



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 3, 2021 – 05:40 PM EDT

PDB ID : 3INT
Title : Structure of UDP-galactopyranose mutase bound to UDP-galactose (reduced)
Authors : Gruber, T.D.; Kiessling, L.L.; Forest, K.T.
Deposited on : 2009-08-12
Resolution : 2.51 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 : 2.23.2
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) : 2.23.2

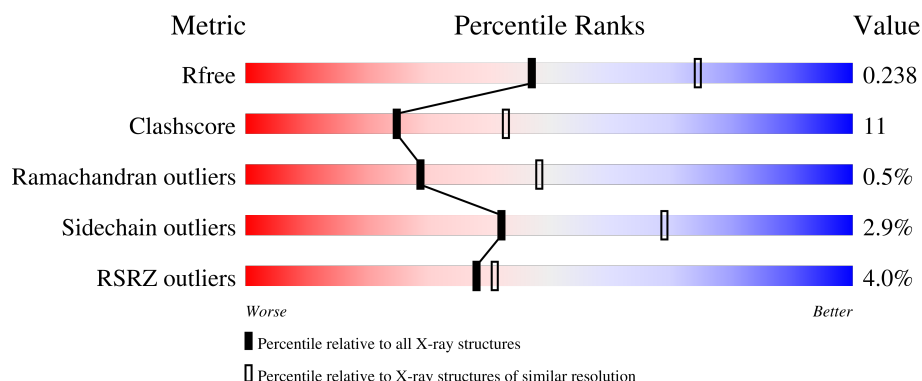
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	390	
1	B	390	

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
4	GDU	B	392	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6525 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.

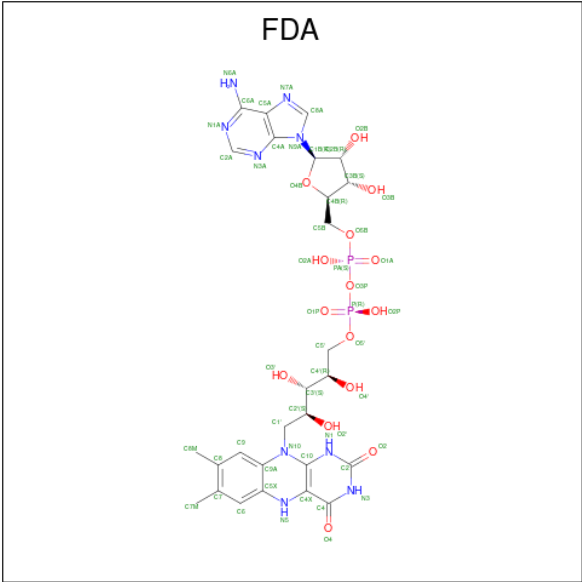
- Molecule 1 is a protein called Probable UDP-galactopyranose mutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	383	Total	C	N	O	S	0	0	0
			3119	1994	525	583	17			
1	B	374	Total	C	N	O	S	0	0	0
			3039	1942	515	565	17			

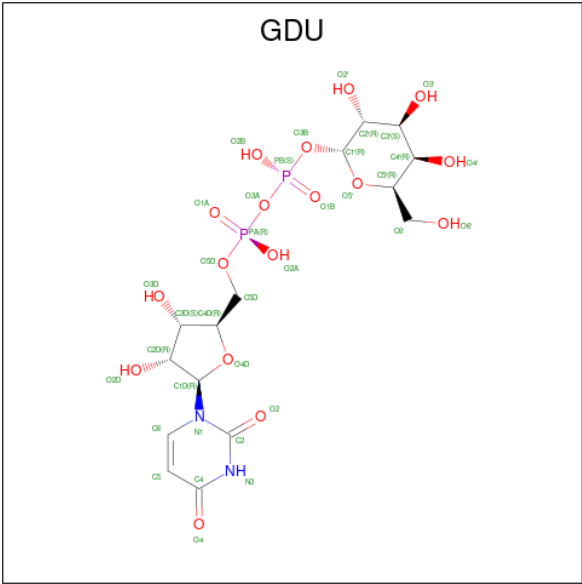
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	73	ILE	VAL	SEE REMARK 999	UNP Q48485
A	222	ASP	GLU	SEE REMARK 999	UNP Q48485
A	258	ILE	THR	SEE REMARK 999	UNP Q48485
A	372	ASP	GLU	SEE REMARK 999	UNP Q48485
A	384	GLY	ARG	engineered mutation	UNP Q48485
A	385	HIS	-	expression tag	UNP Q48485
A	386	HIS	-	expression tag	UNP Q48485
A	387	HIS	-	expression tag	UNP Q48485
A	388	HIS	-	expression tag	UNP Q48485
A	389	HIS	-	expression tag	UNP Q48485
A	390	HIS	-	expression tag	UNP Q48485
B	73	ILE	VAL	SEE REMARK 999	UNP Q48485
B	222	ASP	GLU	SEE REMARK 999	UNP Q48485
B	258	ILE	THR	SEE REMARK 999	UNP Q48485
B	372	ASP	GLU	SEE REMARK 999	UNP Q48485
B	384	GLY	ARG	engineered mutation	UNP Q48485
B	385	HIS	-	expression tag	UNP Q48485
B	386	HIS	-	expression tag	UNP Q48485
B	387	HIS	-	expression tag	UNP Q48485
B	388	HIS	-	expression tag	UNP Q48485
B	389	HIS	-	expression tag	UNP Q48485
B	390	HIS	-	expression tag	UNP Q48485

- Molecule 2 is DIHYDROFLAVINE-ADENINE DINUCLEOTIDE (three-letter code: FDA) (formula: C₂₇H₃₅N₉O₁₅P₂).



