



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 12:21 AM BST

PDB ID : 5U6M  
Title : Crystal structure of UDP-glucosyltransferase, UGT74F2, with UDP and salicylic acid  
Authors : George Thompson, A.M.; Iancu, C.V.; Dean, J.V.; Choe, J.  
Deposited on : 2016-12-08  
Resolution : 2.57 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](mailto: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 : 2.13.1  
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) : 2.13.1

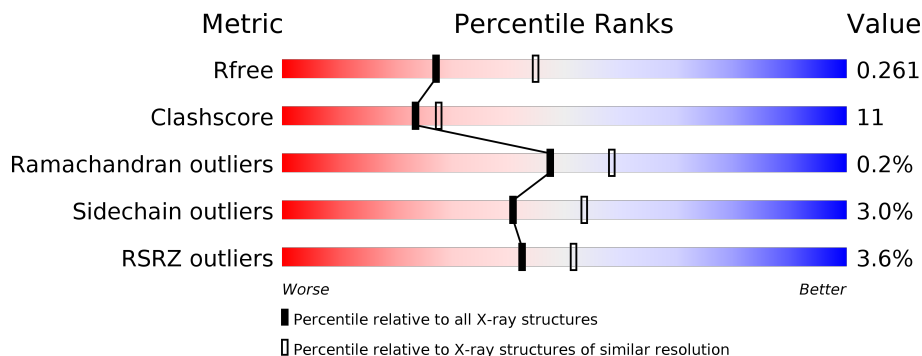
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.57 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	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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  $\geq 3$ , 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	449	 4% 72% 25% ..
1	B	449	 3% 77% 20% ..

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	BGC	B	501	X	-	-	-

## 2 Entry composition [i](#)

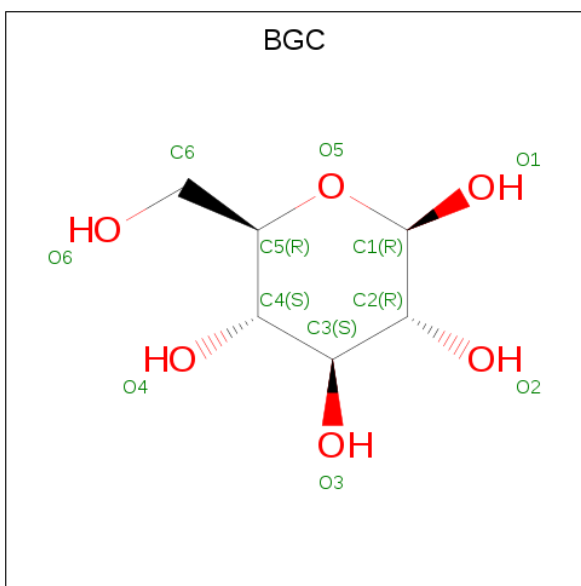
There are 5 unique types of molecules in this entry. The entry contains 7157 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-glycosyltransferase 74F2.

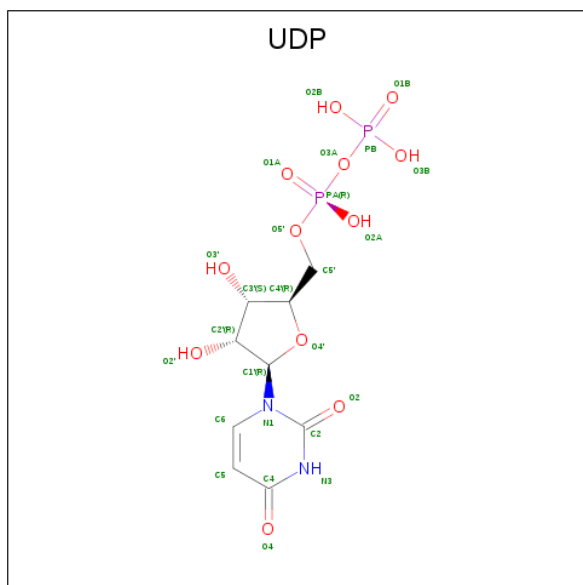
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	442	Total	C	N	O	S	0	0	0
			3517	2265	566	670	16			
1	B	445	Total	C	N	O	S	0	0	0
			3543	2280	573	674	16			

- Molecule 2 is beta-D-glucopyranose (three-letter code: BGC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



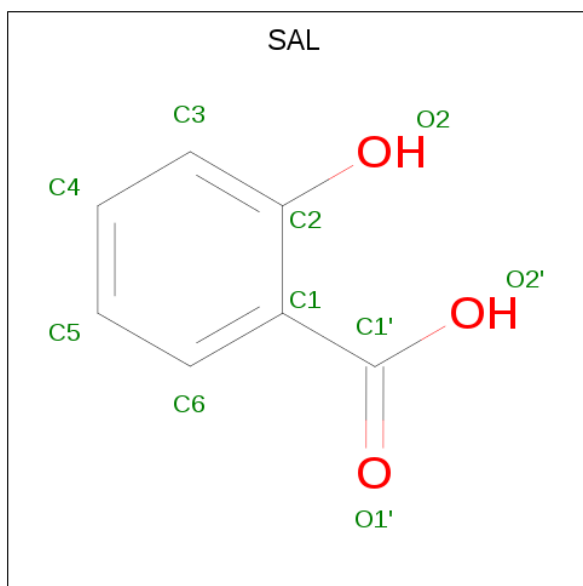
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	Total	C	O	0	0
			11	6	5		
2	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C<sub>9</sub>H<sub>14</sub>N<sub>2</sub>O<sub>12</sub>P<sub>2</sub>).



