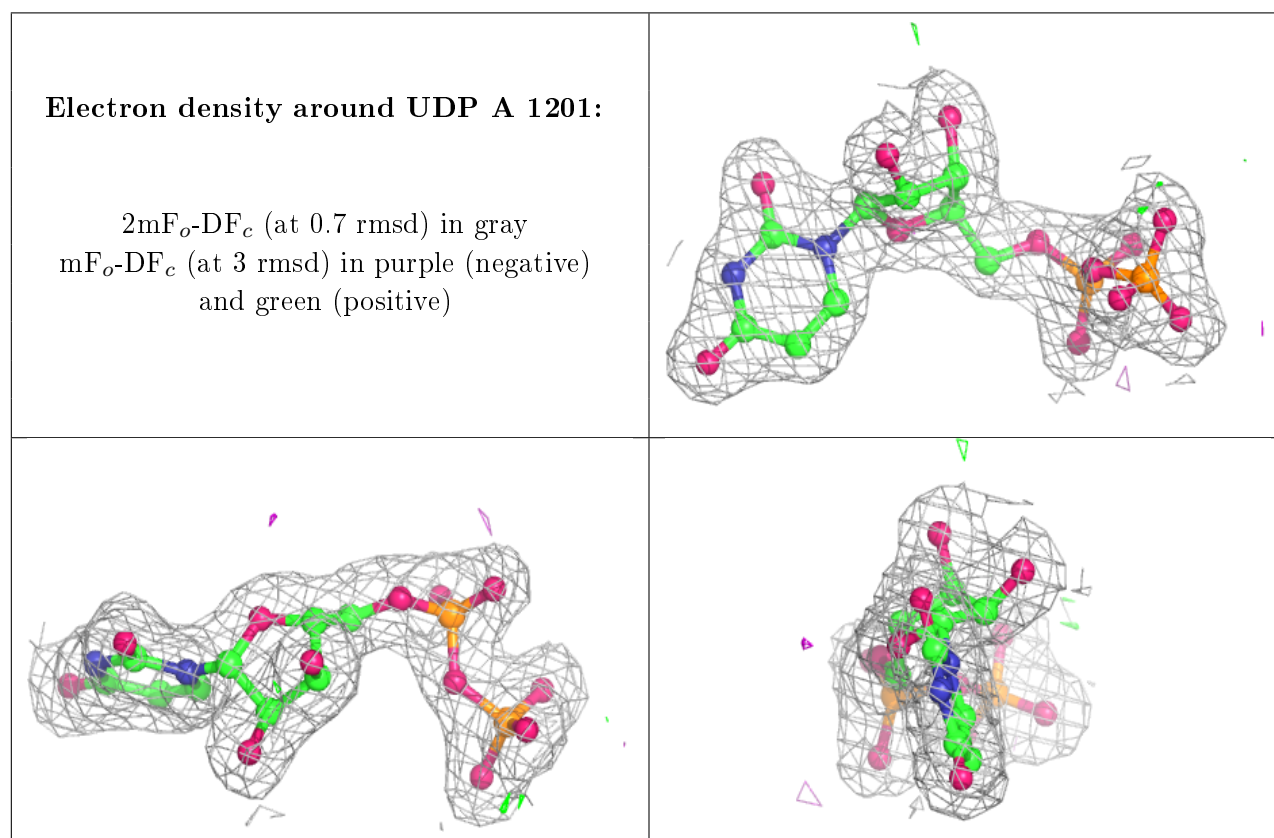
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	A	1202	6/6	0.85	0.18	27,39,43,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	UDP	A	1201	25/25	0.98	0.09	20,24,27,29	0
4	PO4	A	1204	5/5	0.98	0.14	51,55,67,67	0
4	PO4	A	1203	5/5	0.99	0.11	39,39,42,45	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.