(formula: C₁₅H₂₄N₂O₁₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	B	1	36	15	2	17	2	0	0

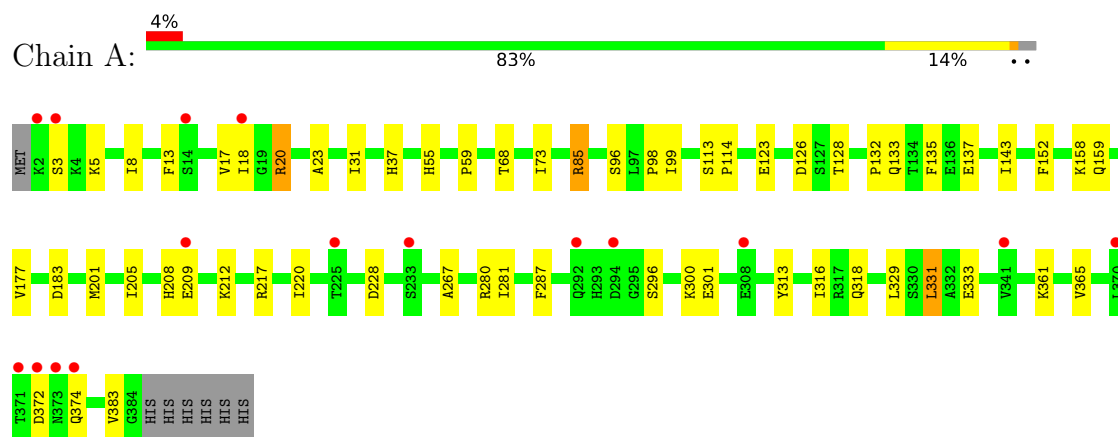
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	104	Total O	0	0
			104 104		
5	B	96	Total O	0	0
			96 96		

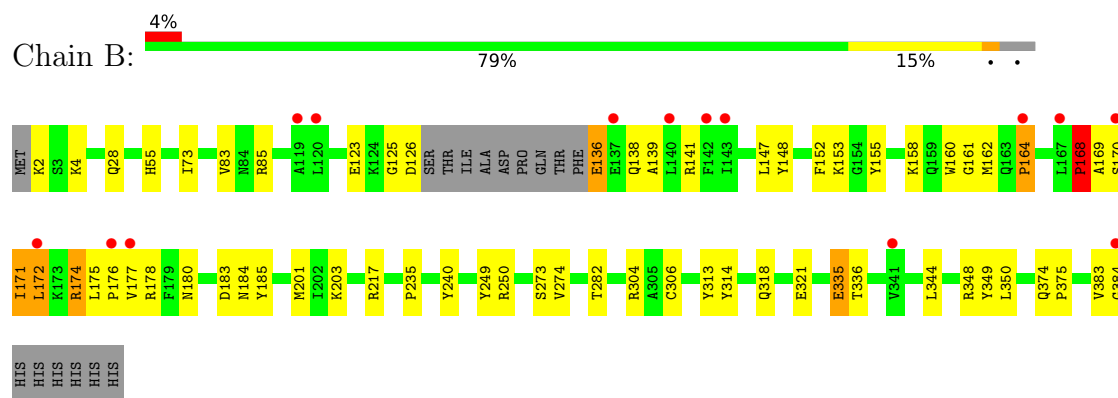
3 Residue-property plots [i](#)

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: Probable UDP-galactopyranose mutase



- Molecule 1: Probable UDP-galactopyranose mutase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	94.08Å 94.08Å 129.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.51 29.75 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.5 (30.00-2.51) 99.5 (29.75-2.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.17 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.194 , 0.247 0.188 , 0.238	Depositor DCC
R_{free} test set	1925 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.324	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.035 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6525	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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: FDA, UDP, GDU

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	0.93	3/3202 (0.1%)	0.85	3/4336 (0.1%)
1	B	0.97	0/3117	0.88	2/4216 (0.0%)
All	All	0.95	3/6319 (0.0%)	0.86	5/8552 (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	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	123	GLU	CG-CD	6.11	1.61	1.51
1	A	301	GLU	CG-CD	5.54	1.60	1.51
1	A	135	PHE	CE2-CZ	5.04	1.47	1.37

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	85	ARG	NE-CZ-NH1	8.31	124.45	120.30
1	B	168	PRO	N-CA-CB	7.19	111.93	103.30
1	A	331	LEU	CB-CG-CD2	-6.95	99.19	111.00
1	B	164	PRO	N-CA-CB	6.10	110.62	103.30
1	A	20	ARG	NE-CZ-NH2	-5.13	117.73	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	174	ARG	Sidechain

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	3119	0	2998	50	0
1	B	3039	0	2906	85	0
2	A	53	0	33	1	0
2	B	53	0	33	7	0
3	A	25	0	11	0	0
4	B	36	0	22	22	0
5	A	104	0	0	12	0
5	B	96	0	0	13	0
All	All	6525	0	6003	133	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 11.