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

- Molecule 4 is 2-HYDROXYBENZOIC ACID (three-letter code: SAL) (formula: C<sub>7</sub>H<sub>6</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	10	7	3	0	0
4	B	1	10	7	3	0	0

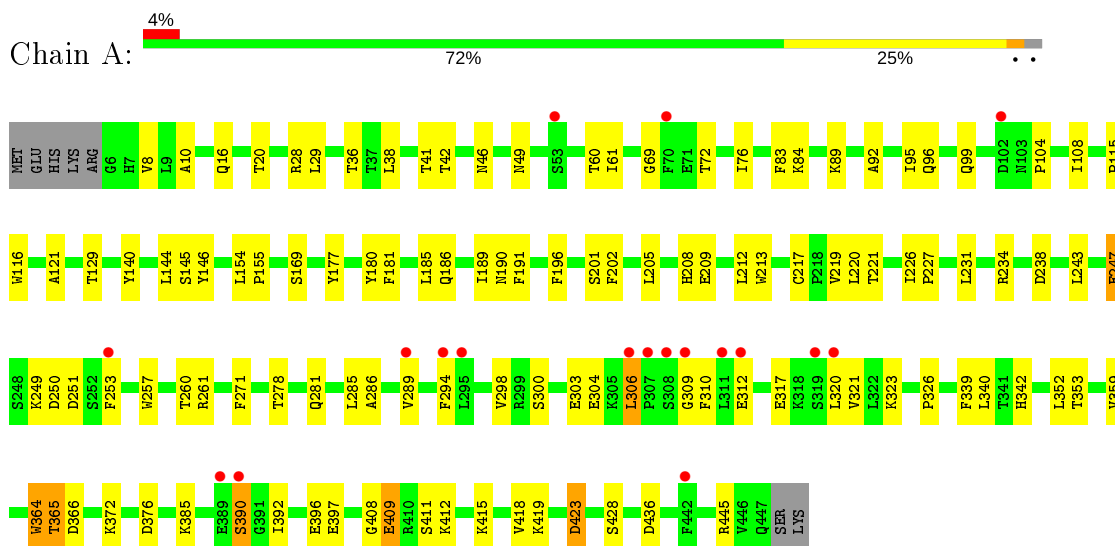
- Molecule 5 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
5	A	3	Total 3	O 3	0	0
5	B	2	Total 2	O 2	0	0

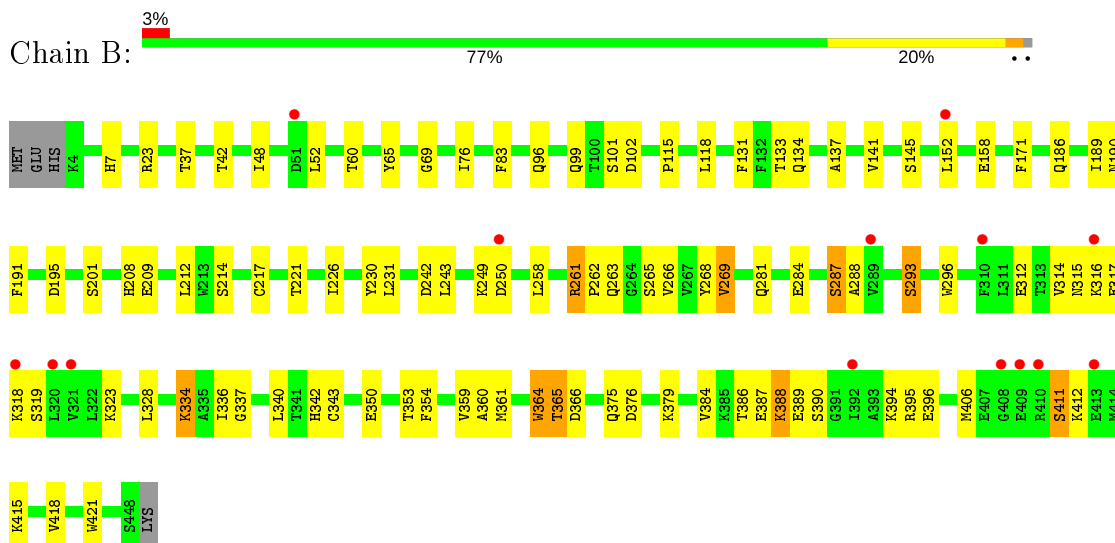
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: UDP-glycosyltransferase 74F2



- Molecule 1: UDP-glycosyltransferase 74F2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.69Å 87.68Å 164.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.84 – 2.57 43.84 – 2.57	Depositor EDS
% Data completeness (in resolution range)	98.7 (43.84-2.57) 98.7 (43.84-2.57)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.63 (at 2.58Å)	Xtrriage
Refinement program	PHENIX (dev_2481: ???)	Depositor
R, $R_{free}$	0.188 , 0.260 0.188 , 0.261	Depositor DCC
$R_{free}$ test set	1541 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	68.9	Xtrriage
Anisotropy	0.433	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 44.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7157	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	81.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: UDP, BGC, SAL

