



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 26, 2022 – 02:08 PM JST

PDB ID : 7VLB
Title : Crystal structure of UGT109A1 from Bacillus
Authors : Chen, L.Q.; Zhang, Y.
Deposited on : 2021-10-02
Resolution : 3.00 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
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) : 2.31.2

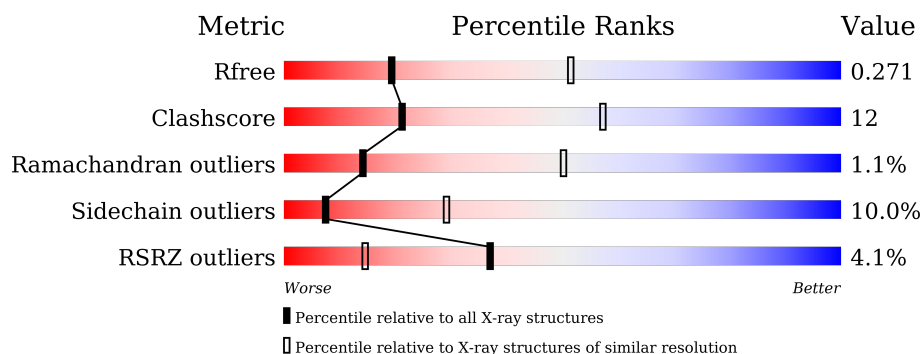
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	392	<div> <div>4%</div> <div>68%</div> <div>28%</div> <div>• •</div> </div>
1	B	392	<div> <div>4%</div> <div>63%</div> <div>21%</div> <div>5%</div> <div>10%</div> </div>

2 Entry composition [i](#)

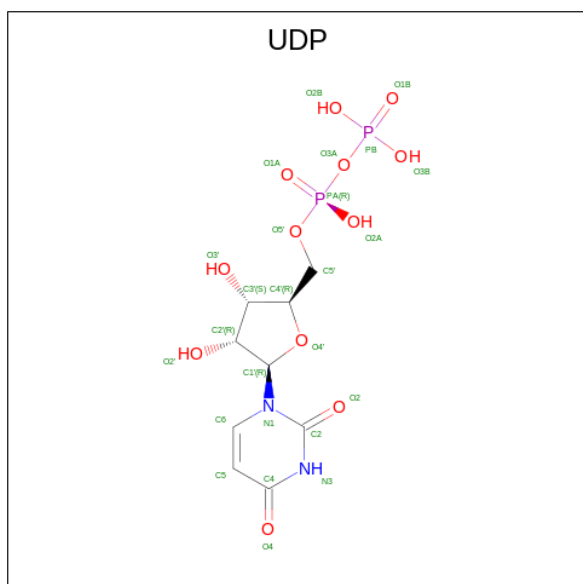
There are 3 unique types of molecules in this entry. The entry contains 11659 atoms, of which 5809 are hydrogens and 0 are deuteriums.

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-glycosyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	385	Total	C	H	N	O	S	68	0	0
			6038	1918	3021	495	586	18			
1	B	353	Total	C	H	N	O	S	63	0	0
			5527	1760	2766	453	532	16			

- Molecule 2 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total 15	O 15	0	0
3	B	7	Total 7	O 7	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	137.87Å 137.87Å 135.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.75 – 3.00 48.75 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.75-3.00) 99.9 (48.75-3.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.231 , 0.275 0.233 , 0.271	Depositor DCC
R_{free} test set	1297 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	87.6	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 71.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -h,-l,-k 0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11659	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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: 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	0.69	0/3072	0.84	2/4152 (0.0%)
1	B	0.67	0/2811	1.00	9/3796 (0.2%)
All	All	0.68	0/5883	0.92	11/7948 (0.1%)

