



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 7, 2020 – 02:50 am BST

PDB ID : 6IJ9  
Title : Crystal Structure of Arabidopsis thaliana UGT89C1 complexed with UDP  
Authors : Zong, G.N.; Wang, X.Q.  
Deposited on : 2018-10-09  
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](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 : 2.11  
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

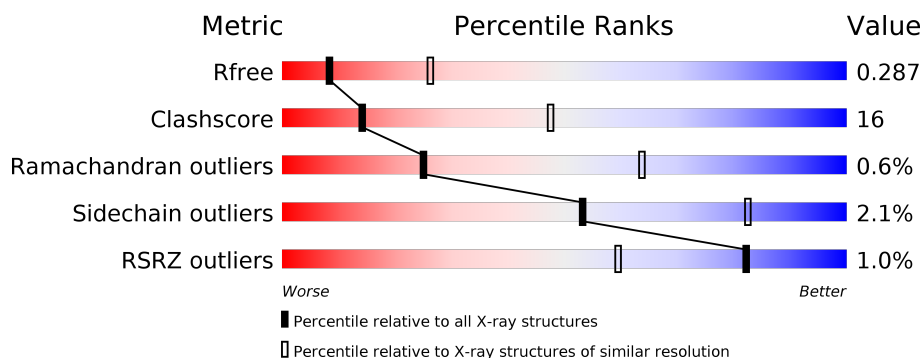
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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 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	435	<div> <div> <div>2%</div> <div>66%</div> <div>27%</div> <div>• 6%</div> </div> </div>
1	B	435	<div> <div>63%</div> <div>30%</div> <div>• 6%</div> </div>
1	C	435	<div> <div> <div>2%</div> <div>65%</div> <div>28%</div> <div>• 5%</div> </div> </div>
1	D	435	<div> <div> <div>2%</div> <div>52%</div> <div>34%</div> <div>• 11%</div> </div> </div>

## 2 Entry composition [i](#)

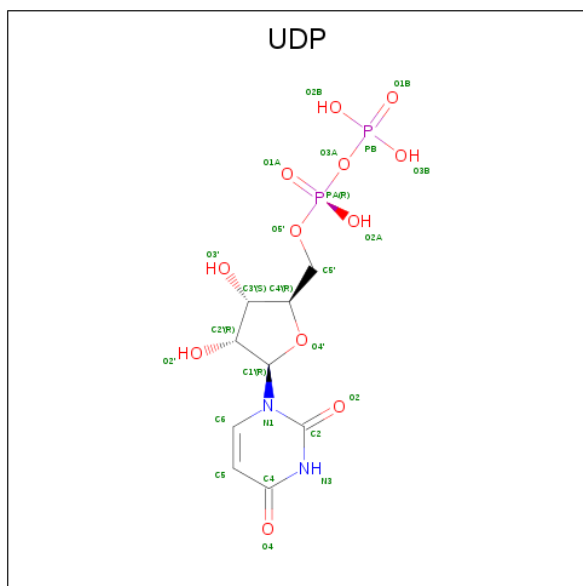
There are 3 unique types of molecules in this entry. The entry contains 12804 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 89C1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	410	Total	C	N	O	S	0	0	0
			3197	2051	549	584	13			
1	B	408	Total	C	N	O	S	0	0	0
			3174	2038	544	580	12			
1	C	413	Total	C	N	O	S	0	0	0
			3232	2074	554	591	13			
1	D	388	Total	C	N	O	S	0	0	0
			3040	1959	518	551	12			

- 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
2	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