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	0.50	0/3606	0.65	1/4898 (0.0%)
1	B	0.54	0/3632	0.71	4/4931 (0.1%)
All	All	0.52	0/7238	0.68	5/9829 (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	3
1	B	0	1
All	All	0	4

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	261	ARG	NE-CZ-NH2	-8.55	116.03	120.30
1	B	334	LYS	CD-CE-NZ	8.18	130.51	111.70
1	B	261	ARG	NE-CZ-NH1	8.16	124.38	120.30
1	A	306	LEU	CB-CG-CD1	-5.62	101.45	111.00
1	B	269	VAL	CG1-CB-CG2	-5.12	102.71	110.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	247	GLU	Peptide

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Group
1	A	364	TRP	Peptide
1	A	409	GLU	Peptide
1	B	364	TRP	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	3517	0	3447	78	0
1	B	3543	0	3478	81	0
2	A	11	0	10	0	0
2	B	11	0	9	0	0
3	A	25	0	11	0	0
3	B	25	0	11	1	0
4	A	10	0	5	0	0
4	B	10	0	5	1	0
5	A	3	0	0	0	0
5	B	2	0	0	1	0
All	All	7157	0	6976	158	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 (158) 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:263:GLN:HG3	1:B:334:LYS:HE3	1.43	0.95
1:A:409:GLU:OE2	1:A:412:LYS:HG2	1.75	0.86
1:A:249:LYS:O	1:A:249:LYS:HD2	1.77	0.84
1:A:309:GLY:O	1:A:312:GLU:HG2	1.80	0.82
1:B:269:VAL:HG13	1:B:296:TRP:CD1	2.14	0.81
1:B:76:ILE:H	1:B:76:ILE:HD12	1.47	0.80
1:B:42:THR:HG22	1:B:60:THR:HB	1.61	0.79
1:A:298:VAL:HG11	1:A:306:LEU:HD11	1.65	0.79
1:A:46:ASN:HD21	1:B:60:THR:H	1.29	0.79
1:B:7:HIS:NE2	1:B:37:THR:HG23	1.99	0.78

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:GLU:OE1	1:B:395:ARG:N	2.18	0.76
1:B:186:GLN:O	1:B:189:ILE:HG13	1.86	0.75
1:A:408:GLY:O	1:A:411:SER:HB3	1.87	0.74
1:B:266:VAL:HG23	1:B:336:ILE:HD13	1.69	0.73
1:B:263:GLN:HG3	1:B:334:LYS:CE	2.18	0.72
1:A:249:LYS:NZ	1:A:253:PHE:HB3	2.05	0.71
1:A:209:GLU:HA	1:A:212:LEU:HD23	1.71	0.71
1:B:328:LEU:HB2	5:B:602:HOH:O	1.92	0.70
1:B:115:PRO:HB2	1:B:189:ILE:HD11	1.73	0.69
1:B:231:LEU:HD21	1:B:353:THR:HB	1.75	0.69
1:A:285:LEU:O	1:A:289:VAL:HG12	1.93	0.69
1:A:42:THR:HG22	1:A:60:THR:HB	1.77	0.65
1:B:365:THR:HG23	1:B:366:ASP:H	1.62	0.65
1:A:208:HIS:CE1	1:A:212:LEU:HD21	2.31	0.64
1:B:337:GLY:C	1:B:406:MET:HE1	2.17	0.64
1:A:249:LYS:HZ1	1:A:253:PHE:HB3	1.63	0.63
1:B:189:ILE:HD12	1:B:190:ASN:H	1.63	0.63
1:B:189:ILE:HD12	1:B:190:ASN:N	2.13	0.63
1:A:340:LEU:HD13	1:A:359:VAL:HG12	1.82	0.62
1:A:69:GLY:O	1:A:72:THR:HG22	2.00	0.62
1:B:284:GLU:OE1	1:B:394:LYS:HD2	1.99	0.61
1:B:214:SER:HA	1:B:217:CYS:O	2.01	0.60
1:A:365:THR:HG23	1:A:366:ASP:H	1.67	0.60
1:A:115:PRO:HB2	1:A:189:ILE:HD11	1.85	0.59
1:A:16:GLN:HE21	1:A:20:THR:CG2	2.16	0.58
1:B:266:VAL:CG2	1:B:336:ILE:HD13	2.34	0.58
1:B:201:SER:O	1:B:221:THR:HB	2.05	0.57
1:B:261:ARG:HG2	1:B:262:PRO:CD	2.35	0.57
1:B:284:GLU:CD	1:B:394:LYS:HD2	2.24	0.57
1:A:303:GLU:HA	1:A:306:LEU:CD1	2.36	0.56
1:B:249:LYS:HG3	1:B:249:LYS:O	2.05	0.56
1:A:217:CYS:SG	1:A:219:VAL:HG23	2.46	0.55
1:A:8:VAL:HB	1:A:36:THR:HG22	1.89	0.55
1:A:146:TYR:OH	1:A:189:ILE:HA	2.06	0.54
1:A:320:LEU:HD12	1:A:321:VAL:H	1.73	0.54
1:B:318:LYS:HG3	1:B:319:SER:OG	2.08	0.54
1:A:231:LEU:HD21	1:A:353:THR:CG2	2.38	0.54
1:B:76:ILE:CD1	1:B:76:ILE:H	2.18	0.54
1:B:145:SER:HG	1:B:171:PHE:HE1	1.56	0.53
1:B:261:ARG:HG3	1:B:261:ARG:HH11	1.73	0.53
1:A:317:GLU:N	1:A:317:GLU:OE1	2.42	0.53