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	A	0	2

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	111	ALA	N-CA-CB	20.06	138.19	110.10
1	B	110	LEU	N-CA-C	-13.18	75.41	111.00
1	B	114	LEU	N-CA-CB	13.05	136.51	110.40
1	B	113	LYS	CB-CA-C	-13.04	84.31	110.40
1	B	111	ALA	N-CA-C	-11.15	80.89	111.00
1	B	114	LEU	N-CA-C	-10.29	83.23	111.00
1	A	57	HIS	CB-CA-C	9.23	128.87	110.40
1	B	113	LYS	N-CA-C	-8.20	88.87	111.00
1	B	367	ASP	CB-CA-C	6.42	123.24	110.40
1	A	367	ASP	CB-CA-C	5.25	120.89	110.40
1	B	190	HIS	CA-CB-CG	5.11	122.28	113.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	272	PHE	Peptide
1	A	353	GLY	Peptide

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	3017	3021	3013	68	0
1	B	2761	2766	2756	73	0
2	A	25	11	11	0	0
2	B	25	11	11	0	0
3	A	15	0	0	0	0
3	B	7	0	0	0	0
All	All	5850	5809	5791	140	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 12.

All (140) 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:113:LYS:O	1:B:114:LEU:HG	1.70	0.90
1:A:83:GLU:O	1:A:87:ILE:HG12	1.75	0.87
1:A:202:PRO:HG3	1:A:300:MET:HB3	1.58	0.83
1:A:75:ASP:HB3	1:A:78:LEU:HD13	1.61	0.82
1:A:68:ARG:O	1:A:71:MET:HB2	1.82	0.79
1:B:240:MET:HE1	1:B:339:VAL:O	1.88	0.74
1:A:372:GLY:HA3	1:A:376:ARG:HB3	1.69	0.73
1:B:353:GLY:HA2	1:B:355:GLN:HE21	1.53	0.73
1:A:56:TYR:CD1	1:A:87:ILE:HD12	2.27	0.70
1:A:372:GLY:HA3	1:A:376:ARG:CB	2.21	0.70
1:A:246:ARG:NH2	1:A:266:ASP:OD2	2.26	0.69
1:B:336:ARG:O	1:B:339:VAL:CG1	2.42	0.67
1:B:90:GLN:O	1:B:94:LEU:HD13	1.94	0.67
1:B:242:ILE:HG23	1:B:269:PRO:HD3	1.76	0.66
1:A:340:THR:O	1:A:344:LEU:HD22	1.96	0.65
1:A:241:CYS:HB2	1:A:254:MET:HE1	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LYS:HD3	1:A:382:GLU:OE1	1.98	0.64
1:A:132:GLN:NE2	1:A:138:LEU:HD22	2.13	0.63
1:B:202:PRO:HG3	1:B:300:MET:HB3	1.82	0.61
1:A:241:CYS:CB	1:A:254:MET:HE1	2.31	0.60
1:B:213:LEU:CD2	1:B:253:ILE:HD13	2.31	0.60
1:B:91:LEU:HA	1:B:94:LEU:HB2	1.82	0.60
1:B:81:LEU:HD12	1:B:110:LEU:HD21	1.83	0.60
1:A:217:LYS:HA	1:A:251:GLN:OE1	2.00	0.60
1:B:188:ILE:HG22	1:B:189:GLN:HG2	1.83	0.60
1:A:338:GLU:O	1:A:343:LYS:HD3	2.02	0.60
1:B:336:ARG:O	1:B:339:VAL:HG12	2.02	0.59
1:B:55:ILE:N	1:B:55:ILE:HD12	2.18	0.59
1:A:367:ASP:OD2	1:B:367:ASP:OD2	2.22	0.58
1:A:104:ILE:CG2	1:A:126:LEU:HD13	2.32	0.58
1:B:5:HIS:CE1	1:B:35:THR:HG21	2.38	0.57
1:B:213:LEU:HD12	1:B:275:ARG:CZ	2.33	0.57
1:A:138:LEU:O	1:A:138:LEU:HD23	2.04	0.57