All (133) 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:161:GLY:HA3	1:B:250:ARG:NH2	1.34	1.38
1:B:161:GLY:CA	1:B:250:ARG:HH21	1.60	1.13
1:B:161:GLY:CA	1:B:250:ARG:NH2	2.14	1.09
1:B:172:LEU:CD1	4:B:392:GDU:H1D	1.84	1.07
1:B:318:GLN:HG2	1:B:321:GLU:OE1	1.61	1.00
1:B:161:GLY:HA3	1:B:250:ARG:HH21	0.87	0.96
1:B:172:LEU:HD12	4:B:392:GDU:H1D	1.50	0.94
1:B:335:GLU:HG3	5:B:478:HOH:O	1.67	0.92
1:A:13:PHE:HE2	1:A:201:MET:HE2	1.45	0.82
4:B:392:GDU:H6	4:B:392:GDU:H5'2	1.61	0.81
1:B:161:GLY:HA3	1:B:250:ARG:HH22	1.46	0.80
1:B:240:TYR:OH	1:B:335:GLU:OE1	2.00	0.80
1:B:125:GLY:O	1:B:126:ASP:HB2	1.82	0.79
1:A:126:ASP:OD1	1:A:128:THR:HB	1.82	0.79
1:B:172:LEU:HD13	4:B:392:GDU:O2D	1.82	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:TRP:O	1:B:250:ARG:NE	2.14	0.79
1:B:172:LEU:HD13	4:B:392:GDU:C1D	2.13	0.79
1:B:318:GLN:CG	1:B:321:GLU:OE1	2.35	0.75
1:B:175:LEU:CD1	4:B:392:GDU:H5'1	2.16	0.74
1:B:172:LEU:CD1	4:B:392:GDU:C1D	2.63	0.74
1:A:85:ARG:HH21	1:B:85:ARG:HH21	1.36	0.73
1:A:159:GLN:NE2	5:A:584:HOH:O	2.06	0.73
1:B:314:TYR:OH	2:B:391:FDA:HM73	1.89	0.72
1:A:96:SER:HB2	5:A:494:HOH:O	1.89	0.71
1:B:172:LEU:HD13	4:B:392:GDU:H1D	1.66	0.71
4:B:392:GDU:H6	4:B:392:GDU:C5D	2.22	0.69
1:B:282:THR:OG1	5:B:541:HOH:O	2.11	0.68
1:B:175:LEU:HD11	4:B:392:GDU:H5'1	1.75	0.68
1:B:152:PHE:CD1	4:B:392:GDU:O2	2.47	0.67
1:A:383:VAL:HG13	5:A:499:HOH:O	1.95	0.67
1:B:178:ARG:NH2	1:B:184:ASN:O	2.24	0.66
1:A:13:PHE:CE2	1:A:201:MET:HE2	2.29	0.66
1:B:55:HIS:CG	5:B:514:HOH:O	2.49	0.65
1:B:73:ILE:HG22	1:B:201:MET:CE	2.28	0.64
1:B:177:VAL:O	1:B:177:VAL:HG12	1.96	0.63
1:B:160:TRP:C	1:B:250:ARG:HH21	2.01	0.63
1:B:168:PRO:O	1:B:169:ALA:HB3	1.98	0.62
1:A:132:PRO:HD2	5:A:496:HOH:O	1.99	0.61
1:B:73:ILE:CG2	1:B:201:MET:HE1	2.30	0.61
1:B:83:VAL:HG11	1:B:85:ARG:CZ	2.29	0.61
1:B:152:PHE:HD1	4:B:392:GDU:O2	1.81	0.61
1:B:349:TYR:CD2	1:B:349:TYR:C	2.74	0.60
1:A:85:ARG:NH2	1:B:85:ARG:HE	2.00	0.59
1:B:203:LYS:HE3	5:B:479:HOH:O	2.02	0.59
1:B:250:ARG:HD3	1:B:304:ARG:O	2.03	0.59
1:B:161:GLY:N	1:B:250:ARG:HH21	2.00	0.58
1:B:282:THR:CB	5:B:541:HOH:O	2.51	0.58
1:A:85:ARG:CZ	1:B:85:ARG:NE	2.66	0.57
1:B:158:LYS:NZ	1:B:273:SER:O	2.36	0.57
1:B:152:PHE:CE1	4:B:392:GDU:C2	2.88	0.57
1:A:85:ARG:NH2	1:B:85:ARG:HH21	2.03	0.56
1:B:73:ILE:HG22	1:B:201:MET:HE1	1.85	0.56
1:A:331:LEU:HD21	5:A:527:HOH:O	2.05	0.56
1:B:136:GLU:HB2	1:B:164:PRO:O	2.05	0.56
4:B:392:GDU:H6	4:B:392:GDU:C4D	2.36	0.55
1:A:8:ILE:HB	1:A:31:ILE:HG12	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:TRP:O	1:B:250:ARG:NH2	2.39	0.55
1:B:170:SER:O	1:B:171:ILE:HB	2.05	0.55
1:A:85:ARG:HE	1:B:85:ARG:NH2	2.05	0.55
1:A:37:HIS:HE1	5:A:455:HOH:O	1.90	0.55
1:B:85:ARG:NH1	1:B:183:ASP:OD1	2.37	0.54
2:B:391:FDA:C8	5:B:486:HOH:O	2.55	0.54
1:B:28:GLN:HG2	5:B:397:HOH:O	2.09	0.53
1:B:55:HIS:CD2	5:B:514:HOH:O	2.62	0.53
1:B:172:LEU:CD1	4:B:392:GDU:O2D	2.54	0.52
1:A:23:ALA:O	1:A:208:HIS:CE1	2.62	0.52
1:A:313:TYR:HD2	2:A:391:FDA:HM82	1.73	0.52
1:B:2:LYS:HE2	1:B:4:LYS:NZ	2.25	0.51
1:B:136:GLU:HG2	1:B:148:TYR:OH	2.11	0.51
1:A:98:PRO:HB2	5:A:394:HOH:O	2.10	0.51
1:A:143:ILE:HG22	1:A:177:VAL:HG21	1.93	0.50
1:A:73:ILE:HG22	1:A:201:MET:HE3	1.94	0.50
1:A:59:PRO:HB3	1:A:287:PHE:CZ	2.47	0.49
1:B:73:ILE:CG2	1:B:201:MET:CE	2.89	0.49
1:A:5:LYS:N	1:A:228:ASP:OD2	2.37	0.49
1:A:331:LEU:HD11	5:A:527:HOH:O	2.12	0.49
1:B:152:PHE:HE1	4:B:392:GDU:C2	2.25	0.49
1:B:160:TRP:O	1:B:250:ARG:CZ	2.61	0.49
1:A:85:ARG:NH1	1:A:183:ASP:OD1	2.26	0.49
1:A:13:PHE:HE2	1:A:201:MET:CE	2.22	0.48
1:B:160:TRP:C	1:B:250:ARG:NH2	2.65	0.48
1:B:313:TYR:HD2	2:B:391:FDA:HM82	1.78	0.48
1:A:372:ASP:O	1:A:374:GLN:HG2	2.13	0.48
1:A:85:ARG:NE	1:B:85:ARG:CZ	2.77	0.47
1:B:158:LYS:HD3	1:B:274:VAL:HA	1.96	0.47
1:B:168:PRO:O	1:B:169:ALA:CB	2.61	0.47
1:A:37:HIS:CE1	5:A:455:HOH:O	2.66	0.47
1:A:99:ILE:HD11	1:A:152:PHE:HE1	1.79	0.47
1:B:384:GLY:HA3	5:B:460:HOH:O	2.13	0.47
1:B:174:ARG:NH2	4:B:392:GDU:O2A	2.47	0.47
1:B:203:LYS:CE	5:B:479:HOH:O	2.63	0.47
1:B:374:GLN:HB3	1:B:375:PRO:HD2	1.97	0.46
1:A:143:ILE:HG22	1:A:177:VAL:CG2	2.46	0.46
1:A:281:ILE:HG12	1:A:300:LYS:HG2	1.98	0.46
1:B:314:TYR:CD1	1:B:314:TYR:N	2.83	0.46
1:B:83:VAL:HG11	1:B:85:ARG:NH2	2.31	0.46
1:A:208:HIS:ND1	1:A:209:GLU:N	2.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:391:FDA:HM82	5:B:486:HOH:O	2.15	0.46
1:A:99:ILE:HD11	1:A:152:PHE:CE1	2.50	0.46
1:A:133:GLN:N	1:A:137:GLU:OE1	2.35	0.45
1:A:17:VAL:O	1:A:18:ILE:C	2.53	0.45
1:B:250:ARG:HH11	1:B:304:ARG:C	2.20	0.45
1:A:85:ARG:NH2	1:B:85:ARG:NE	2.64	0.44
1:B:235:PRO:HA	1:B:344:LEU:O	2.18	0.44
1:B:160:TRP:C	1:B:250:ARG:HE	2.16	0.44
1:B:172:LEU:HD13	4:B:392:GDU:C2D	2.47	0.44
1:B:55:HIS:CE1	5:B:514:HOH:O	2.69	0.44
1:B:250:ARG:NH1	1:B:304:ARG:C	2.71	0.44
1:A:85:ARG:NH2	1:B:85:ARG:NH2	2.65	0.43
1:A:85:ARG:NE	1:B:85:ARG:NH2	2.66	0.43
1:A:217:ARG:HD2	5:A:443:HOH:O	2.17	0.43
1:B:155:TYR:CD2	4:B:392:GDU:C2	3.02	0.43
1:B:249:TYR:O	1:B:306:CYS:CB	2.66	0.43
1:A:20:ARG:HD3	1:A:205:ILE:O	2.19	0.42
1:B:180:ASN:OD1	1:B:180:ASN:C	2.58	0.42
1:A:55:HIS:HD2	5:A:471:HOH:O	2.02	0.42
2:B:391:FDA:H1'1	2:B:391:FDA:H9	1.75	0.42
2:B:391:FDA:C8M	5:B:486:HOH:O	2.68	0.42
1:B:160:TRP:CE3	1:B:160:TRP:HA	2.54	0.41
1:B:314:TYR:HE2	4:B:392:GDU:O2B	2.02	0.41
1:A:85:ARG:CZ	1:B:85:ARG:CZ	2.98	0.41
1:A:99:ILE:CD1	1:A:152:PHE:CE1	3.03	0.41
1:A:329:LEU:HD23	1:A:329:LEU:HA	1.88	0.41
1:A:361:LYS:O	1:A:365:VAL:HG23	2.20	0.41
1:B:176:PRO:HG2	1:B:178:ARG:HH21	1.84	0.41
1:A:113:SER:HB2	1:A:114:PRO:HD2	2.02	0.41
2:B:391:FDA:O4	4:B:392:GDU:O4'	2.36	0.41
1:A:316:ILE:HG22	1:A:318:GLN:HG3	2.03	0.41
1:A:280:ARG:NH1	5:A:583:HOH:O	2.41	0.41
1:B:152:PHE:CE1	4:B:392:GDU:O2	2.74	0.41
1:A:99:ILE:CD1	1:A:152:PHE:HE1	2.35	0.40
1:A:267:ALA:HB1	1:A:287:PHE:CE2	2.55	0.40
1:B:348:ARG:NH2	1:B:350:LEU:HD11	2.36	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	381/390 (98%)	368 (97%)	13 (3%)	0	100	100
1	B	370/390 (95%)	346 (94%)	20 (5%)	4 (1%)	14	26
All	All	751/780 (96%)	714 (95%)	33 (4%)	4 (0%)	29	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171	ILE
1	B	153	LYS
1	B	139	ALA
1	B	168	PRO