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:287:SER:OG	1:B:395:ARG:HD3	2.09	0.52
1:A:212:LEU:H	1:A:212:LEU:HD22	1.74	0.52
1:A:409:GLU:HA	1:A:409:GLU:OE2	2.09	0.52
1:B:337:GLY:O	1:B:406:MET:HE1	2.09	0.52
1:B:284:GLU:CG	1:B:394:LYS:HA	2.40	0.51
1:B:261:ARG:HG3	1:B:261:ARG:NH1	2.25	0.51
1:B:411:SER:O	1:B:415:LYS:HE2	2.10	0.51
1:A:385:LYS:HB2	1:A:397:GLU:CG	2.41	0.51
1:B:268:TYR:HB2	1:B:336:ILE:HD12	1.92	0.51
1:B:314:VAL:HA	1:B:318:LYS:HZ2	1.76	0.51
1:A:300:SER:OG	1:A:323:LYS:HG2	2.12	0.50
1:A:76:ILE:HD13	1:A:180:TYR:CE2	2.46	0.50
1:A:186:GLN:O	1:A:189:ILE:HG13	2.12	0.50
1:B:340:LEU:HD13	1:B:359:VAL:HG13	1.93	0.50
1:B:52:LEU:HD22	1:B:52:LEU:N	2.27	0.50
1:B:48:ILE:HD13	1:B:243:LEU:HD22	1.94	0.50
1:B:315:ASN:OD1	1:B:318:LYS:HB3	2.12	0.49
1:A:231:LEU:HD21	1:A:353:THR:HG23	1.93	0.49
1:A:289:VAL:HG22	1:A:294:PHE:CD2	2.47	0.49
1:B:266:VAL:HG12	1:B:293:SER:OG	2.11	0.49
1:B:343:CYS:SG	1:B:360:ALA:HB1	2.52	0.49
1:B:261:ARG:HD2	1:B:266:VAL:HG12	1.94	0.49
1:A:177:TYR:HB3	1:A:180:TYR:HB2	1.95	0.49
1:A:251:ASP:HB2	1:A:326:PRO:HG3	1.95	0.49
1:A:385:LYS:HB2	1:A:397:GLU:HG2	1.94	0.49
1:B:288:ALA:HB2	1:B:395:ARG:HB2	1.95	0.49
1:B:412:LYS:HD3	1:B:415:LYS:HE3	1.95	0.48
1:B:262:PRO:HD2	1:B:265:SER:OG	2.13	0.48
1:A:10:ALA:O	1:A:38:LEU:HD12	2.14	0.48
1:B:249:LYS:NZ	1:B:250:ASP:HB2	2.27	0.48
1:B:375:GLN:HG3	1:B:376:ASP:OD1	2.12	0.48
1:A:340:LEU:HD13	1:A:359:VAL:CG1	2.44	0.48
1:B:418:VAL:HA	1:B:421:TRP:HD1	1.79	0.48
1:A:226:ILE:HG22	1:A:227:PRO:HD2	1.96	0.47
1:B:158:GLU:N	1:B:158:GLU:OE1	2.46	0.47
1:A:390:SER:OG	1:A:392:ILE:HG12	2.14	0.47
1:B:281:GLN:O	1:B:281:GLN:NE2	2.45	0.47
1:B:318:LYS:HG2	1:B:319:SER:N	2.29	0.47
1:A:189:ILE:HD12	1:A:190:ASN:N	2.30	0.46
1:A:108:ILE:HD12	1:A:121:ALA:HB2	1.97	0.46
1:A:320:LEU:HD12	1:A:321:VAL:N	2.30	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:PRO:CB	1:B:189:ILE:HD11	2.42	0.46
1:A:41:THR:HA	1:A:61:ILE:O	2.15	0.46
1:A:95:ILE:HD11	1:A:108:ILE:HD11	1.98	0.46
1:A:298:VAL:HG21	1:A:306:LEU:HD21	1.97	0.46
1:B:261:ARG:HG2	1:B:262:PRO:HD2	1.97	0.46
1:B:65:TYR:O	1:B:69:GLY:N	2.42	0.46
1:B:396:GLU:OE2	1:B:396:GLU:N	2.47	0.45
1:A:16:GLN:O	1:A:20:THR:HG23	2.17	0.45
1:A:181:PHE:CZ	1:A:185:LEU:HD11	2.51	0.45
1:A:304:GLU:H	1:A:304:GLU:CD	2.16	0.45
1:A:249:LYS:HG3	1:A:251:ASP:HA	1.99	0.45
1:A:201:SER:O	1:A:221:THR:HB	2.17	0.45
1:A:271:PHE:CD2	1:A:342:HIS:HB3	2.52	0.45
1:B:137:ALA:HB2	1:B:209:GLU:HG2	1.99	0.45
1:B:209:GLU:H	1:B:209:GLU:CD	2.20	0.44
1:B:386:THR:OG1	1:B:387:GLU:N	2.49	0.44
1:A:49:ASN:HD22	1:A:247:GLU:CD	2.20	0.44
1:B:131:PHE:CE1	1:B:133:THR:HG22	2.53	0.44
1:A:129:THR:HG22	1:A:196:PHE:H	1.83	0.44
1:A:243:LEU:HD23	1:A:243:LEU:HA	1.82	0.44
1:B:141:VAL:HG13	1:B:152:LEU:HD11	2.00	0.43
1:B:261:ARG:CG	1:B:262:PRO:HD3	2.48	0.43
1:B:316:LYS:HE3	1:B:316:LYS:HB3	1.56	0.43
1:A:140:TYR:CE2	1:A:144:LEU:HD11	2.54	0.43
1:A:202:PHE:CE1	1:A:205:LEU:HD22	2.53	0.43
1:A:169:SER:HB3	1:A:364:TRP:HA	2.01	0.43
1:A:99:GLN:HA	1:A:104:PRO:HB3	2.00	0.43
1:A:253:PHE:HE2	1:A:320:LEU:HD21	1.84	0.43
1:B:23:ARG:NE	1:B:242:ASP:OD1	2.52	0.43
1:B:243:LEU:HD23	1:B:243:LEU:HA	1.83	0.43
1:A:16:GLN:HE21	1:A:20:THR:HG22	1.84	0.43
1:B:118:LEU:HD23	1:B:190:ASN:CG	2.40	0.42
1:B:312:GLU:OE1	1:B:312:GLU:N	2.52	0.42
1:A:227:PRO:HA	1:A:238:ASP:OD2	2.19	0.42
1:B:226:ILE:HD11	1:B:350:GLU:HA	2.01	0.42
1:B:364:TRP:CG	1:B:365:THR:N	2.87	0.42
1:A:352:LEU:HD23	1:A:418:VAL:HG23	2.01	0.42
1:B:249:LYS:HG3	1:B:249:LYS:HZ2	1.68	0.42
1:A:415:LYS:O	1:A:419:LYS:HG3	2.19	0.42
1:A:145:SER:HB2	1:A:185:LEU:HD21	2.02	0.42
1:A:29:LEU:HD23	1:A:29:LEU:HA	1.84	0.42