*Continued on next page...*

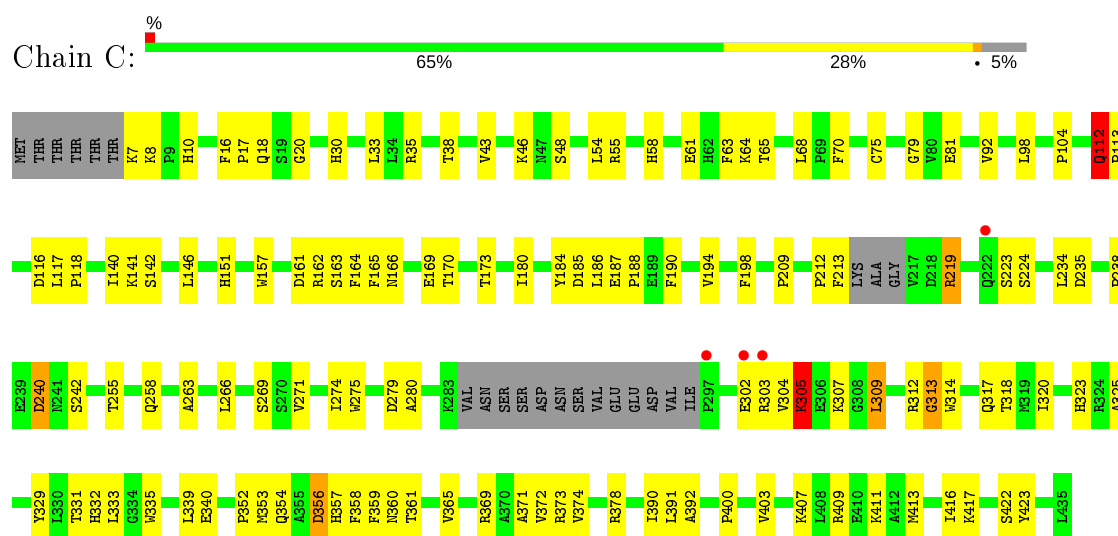
*Continued from previous page...*

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

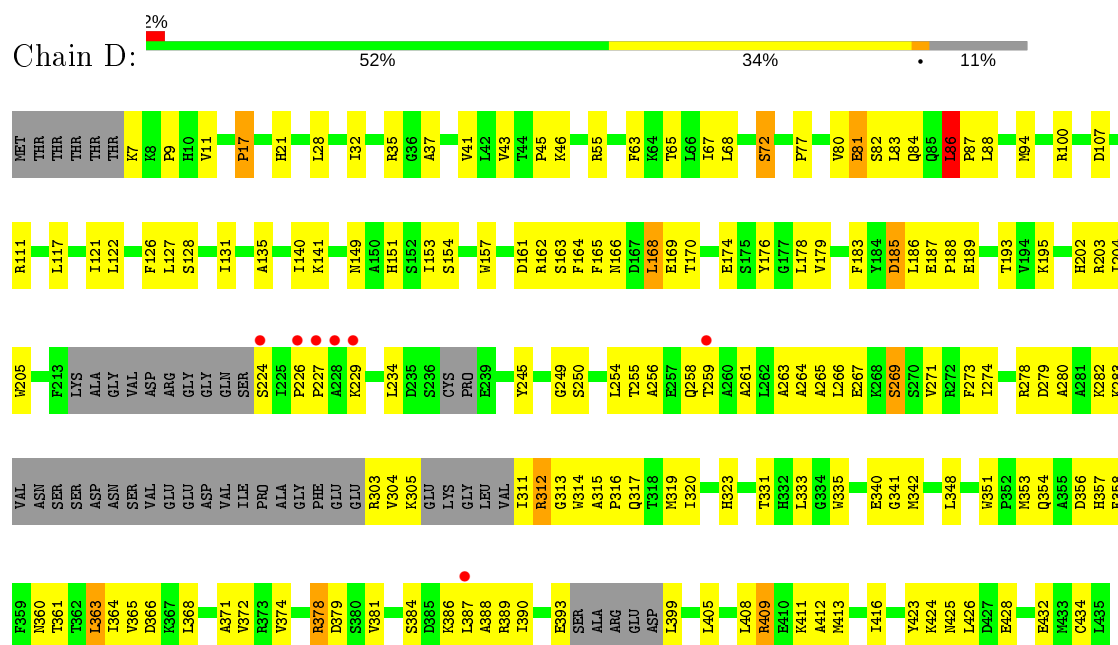
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	14	Total	O	0	0
			14	14		
3	B	18	Total	O	0	0
			18	18		
3	C	14	Total	O	0	0
			14	14		
3	D	15	Total	O	0	0
			15	15		





• Molecule 1: UDP-glycosyltransferase 89C1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.98Å 84.14Å 129.15Å 90.00° 108.90° 90.00°	Depositor
Resolution (Å)	29.93 – 3.00 29.93 – 3.00	Depositor EDS
% Data completeness (in resolution range)	83.1 (29.93-3.00) 83.2 (29.93-3.00)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.86 (at 3.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.217 , 0.274 0.228 , 0.287	Depositor DCC
$R_{free}$ test set	1575 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.4	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 41.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	12804	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.04 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.5595e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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