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	336/343 (98%)	329 (98%)	7 (2%)	53	78
1	B	323/343 (94%)	311 (96%)	12 (4%)	34	60
All	All	659/686 (96%)	640 (97%)	19 (3%)	42	69

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	68	THR

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Mol	Chain	Res	Type
1	A	158	LYS
1	A	212	LYS
1	A	220	ILE
1	A	296	SER
1	A	333	GLU
1	B	123	GLU
1	B	136	GLU
1	B	138	GLN
1	B	141	ARG
1	B	147	LEU
1	B	162	MET
1	B	172	LEU
1	B	185	TYR
1	B	217	ARG
1	B	335	GLU
1	B	336	THR
1	B	383	VAL

Sometimes 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 monosaccharides 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	$\# Z > 2$
3	UDP	A	392	-	20,26,26	1.47	3 (15%)	25,40,40	1.21	2 (8%)
2	FDA	A	391	-	51,58,58	1.90	10 (19%)	60,89,89	2.04	16 (26%)
2	FDA	B	391	-	51,58,58	1.85	7 (13%)	60,89,89	2.04	8 (13%)
4	GDU	B	392	-	31,38,38	0.84	2 (6%)	41,58,58	1.07	2 (4%)

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
3	UDP	A	392	-	-	2/14/32/32	0/2/2/2
2	FDA	A	391	-	-	2/30/50/50	0/6/6/6
2	FDA	B	391	-	-	2/30/50/50	0/6/6/6
4	GDU	B	392	-	-	9/21/59/59	0/3/3/3

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	391	FDA	C4X-N5	7.04	1.43	1.33
2	B	391	FDA	C10-N1	6.46	1.41	1.33
2	B	391	FDA	C4X-N5	5.97	1.41	1.33
2	A	391	FDA	C4-N3	4.97	1.41	1.33
2	B	391	FDA	C4-N3	4.64	1.41	1.33
2	B	391	FDA	C2A-N3A	4.55	1.39	1.32
2	A	391	FDA	C10-N1	4.37	1.38	1.33
2	A	391	FDA	C2A-N3A	4.17	1.38	1.32
2	A	391	FDA	C5X-N5	3.86	1.41	1.35
2	B	391	FDA	C5X-N5	3.66	1.41	1.35
3	A	392	UDP	C6-N1	3.23	1.39	1.35
3	A	392	UDP	C4-N3	3.12	1.38	1.33
2	A	391	FDA	C9A-N10	3.01	1.42	1.38
3	A	392	UDP	O4'-C4'	-2.45	1.39	1.45
2	A	391	FDA	C2A-N1A	2.37	1.38	1.33
2	A	391	FDA	C1'-N10	2.34	1.50	1.48
4	B	392	GDU	C6-N1	-2.33	1.33	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	391	FDA	C5'-C4'	2.26	1.55	1.51
2	B	391	FDA	C2B-C1B	-2.23	1.50	1.53
4	B	392	GDU	C2-N3	-2.18	1.33	1.38
2	A	391	FDA	C2B-C1B	-2.10	1.50	1.53
2	B	391	FDA	C1'-N10	2.01	1.50	1.48