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ALA:O	1:A:96:GLN:HG3	2.20	0.42
1:B:388:LYS:HD2	1:B:388:LYS:H	1.84	0.42
1:A:423:ASP:N	1:A:423:ASP:OD1	2.52	0.42
1:B:96:GLN:O	1:B:99:GLN:HB2	2.20	0.42
1:A:257:TRP:O	1:A:260:THR:HB	2.20	0.41
1:B:342:HIS:HE1	3:B:502:UDP:O2B	2.02	0.41
1:A:396:GLU:CD	1:A:396:GLU:H	2.23	0.41
1:A:84:LYS:HE3	1:A:116:TRP:HB3	2.01	0.41
1:A:281:GLN:HG2	1:A:392:ILE:HD12	2.02	0.41
1:B:261:ARG:CZ	1:B:317:GLU:O	2.68	0.41
1:A:372:LYS:HE3	1:A:376:ASP:HB2	2.03	0.41
1:B:230:TYR:HB3	1:B:354:PHE:HE1	1.85	0.41
1:A:278:THR:OG1	1:A:281:GLN:HG3	2.21	0.41
1:A:286:ALA:HB1	1:A:310:PHE:CE1	2.55	0.41
1:B:208:HIS:NE2	1:B:212:LEU:HD11	2.35	0.41
1:B:134:GLN:HE22	4:B:503:SAL:H5	1.86	0.41
1:A:213:TRP:CE3	1:A:219:VAL:HG21	2.56	0.41
1:A:196:PHE:HB3	1:A:220:LEU:HD13	2.03	0.40
1:B:261:ARG:CG	1:B:262:PRO:CD	2.99	0.40
1:A:154:LEU:HD22	1:A:155:PRO:HA	2.02	0.40
1:A:234:ARG:HH12	1:A:436:ASP:CG	2.24	0.40
1:B:258:LEU:HD22	1:B:266:VAL:HG21	2.03	0.40
1:B:269:VAL:CG1	1:B:296:TRP:CD1	2.95	0.40
1:B:375:GLN:O	1:B:379:LYS:HA	2.21	0.40
1:B:360:ALA:O	1:B:384:VAL:HG23	2.22	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	440/449 (98%)	421 (96%)	18 (4%)	1 (0%)	47 58

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	443/449 (99%)	417 (94%)	25 (6%)	1 (0%)	47	58
All	All	883/898 (98%)	838 (95%)	43 (5%)	2 (0%)	47	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	365	THR
1	B	365	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	394/401 (98%)	383 (97%)	11 (3%)	43	56
1	B	397/401 (99%)	384 (97%)	13 (3%)	38	50
All	All	791/802 (99%)	767 (97%)	24 (3%)	41	53