1:B:134:GLU:H	1:B:134:GLU:CD	2.08	0.57
1:B:194:ASP:OD1	1:B:196:ARG:CG	2.53	0.56
1:A:147:ILE:HG22	1:A:147:ILE:O	2.05	0.56
1:B:5:HIS:CE1	1:B:35:THR:CG2	2.88	0.56
1:B:335:GLN:O	1:B:337:GLU:O	2.25	0.55
1:A:338:GLU:O	1:A:343:LYS:CD	2.55	0.55
1:B:119:LEU:HB3	1:B:121:VAL:HG12	1.88	0.55
1:A:4:HIS:HB2	1:A:32:HIS:HD2	1.71	0.54
1:A:336:ARG:C	1:A:338:GLU:H	2.11	0.54
1:B:114:LEU:O	1:B:115:PHE:C	2.46	0.54
1:B:244:ALA:HA	1:B:345:GLN:HE21	1.74	0.53
1:A:147:ILE:O	1:A:147:ILE:CG2	2.56	0.53
1:B:194:ASP:OD1	1:B:196:ARG:HG3	2.09	0.53
1:B:81:LEU:HD13	1:B:84:SER:H	1.74	0.53
1:A:174:PRO:HG2	1:A:196:ARG:NH1	2.24	0.53
1:B:296:MET:CE	1:B:300:MET:HG2	2.39	0.52
1:A:25:GLU:O	1:A:29:GLU:HG3	2.10	0.52
1:A:38:THR:O	1:A:55:ILE:HA	2.10	0.52
1:A:174:PRO:HG2	1:A:196:ARG:CZ	2.40	0.51
1:B:115:PHE:O	1:B:119:LEU:HB2	2.09	0.51
1:B:213:LEU:HD21	1:B:253:ILE:HD13	1.92	0.51
1:B:6:ILE:HD13	1:B:27:LEU:HD13	1.92	0.51
1:B:181:PHE:HA	1:B:201:GLY:O	2.10	0.51
1:B:114:LEU:O	1:B:118:LYS:N	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:GLU:N	1:A:41:GLU:OE2	2.44	0.51
1:B:88:LEU:N	1:B:89:PRO:HD2	2.26	0.50
1:A:314:GLN:OE1	1:A:336:ARG:NH1	2.44	0.50
1:A:194:ASP:OD1	1:A:196:ARG:HG3	2.11	0.50
1:B:269:PRO:HB2	1:B:271:ASN:OD1	2.11	0.50
1:A:191:GLU:H	1:A:191:GLU:CD	2.14	0.49
1:A:184:LYS:O	1:A:190:HIS:HB3	2.12	0.49
1:B:144:LEU:O	1:B:147:ILE:HG13	2.12	0.49
1:B:194:ASP:OD1	1:B:196:ARG:HG2	2.13	0.49
1:A:211:GLU:OE2	1:A:275:ARG:HD2	2.13	0.48
1:B:83:GLU:O	1:B:87:ILE:HD13	2.13	0.48
1:B:36:TYR:HB3	1:B:46:VAL:HG11	1.95	0.48
1:A:100:PRO:O	1:A:121:VAL:HG21	2.14	0.48
1:B:213:LEU:HD12	1:B:275:ARG:NH1	2.29	0.48
1:A:268:THR:HG22	1:A:274:ILE:HD12	1.96	0.48
1:A:187:GLN:HB2	1:A:199:PHE:CE2	2.49	0.47
1:B:92:GLU:HG2	1:B:115:PHE:CE1	2.49	0.47
1:A:261:ASP:HB3	1:A:264:SER:HB3	1.96	0.47
1:B:296:MET:HE2	1:B:300:MET:HG2	1.95	0.47
1:A:88:LEU:HB3	1:A:89:PRO:HD3	1.97	0.46
1:A:158:LEU:O	1:A:158:LEU:HG	2.15	0.46
1:B:92:GLU:OE2	1:B:118:LYS:HD2	2.14	0.46
1:B:353:GLY:HA2	1:B:355:GLN:NE2	2.27	0.46
1:B:356:GLU:HG3	1:B:357:LEU:N	2.30	0.46
1:A:43:ALA:N	1:A:44:PRO:CD	2.78	0.46
1:B:324:ARG:O	1:B:328:LEU:HG	2.15	0.46
1:A:18:ASN:N	1:A:19:PRO:CD	2.78	0.46
1:A:224:LEU:HD13	1:A:283:VAL:HG11	1.98	0.46
1:B:364:MET:O	1:B:368:VAL:HB	2.16	0.46
1:A:372:GLY:CA	1:A:376:ARG:CB	2.92	0.46
1:B:55:ILE:N	1:B:55:ILE:CD1	2.79	0.46
1:B:225:ILE:HG21	1:B:254:MET:HE2	1.98	0.45