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	391	FDA	N3A-C2A-N1A	-7.09	117.60	128.68
2	A	391	FDA	N3A-C2A-N1A	-6.88	117.93	128.68
2	B	391	FDA	C5X-C9A-N10	6.59	122.49	117.72
2	A	391	FDA	C10-C4X-N5	-6.56	116.72	121.26
2	B	391	FDA	C4-N3-C2	5.98	120.19	115.14
2	B	391	FDA	C9A-N10-C10	-4.86	115.55	121.91
2	A	391	FDA	C4X-N5-C5X	4.38	121.15	116.77
2	B	391	FDA	C1'-N10-C9A	4.21	121.60	118.29
3	A	392	UDP	O3B-PB-O3A	3.79	117.34	104.64
2	A	391	FDA	C7-C6-C5X	-3.68	116.01	121.22
2	A	391	FDA	O4B-C1B-C2B	-3.29	102.11	106.93
2	A	391	FDA	C4-C4X-N5	3.25	122.31	118.60
2	A	391	FDA	P-O3P-PA	-3.21	121.80	132.83
2	A	391	FDA	C6-C5X-C9A	3.07	123.07	119.05
2	B	391	FDA	C1'-N10-C10	3.05	121.14	118.41
2	A	391	FDA	C4-N3-C2	2.98	117.66	115.14
2	A	391	FDA	C5X-C9A-N10	2.92	119.83	117.72
4	B	392	GDU	PB-O3A-PA	-2.89	122.90	132.83
3	A	392	UDP	PA-O3A-PB	-2.87	122.99	132.83
2	A	391	FDA	C9A-N10-C10	-2.83	118.20	121.91
2	A	391	FDA	C1'-N10-C10	2.82	120.93	118.41
2	A	391	FDA	C9A-C5X-N5	-2.64	118.24	122.36
2	A	391	FDA	C4X-C4-N3	-2.61	119.87	123.43
2	B	391	FDA	C3B-C2B-C1B	2.53	104.79	100.98
2	A	391	FDA	O3'-C3'-C2'	-2.23	103.42	108.81
2	A	391	FDA	C4X-C10-N10	2.23	122.59	120.30
4	B	392	GDU	O3A-PB-O3B	2.22	106.95	102.48
2	B	391	FDA	P-O3P-PA	-2.12	125.54	132.83

There are no chirality outliers.

All (15) torsion outliers are listed below:

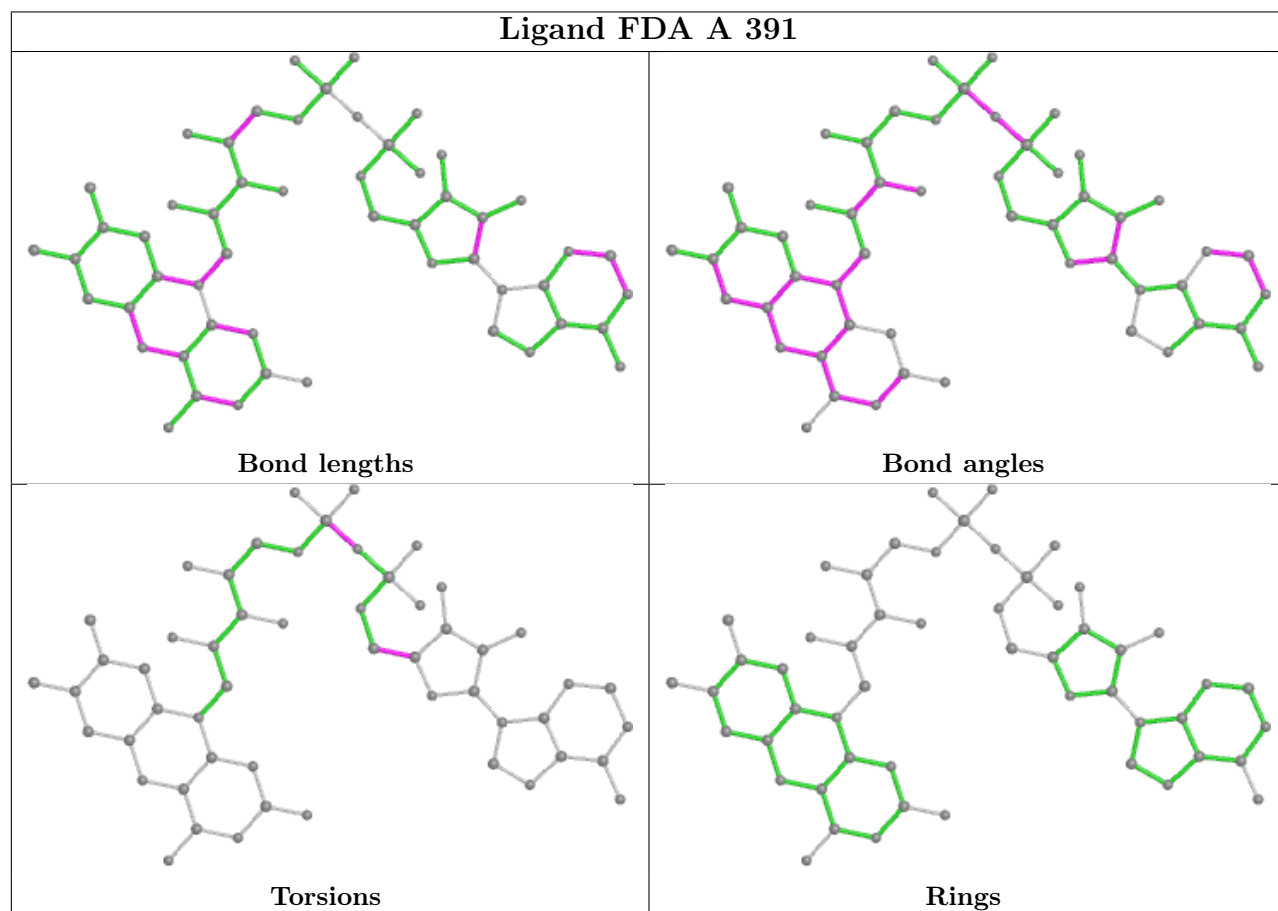
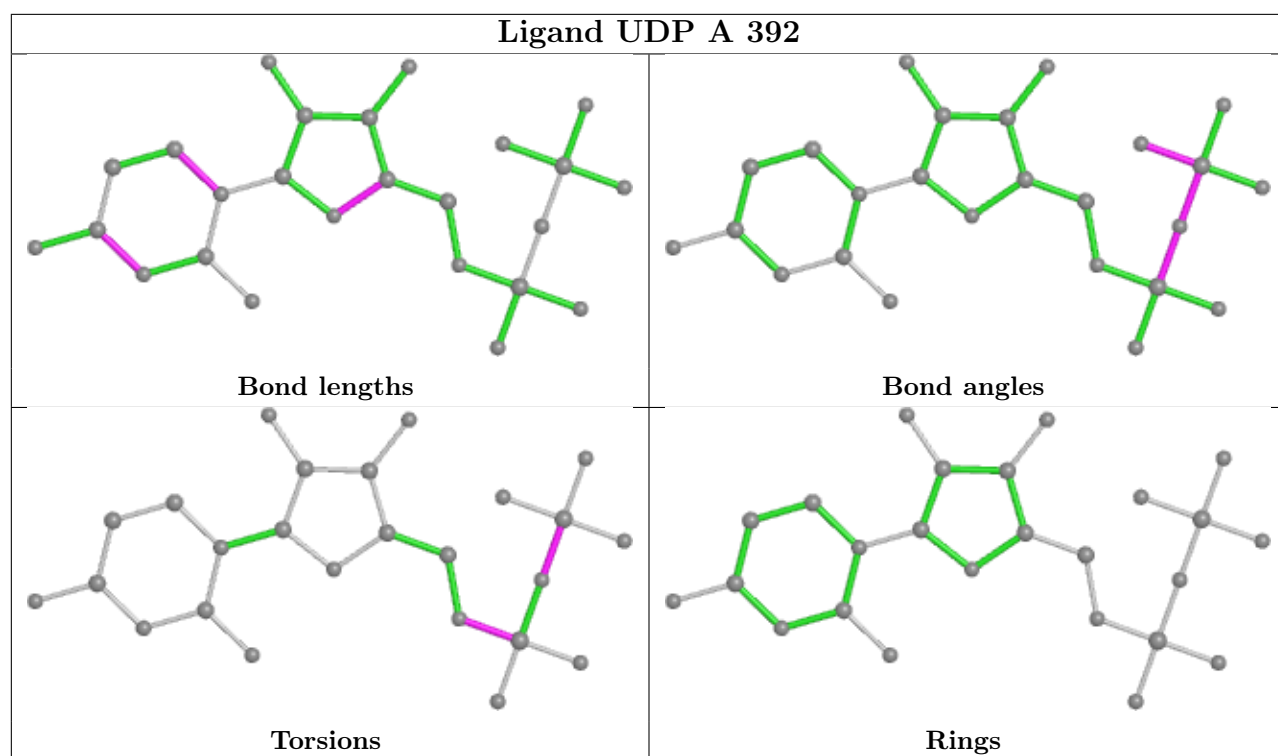
Mol	Chain	Res	Type	Atoms
2	A	391	FDA	PA-O3P-P-O5'
3	A	392	UDP	PA-O3A-PB-O2B
4	B	392	GDU	C5D-O5D-PA-O1A
4	B	392	GDU	C5D-O5D-PA-O2A
4	B	392	GDU	O5'-C1'-O3B-PB
2	B	391	FDA	O4B-C4B-C5B-O5B
4	B	392	GDU	O5'-C5'-C6'-O6'
2	B	391	FDA	C3B-C4B-C5B-O5B
4	B	392	GDU	C4'-C5'-C6'-O6'
4	B	392	GDU	PA-O3A-PB-O3B
4	B	392	GDU	C5D-O5D-PA-O3A
4	B	392	GDU	O4D-C4D-C5D-O5D
2	A	391	FDA	O4B-C4B-C5B-O5B
3	A	392	UDP	C5'-O5'-PA-O1A
4	B	392	GDU	C3D-C4D-C5D-O5D