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ARG
1	A	83	PHE
1	A	89	LYS
1	A	191	PHE
1	A	250	ASP
1	A	261	ARG
1	A	339	PHE
1	A	390	SER
1	A	423	ASP
1	A	428	SER
1	A	445	ARG
1	B	83	PHE
1	B	101	SER
1	B	102	ASP

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	191	PHE
1	B	195	ASP
1	B	287	SER
1	B	293	SER
1	B	323	LYS
1	B	361	MET
1	B	388	LYS
1	B	389	GLU
1	B	390	SER
1	B	411	SER

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

Mol	Chain	Res	Type
1	A	46	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

6 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	BGC	B	501	1	11,11,12	1.06	0	15,15,17	2.58	6 (40%)
2	BGC	A	501	1	11,11,12	1.40	3 (27%)	15,15,17	1.18	1 (6%)
3	UDP	A	502	-	20,26,26	3.74	10 (50%)	25,40,40	1.20	3 (12%)
4	SAL	B	503	-	8,10,10	4.28	4 (50%)	9,13,13	1.36	1 (11%)
4	SAL	A	503	-	8,10,10	4.32	3 (37%)	9,13,13	1.10	1 (11%)
3	UDP	B	502	-	20,26,26	3.80	9 (45%)	25,40,40	1.28	2 (8%)

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	BGC	B	501	1	1/1/4/5	2/2/19/22	0/1/1/1
2	BGC	A	501	1	-	2/2/19/22	0/1/1/1
3	UDP	A	502	-	-	3/14/32/32	0/2/2/2
4	SAL	B	503	-	-	0/0/4/4	0/1/1/1
4	SAL	A	503	-	-	0/0/4/4	0/1/1/1
3	UDP	B	502	-	-	3/14/32/32	0/2/2/2

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	503	SAL	C1-C1'	10.96	1.58	1.47
4	B	503	SAL	C1-C1'	10.81	1.57	1.47
3	B	502	UDP	C6-N1	8.22	1.46	1.35
3	A	502	UDP	C6-N1	7.16	1.44	1.35
3	B	502	UDP	C4-N3	7.12	1.45	1.33
3	A	502	UDP	C4-N3	6.85	1.45	1.33
3	A	502	UDP	O4'-C1'	-6.57	1.31	1.41
3	A	502	UDP	C3'-C4'	-6.12	1.37	1.53
3	B	502	UDP	O4'-C1'	-6.02	1.32	1.41
3	B	502	UDP	C3'-C4'	-5.90	1.37	1.53
3	B	502	UDP	O4'-C4'	5.78	1.57	1.45
3	A	502	UDP	O4'-C4'	5.75	1.57	1.45
3	B	502	UDP	C6-C5	5.37	1.49	1.38
3	A	502	UDP	C6-C5	4.89	1.48	1.38
3	A	502	UDP	C2-N3	4.13	1.46	1.38
3	B	502	UDP	C2-N3	3.98	1.46	1.38
4	A	503	SAL	C6-C1	-3.90	1.33	1.40

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	503	SAL	C6-C1	-3.57	1.34	1.40
4	A	503	SAL	C5-C6	2.53	1.44	1.38
2	A	501	BGC	O5-C1	-2.51	1.39	1.43
3	A	502	UDP	O2'-C2'	-2.48	1.37	1.43
4	B	503	SAL	O2-C2	2.46	1.41	1.36
4	B	503	SAL	C5-C6	2.43	1.44	1.38
3	A	502	UDP	O3'-C3'	2.37	1.48	1.43
3	B	502	UDP	O3'-C3'	2.33	1.48	1.43
3	A	502	UDP	O4-C4	-2.12	1.19	1.24
3	B	502	UDP	O4-C4	-2.08	1.19	1.24
2	A	501	BGC	C6-C5	2.07	1.58	1.51
2	A	501	BGC	C1-C2	-2.02	1.47	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	BGC	C1-C2-C3	7.49	118.87	109.67
3	B	502	UDP	C2'-C3'-C4'	3.45	109.34	102.64
2	B	501	BGC	O5-C5-C6	-3.40	101.87	107.20
4	B	503	SAL	C6-C1-C1'	-3.39	115.16	120.20
2	A	501	BGC	C1-C2-C3	3.33	113.76	109.67
3	A	502	UDP	O3B-PB-O3A	-3.00	94.56	104.64
4	A	503	SAL	C6-C1-C1'	-2.74	116.12	120.20
3	A	502	UDP	C2'-C3'-C4'	2.64	107.77	102.64
2	B	501	BGC	O2-C2-C3	-2.61	104.91	110.14
2	B	501	BGC	O5-C1-C2	2.48	114.59	110.77
3	B	502	UDP	O2B-PB-O3A	2.33	112.45	104.64
2	B	501	BGC	C1-O5-C5	2.20	115.18	112.19
3	A	502	UDP	O3B-PB-O2B	2.10	115.68	107.64
2	B	501	BGC	O3-C3-C4	-2.04	105.64	110.35