1:A:224:LEU:HD21	1:A:280:GLN:HG3	1.98	0.45
1:B:110:LEU:H	1:B:110:LEU:HD23	1.82	0.45
1:A:24:VAL:HG22	1:A:34:VAL:HG11	1.99	0.45
1:B:242:ILE:HG23	1:B:269:PRO:CD	2.44	0.45
1:A:5:HIS:ND1	1:A:100:PRO:HA	2.32	0.45
1:A:157:TYR:CE1	1:A:163:LEU:HG	2.52	0.45
1:A:13:ALA:O	1:A:17:VAL:HG13	2.17	0.45
1:B:194:ASP:OD1	1:B:194:ASP:C	2.55	0.45
1:B:43:ALA:HB3	1:B:44:PRO:HD3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:HIS:CD2	1:B:16:HIS:N	2.85	0.44
1:B:113:LYS:O	1:B:114:LEU:CG	2.55	0.44
1:B:115:PHE:O	1:B:118:LYS:HG3	2.17	0.44
1:A:241:CYS:CB	1:A:254:MET:CE	2.96	0.43
1:B:355:GLN:O	1:B:356:GLU:C	2.56	0.43
1:A:334:LEU:HD21	1:A:344:LEU:HD13	2.01	0.43
1:B:184:LYS:HG3	1:B:199:PHE:CD1	2.53	0.43
1:B:210:GLN:O	1:B:211:GLU:C	2.56	0.43
1:A:145:LYS:HG3	1:A:148:LYS:CE	2.48	0.43
1:A:382:GLU:O	1:A:385:MET:HB2	2.18	0.43
1:A:234:TRP:CE3	1:A:237:PHE:HB2	2.54	0.43
1:B:35:THR:HG21	1:B:95:TYR:HE1	1.82	0.43
1:A:335:GLN:O	1:A:339:VAL:HG23	2.19	0.43
1:A:9:ILE:HG22	1:A:109:ALA:HB1	2.01	0.43
1:A:90:GLN:O	1:A:94:LEU:CD2	2.67	0.43
1:A:222:LEU:HA	1:A:251:GLN:O	2.20	0.42
1:A:349:GLN:O	1:A:352:SER:O	2.37	0.42
1:B:371:ALA:O	1:B:376:ARG:NH1	2.50	0.42
1:B:103:ILE:HD12	1:B:116:ALA:HB2	2.01	0.42
1:B:116:ALA:C	1:B:118:LYS:H	2.22	0.42
1:B:265:LEU:O	1:B:268:THR:HG22	2.20	0.42
1:A:372:GLY:CA	1:A:376:ARG:HB2	2.49	0.42
1:A:18:ASN:N	1:A:19:PRO:HD2	2.34	0.41
1:A:115:PHE:O	1:A:119:LEU:HD13	2.20	0.41
1:A:187:GLN:HB2	1:A:199:PHE:HE2	1.85	0.41
1:A:358:LEU:O	1:A:362:LYS:HG2	2.20	0.41
1:A:155:LYS:O	1:A:158:LEU:CD2	2.69	0.41
1:B:56:TYR:CD1	1:B:87:ILE:HG13	2.56	0.41
1:B:312:ILE:O	1:B:314:GLN:NE2	2.53	0.41
1:A:103:ILE:O	1:A:123:VAL:HA	2.21	0.41
1:A:182:MET:CA	1:A:201:GLY:HA2	2.51	0.41
1:B:252:VAL:HG11	1:B:272:PHE:CE2	2.56	0.41
1:B:124:ILE:HD13	1:B:381:ILE:HG12	2.02	0.40
1:B:296:MET:HE2	1:B:300:MET:CG	2.51	0.40
1:A:194:ASP:OD1	1:A:196:ARG:CG	2.70	0.40
1:A:204:LEU:HD13	1:A:204:LEU:N	2.36	0.40
1:B:5:HIS:CE1	1:B:35:THR:HG22	2.56	0.40
1:B:14:TYR:HD1	1:B:18:ASN:HD21	1.70	0.40
1:B:110:LEU:O	1:B:110:LEU:HG	2.19	0.40
1:B:200:VAL:HA	1:B:371:ALA:O	2.21	0.40
1:B:252:VAL:HG11	1:B:272:PHE:CZ	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:LEU:HD23	1:B:320:LEU:HA	1.91	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	383/392 (98%)	346 (90%)	35 (9%)	2 (0%)	29	68
1	B	347/392 (88%)	303 (87%)	38 (11%)	6 (2%)	9	39
All	All	730/784 (93%)	649 (89%)	73 (10%)	8 (1%)	14	50