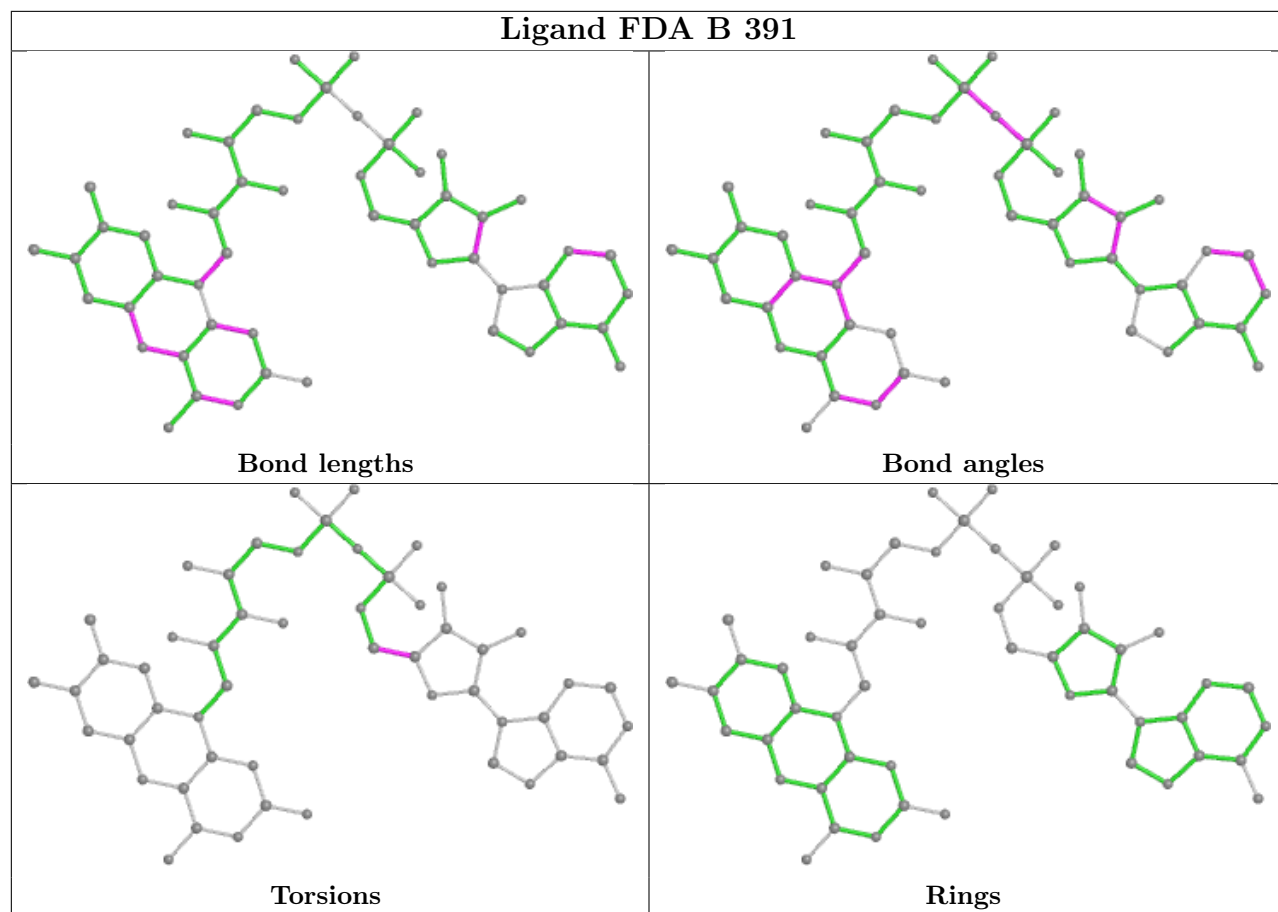
There are no ring outliers.