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	501	BGC	C1

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	UDP	O4'-C4'-C5'-O5'
2	B	501	BGC	O5-C5-C6-O6
2	A	501	BGC	O5-C5-C6-O6

*Continued on next page...*

*Continued from previous page...*

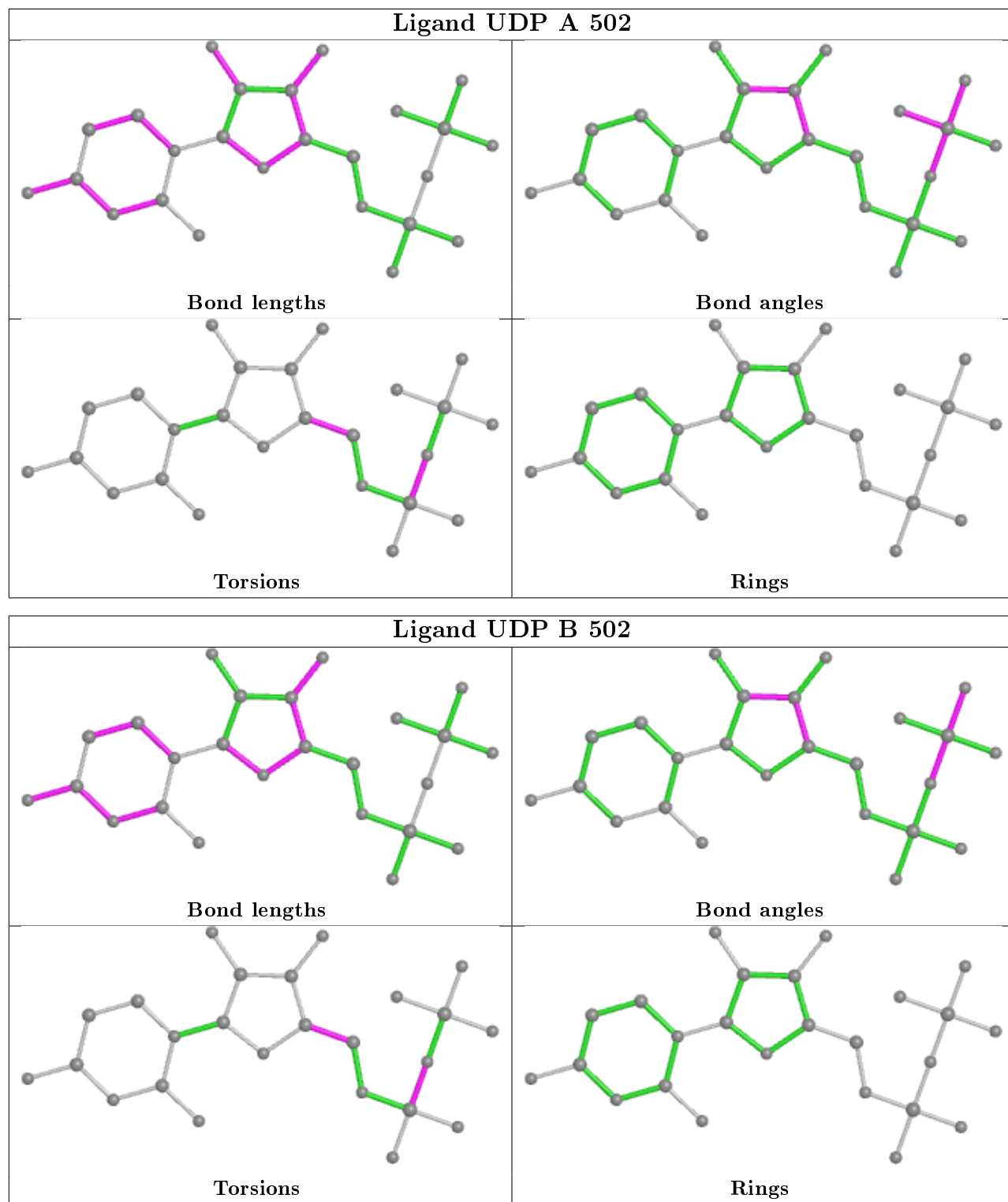
Mol	Chain	Res	Type	Atoms
2	B	501	BGC	C4-C5-C6-O6
3	A	502	UDP	C3'-C4'-C5'-O5'
2	A	501	BGC	C4-C5-C6-O6
3	B	502	UDP	C3'-C4'-C5'-O5'
3	B	502	UDP	O4'-C4'-C5'-O5'
3	A	502	UDP	PB-O3A-PA-O5'
3	B	502	UDP	PB-O3A-PA-O5'