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	111	ALA
1	B	114	LEU
1	A	215	ILE
1	B	355	GLN
1	B	12	PRO
1	B	17	VAL
1	B	386	LYS
1	A	208	THR

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	331/337 (98%)	302 (91%)	29 (9%)	10	36
1	B	301/337 (89%)	267 (89%)	34 (11%)	6	24
All	All	632/674 (94%)	569 (90%)	63 (10%)	7	29

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	MET
1	A	23	LEU
1	A	41	GLU
1	A	63	ASP
1	A	66	GLN
1	A	69	GLU
1	A	70	MET
1	A	71	MET
1	A	72	GLU
1	A	79	SER
1	A	94	LEU
1	A	128	SER
1	A	137	GLN
1	A	138	LEU
1	A	149	GLU
1	A	155	LYS
1	A	158	LEU
1	A	171	LEU
1	A	190	HIS
1	A	204	LEU
1	A	213	LEU
1	A	239	LYS
1	A	249	SER
1	A	255	SER
1	A	264	SER
1	A	267	ASP
1	A	344	LEU
1	A	361	VAL
1	A	364	MET
1	B	18	ASN
1	B	35	THR
1	B	36	TYR
1	B	80	LEU
1	B	81	LEU
1	B	87	ILE

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Mol	Chain	Res	Type
1	B	91	LEU
1	B	97	ASP
1	B	110	LEU
1	B	115	PHE
1	B	118	LYS
1	B	121	VAL
1	B	128	SER
1	B	142	ASP
1	B	145	LYS
1	B	151	GLU
1	B	170	GLN
1	B	178	ASN
1	B	189	GLN
1	B	190	HIS
1	B	191	GLU
1	B	208	THR
1	B	209	GLU
1	B	210	GLN
1	B	211	GLU
1	B	213	LEU
1	B	214	LEU
1	B	255	SER
1	B	268	THR
1	B	339	VAL
1	B	342	SER
1	B	361	VAL
1	B	368	VAL
1	B	388	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	61	ASN
1	A	120	ASN
1	A	132	GLN
1	A	160	GLN
1	A	210	GLN
1	B	18	ASN
1	B	48	GLN
1	B	99	GLN
1	B	189	GLN

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Mol	Chain	Res	Type
1	B	190	HIS
1	B	345	GLN
1	B	355	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	UDP	B	401	-	24,26,26	1.23	4 (16%)	37,40,40	1.68	6 (16%)
2	UDP	A	401	-	24,26,26	1.05	1 (4%)	37,40,40	1.67	7 (18%)

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	B	401	-	-	3/16/32/32	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	UDP	A	401	-	-	3/16/32/32	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	UDP	C2-N1	2.55	1.42	1.38
2	B	401	UDP	O4'-C4'	-2.40	1.39	1.45
2	B	401	UDP	C4-N3	-2.10	1.34	1.38
2	B	401	UDP	C6-C5	2.08	1.39	1.35
2	A	401	UDP	C5-C4	-2.03	1.39	1.43

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	UDP	C4-N3-C2	-4.82	120.22	126.58
2	A	401	UDP	N3-C2-N1	4.33	120.64	114.89
2	B	401	UDP	N3-C2-N1	4.24	120.52	114.89
2	A	401	UDP	C4-N3-C2	-3.91	121.42	126.58
2	A	401	UDP	O2-C2-N1	-3.85	117.66	122.79
2	B	401	UDP	C5-C4-N3	3.84	120.59	114.84
2	A	401	UDP	O4-C4-C5	-3.51	118.98	125.16
2	B	401	UDP	O4-C4-C5	-3.42	119.15	125.16
2	B	401	UDP	O2-C2-N1	-2.41	119.58	122.79
2	B	401	UDP	C2'-C3'-C4'	2.41	107.33	102.64
2	A	401	UDP	O4-C4-N3	2.41	122.85	119.31
2	A	401	UDP	PA-O3A-PB	-2.28	125.00	132.83
2	A	401	UDP	C5-C4-N3	2.22	118.17	114.84

There are no chirality outliers.