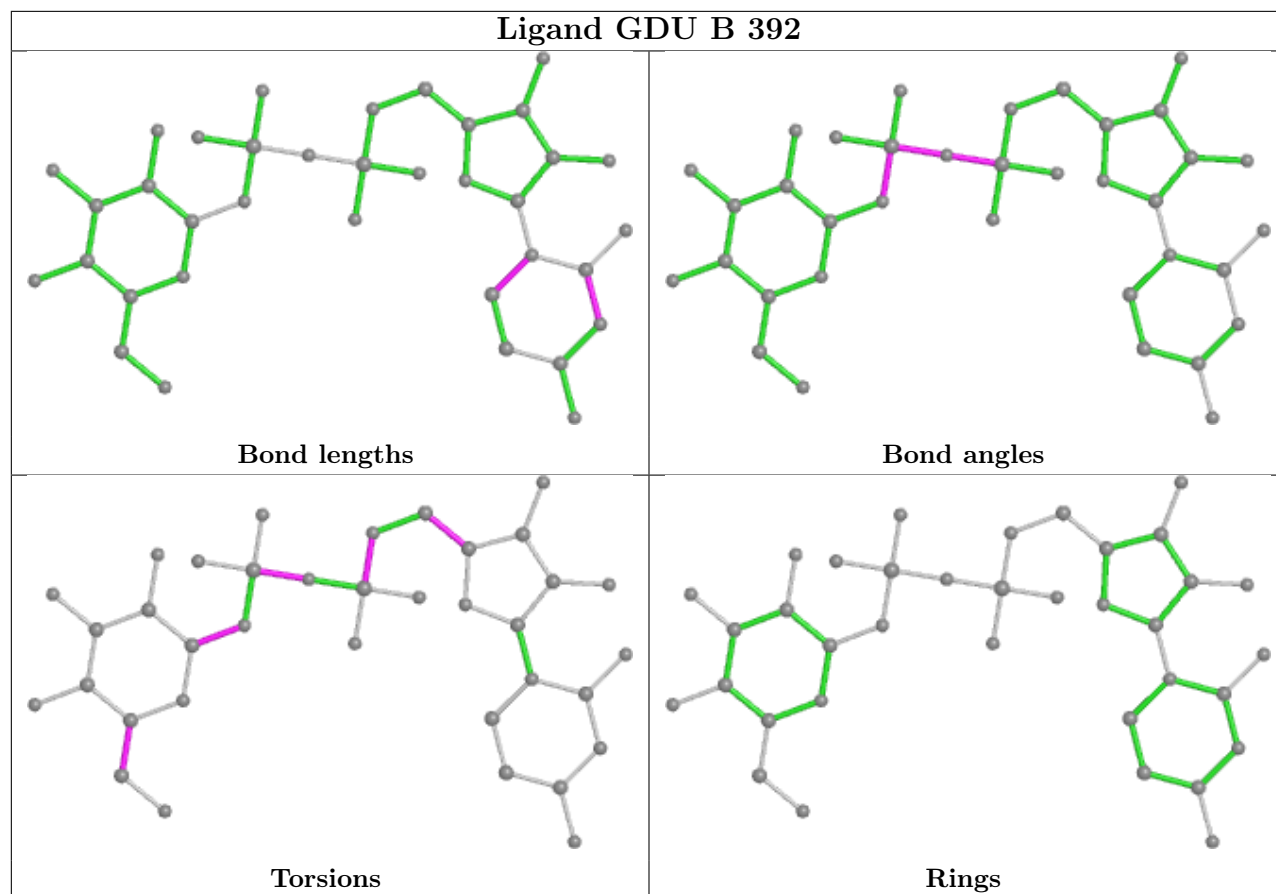
3 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	391	FDA	1	0
2	B	391	FDA	7	0
4	B	392	GDU	22	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

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	383/390 (98%)	-0.01	16 (4%) 36 39	24, 40, 68, 81	0
1	B	374/390 (95%)	0.06	14 (3%) 41 45	24, 42, 80, 93	0
All	All	757/780 (97%)	0.02	30 (3%) 38 41	24, 41, 73, 93	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	373	ASN	4.7
1	B	142	PHE	4.7
1	A	371	THR	4.2
1	A	2	LYS	3.9
1	A	374	GLN	3.7
1	B	170	SER	3.7
1	B	167	LEU	3.7
1	B	140	LEU	3.3
1	B	137	GLU	3.2
1	A	294	ASP	3.1
1	A	292	GLN	3.0
1	A	372	ASP	3.0
1	A	209	GLU	2.9
1	B	176	PRO	2.9
1	A	233	SER	2.9
1	A	14	SER	2.7
1	B	119	ALA	2.6
1	B	341	VAL	2.5
1	A	370	LEU	2.4
1	B	143	ILE	2.4
1	A	308	GLU	2.4
1	B	120	LEU	2.3
1	B	177	VAL	2.2
1	B	164	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	384	GLY	2.2
1	A	341	VAL	2.2
1	A	225	THR	2.1
1	A	3	SER	2.1
1	A	18	ILE	2.1
1	B	172	LEU	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](#)