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	503	SAL	1	0
3	B	502	UDP	1	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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	442/449 (98%)	0.05	18 (4%) 37 46	49, 79, 125, 165	0
1	B	445/449 (99%)	0.02	14 (3%) 49 58	44, 74, 124, 177	0
All	All	887/898 (98%)	0.04	32 (3%) 42 51	44, 77, 126, 177	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	53	SER	6.5
1	B	408	GLY	5.8
1	B	410	ARG	4.6
1	B	321	VAL	4.2
1	B	409	GLU	4.0
1	A	320	LEU	3.9
1	A	309	GLY	3.5
1	A	306	LEU	3.4
1	B	289	VAL	3.3
1	A	308	SER	2.9
1	A	389	GLU	2.9
1	A	253	PHE	2.8
1	A	70	PHE	2.7
1	B	310	PHE	2.7
1	A	294	PHE	2.7
1	A	295	LEU	2.6
1	B	413	GLU	2.6
1	B	392	ILE	2.5
1	A	319	SER	2.4
1	A	390	SER	2.4
1	A	311	LEU	2.3
1	A	289	VAL	2.2
1	B	51	ASP	2.1
1	B	250	ASP	2.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	442	PHE	2.1
1	A	307	PRO	2.1
1	A	312	GLU	2.1
1	B	318	LYS	2.1
1	B	320	LEU	2.1
1	B	316	LYS	2.1
1	A	102	ASP	2.0
1	B	152	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](#)

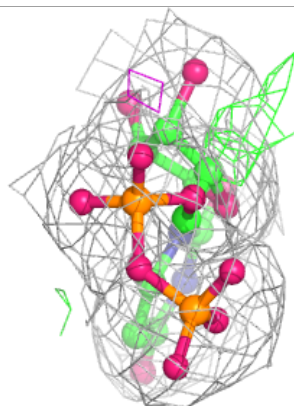
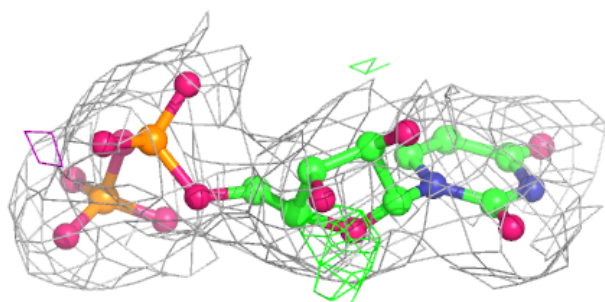
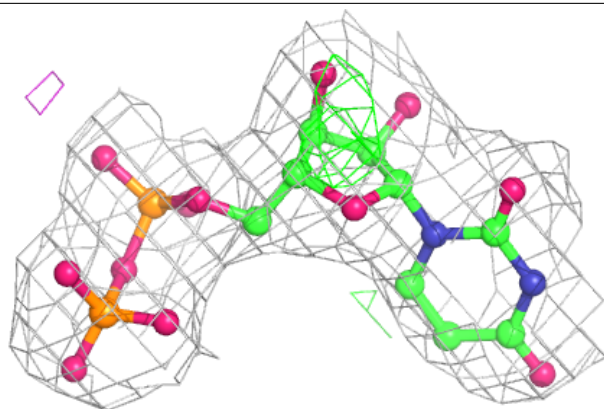
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	SAL	B	503	10/10	0.73	0.36	101,116,120,121	0
4	SAL	A	503	10/10	0.88	0.27	83,90,101,101	0
2	BGC	B	501	11/12	0.90	0.15	75,91,102,102	0
2	BGC	A	501	11/12	0.96	0.12	92,96,97,98	0
3	UDP	A	502	25/25	0.97	0.15	47,57,68,75	0
3	UDP	B	502	25/25	0.97	0.18	50,68,73,74	0

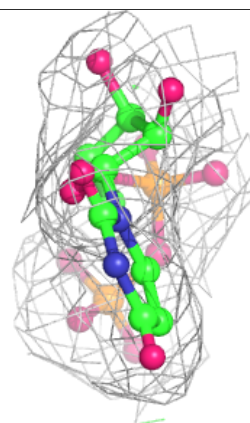
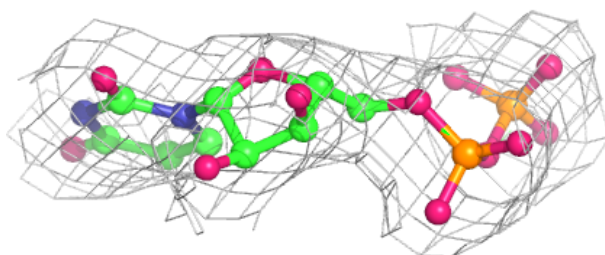
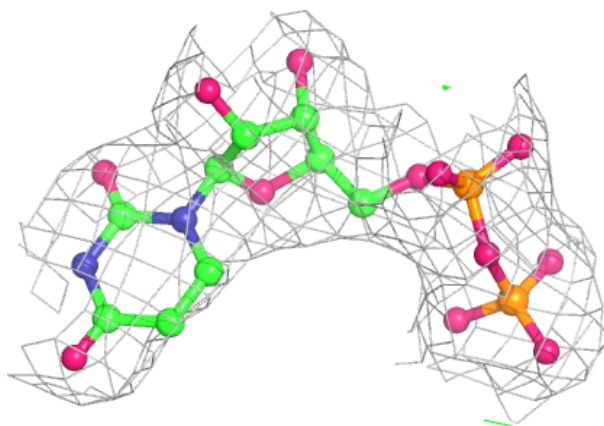
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around UDP A 502:**

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

**Electron density around UDP B 502:**

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





## 6.5 Other polymers

There are no such residues in this entry.