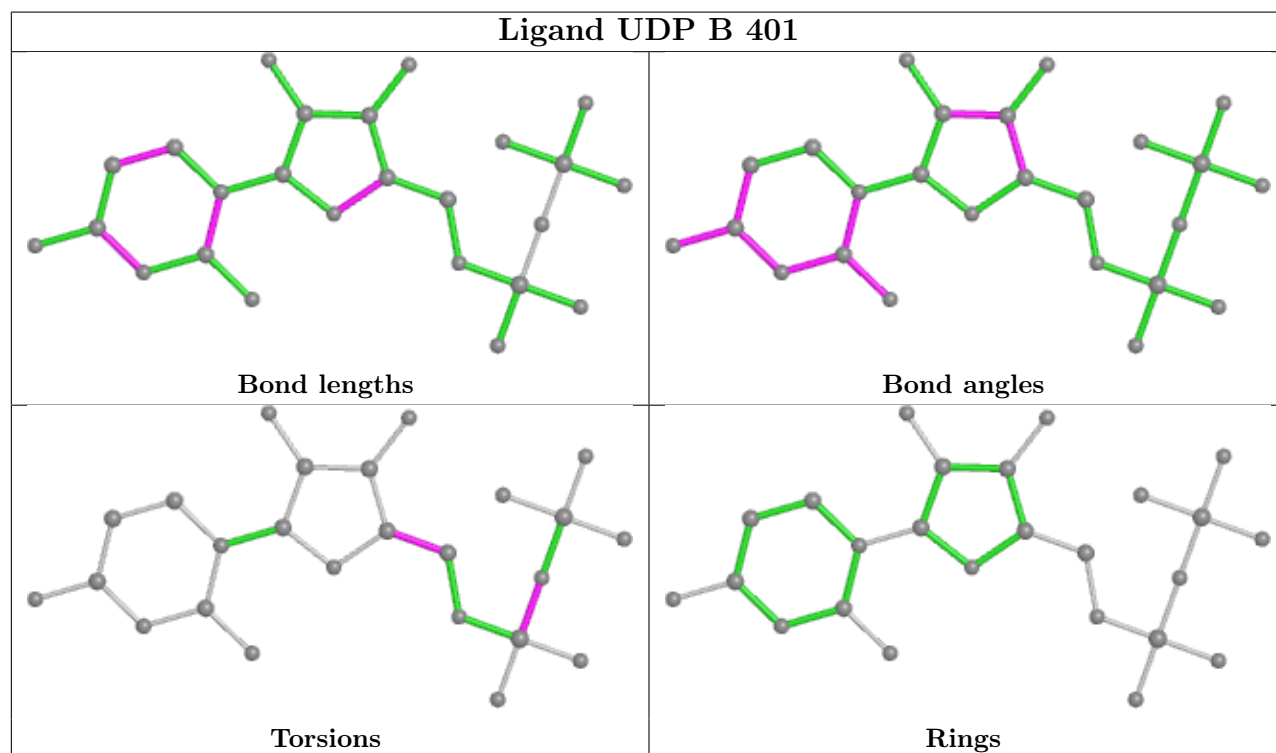
All (6) torsion outliers are listed below:

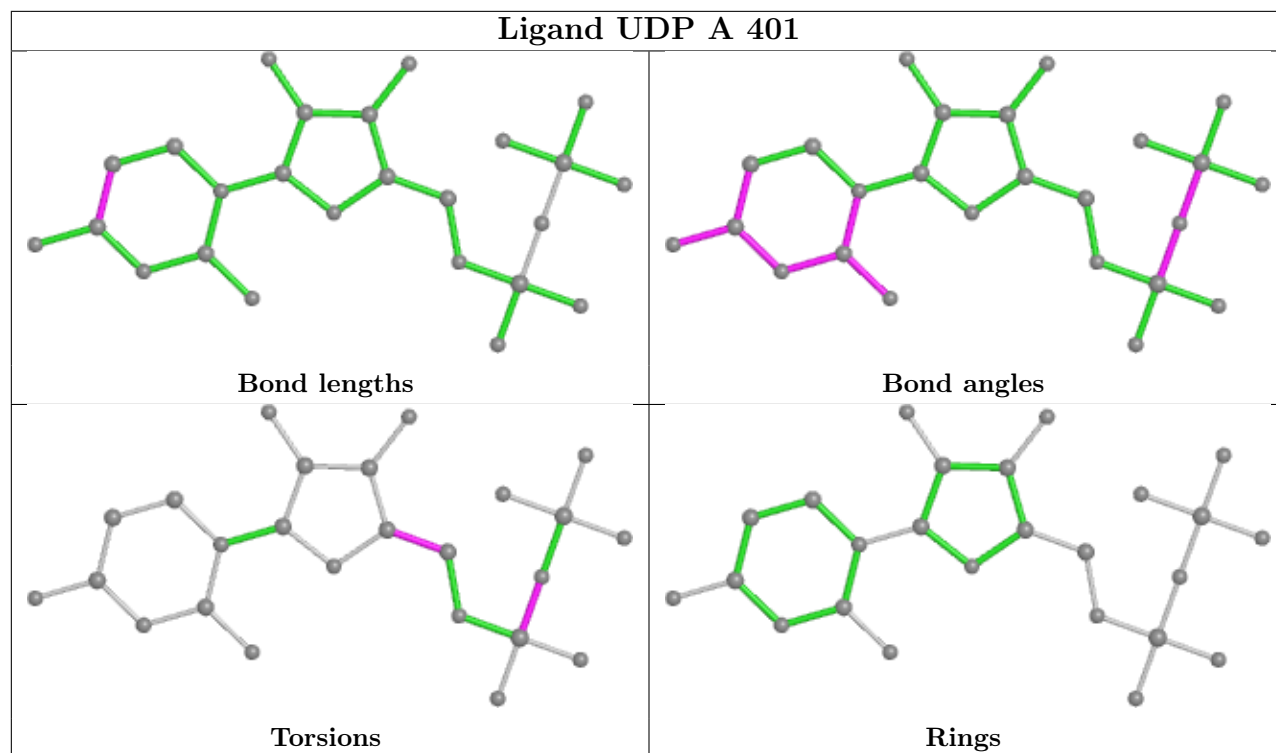
Mol	Chain	Res	Type	Atoms
2	A	401	UDP	O4'-C4'-C5'-O5'
2	A	401	UDP	PB-O3A-PA-O5'
2	B	401	UDP	PB-O3A-PA-O5'
2	A	401	UDP	C3'-C4'-C5'-O5'
2	B	401	UDP	O4'-C4'-C5'-O5'
2	B	401	UDP	C3'-C4'-C5'-O5'

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 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

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	385/392 (98%)	0.31	15 (3%)	39 15	52, 80, 173, 228	0
1	B	353/392 (90%)	0.30	15 (4%)	36 14	51, 86, 166, 209	0
All	All	738/784 (94%)	0.31	30 (4%)	37 14	51, 82, 171, 228	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	164	PRO	7.6
1	A	170	GLN	7.5
1	B	100	PRO	6.4
1	A	75	ASP	6.2
1	A	162	GLN	5.4
1	A	166	VAL	5.3
1	A	163	LEU	5.2
1	B	171	LEU	4.8
1	B	141	GLU	4.7
1	A	165	ALA	4.6
1	A	160	GLN	4.2
1	A	167	SER	4.1
1	A	158	LEU	4.0
1	B	168	PHE	3.5
1	B	152	ALA	3.4
1	B	172	ALA	2.9
1	A	153	GLU	2.9
1	B	103	ILE	2.8
1	B	102	LEU	2.7
1	A	169	GLU	2.6
1	A	171	LEU	2.6
1	B	6	ILE	2.5
1	B	154	PHE	2.5
1	B	101	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	385	MET	2.4
1	B	142	ASP	2.3
1	A	157	TYR	2.1
1	B	9	ILE	2.1
1	A	388	SER	2.1
1	B	140	ASN	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 monosaccharides in this entry.