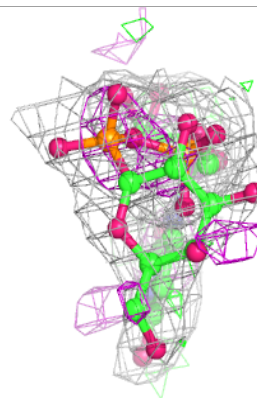
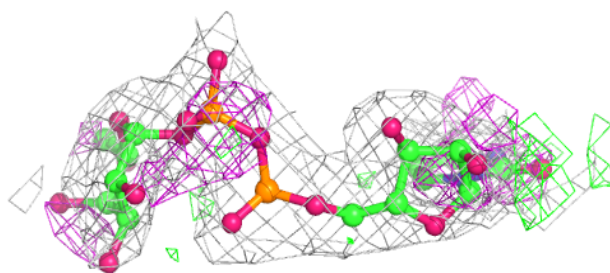
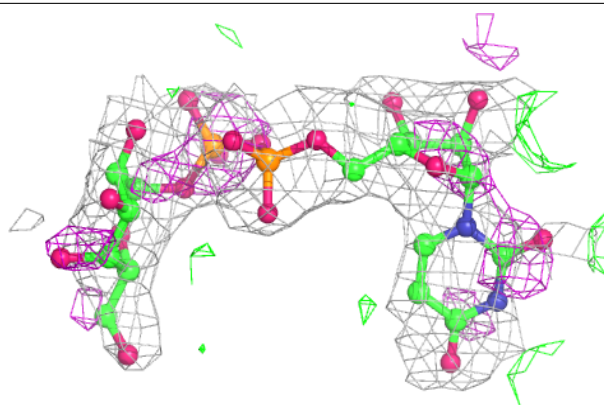
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	GDU	B	392	36/36	0.89	0.18	55,65,69,72	0
3	UDP	A	392	25/25	0.96	0.14	24,32,57,60	0
2	FDA	A	391	53/53	0.96	0.18	26,32,35,37	0
2	FDA	B	391	53/53	0.98	0.19	22,30,45,47	0

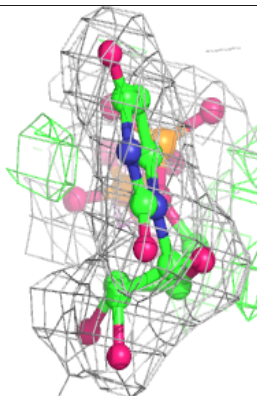
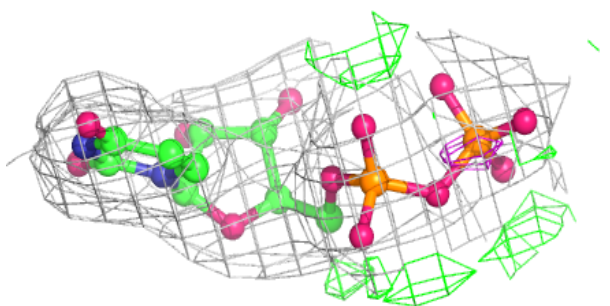
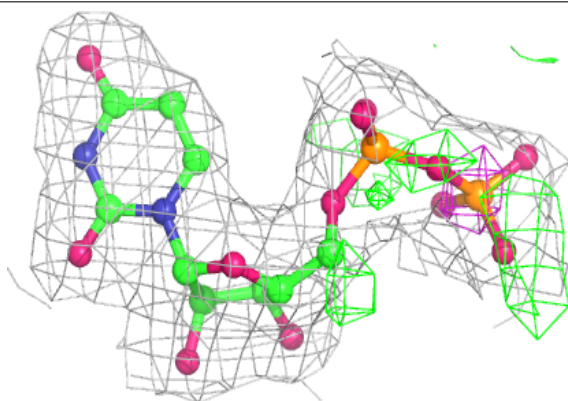
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GDU B 392:

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

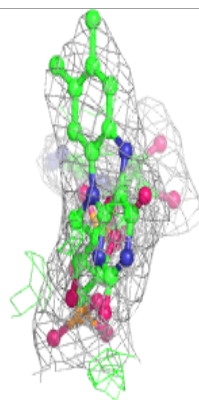
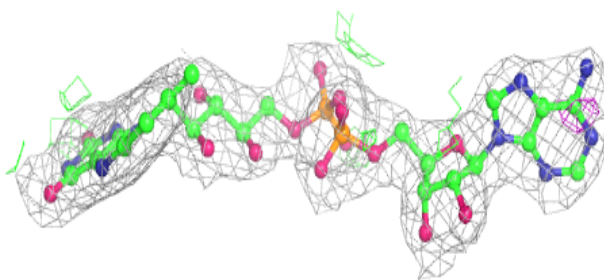
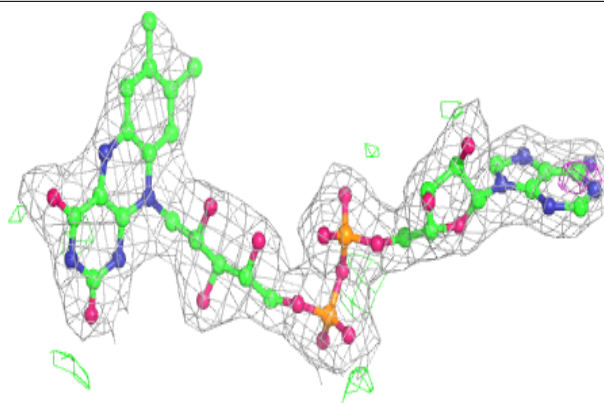
**Electron density around UDP A 392:**

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

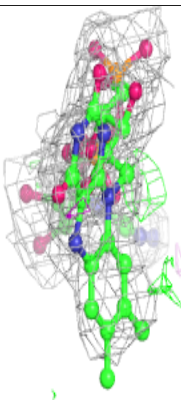
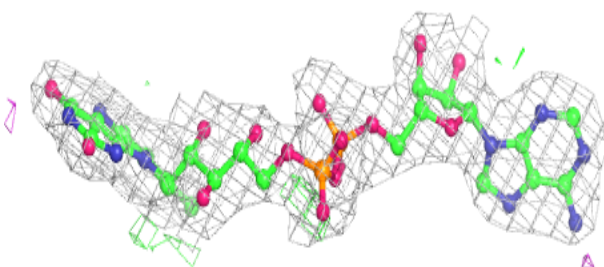
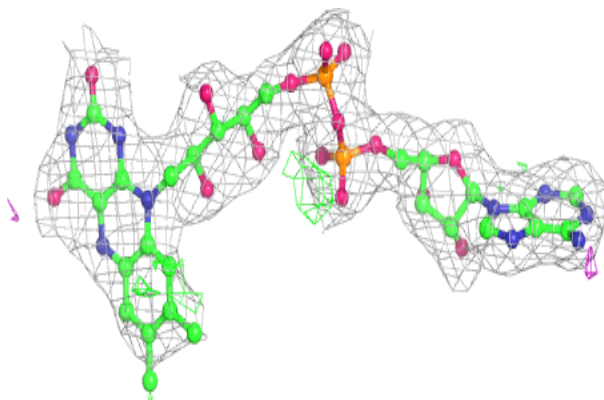


Electron density around FDA A 391:

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

**Electron density around FDA B 391:**

$2mF_o-DF_c$ (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.