6.4 Ligands [i](#)

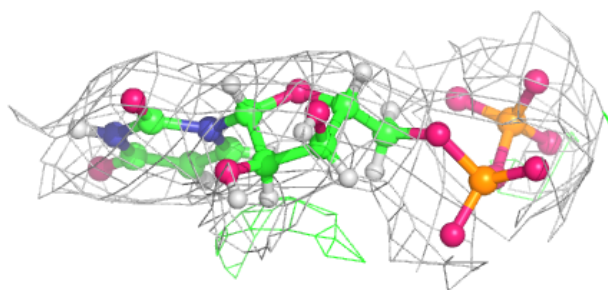
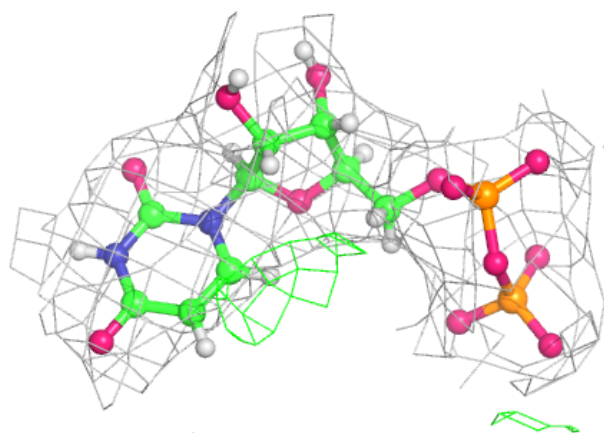
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	UDP	B	401	25/25	0.97	0.20	56,63,73,74	2
2	UDP	A	401	25/25	0.98	0.23	48,56,62,63	2

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.

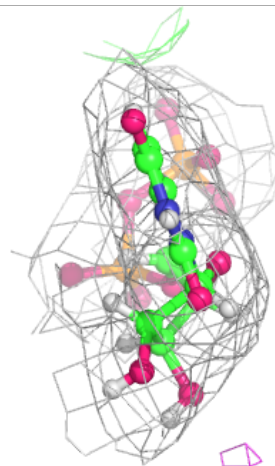
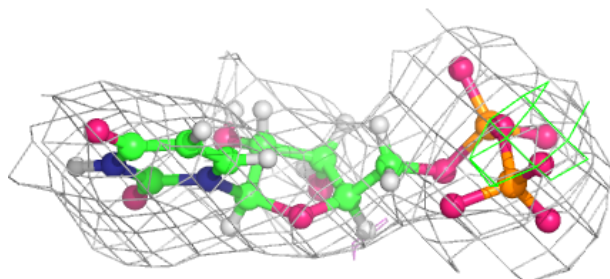
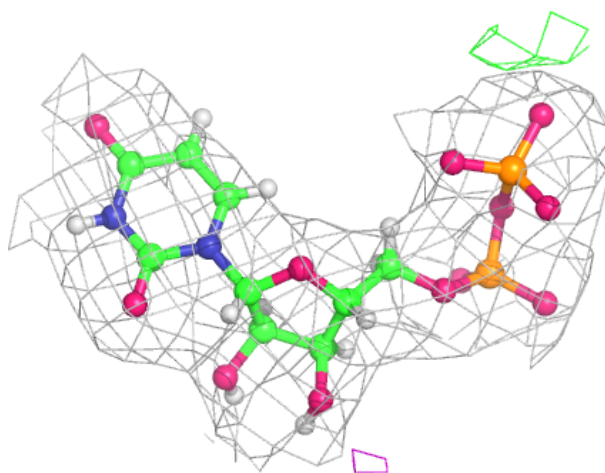
Electron density around UDP B 401:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around UDP A 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.