### 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	0.60	0/3277	0.77	3/4458 (0.1%)
1	B	0.55	1/3253 (0.0%)	0.84	5/4426 (0.1%)
1	C	0.63	1/3313 (0.0%)	0.84	7/4504 (0.2%)
1	D	0.56	0/3115	0.90	8/4235 (0.2%)
All	All	0.59	2/12958 (0.0%)	0.84	23/17623 (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	3
1	C	0	2
1	D	0	2
All	All	0	10

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	398	ASP	C-N	7.38	1.51	1.34
1	C	356	ASP	CB-CG	-7.17	1.36	1.51

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	425	ASN	CB-CG-OD1	20.45	162.51	121.60
1	B	85	GLN	CG-CD-OE1	18.57	158.73	121.60
1	D	425	ASN	CB-CG-ND2	-17.50	74.70	116.70

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	85	GLN	CG-CD-NE2	-17.40	74.94	116.70
1	C	112	GLN	CG-CD-OE1	14.76	151.11	121.60
1	C	112	GLN	CG-CD-NE2	-14.50	81.89	116.70
1	D	425	ASN	OD1-CG-ND2	-11.25	96.03	121.90
1	C	112	GLN	OE1-CD-NE2	-10.09	98.69	121.90
1	B	85	GLN	OE1-CD-NE2	-10.01	98.87	121.90
1	C	112	GLN	CA-CB-CG	-9.37	92.78	113.40
1	D	86	LEU	CA-CB-CG	9.16	136.37	115.30
1	B	168	LEU	CA-CB-CG	8.41	134.65	115.30
1	C	112	GLN	CB-CG-CD	8.08	132.61	111.60
1	A	360	ASN	N-CA-CB	7.97	124.95	110.60
1	A	359	PHE	C-N-CA	-7.93	101.88	121.70
1	C	356	ASP	CB-CA-C	-6.66	97.08	110.40
1	D	425	ASN	N-CA-CB	6.46	122.22	110.60
1	A	360	ASN	CB-CG-OD1	-6.44	108.72	121.60
1	D	168	LEU	CA-CB-CG	6.16	129.48	115.30
1	B	168	LEU	CB-CG-CD2	-5.99	100.82	111.00
1	C	313	GLY	N-CA-C	5.97	128.01	113.10
1	D	363	LEU	CA-CB-CG	5.82	128.69	115.30
1	D	424	LYS	C-N-CA	-5.46	108.04	121.70

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	303	ARG	Peptide
1	A	306	GLU	Peptide
1	A	360	ASN	Sidechain
1	B	307	LYS	Peptide
1	B	308	GLY	Peptide
1	B	85	GLN	Sidechain
1	C	112	GLN	Sidechain
1	C	309	LEU	Peptide
1	D	185	ASP	Peptide
1	D	81	GLU	Peptide

## 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	3197	0	3195	86	0
1	B	3174	0	3164	93	0
1	C	3232	0	3244	104	0
1	D	3040	0	3051	130	0
2	A	25	0	11	3	0
2	B	25	0	11	1	0
2	C	25	0	11	3	0
2	D	25	0	11	1	0
3	A	14	0	0	0	0
3	B	18	0	0	1	0
3	C	14	0	0	1	0
3	D	15	0	0	2	0
All	All	12804	0	12698	396	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 16.

All (396) 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:C:112:GLN:HA	1:C:112:GLN:HE21	1.20	1.01
1:D:224:SER:N	3:D:1001:HOH:O	1.93	1.01
1:C:112:GLN:HG3	1:C:113:PRO:HD2	1.43	0.96
1:D:312:ARG:NH1	1:D:315:ALA:HA	1.88	0.89
1:B:166:ASN:ND2	1:D:379:ASP:OD2	2.08	0.87
1:B:379:ASP:OD2	1:D:166:ASN:ND2	2.09	0.85
1:C:219:ARG:HH11	1:C:318:THR:HG21	1.40	0.85
1:D:179:VAL:HA	1:D:205:TRP:HB2	1.59	0.84
1:C:112:GLN:CA	1:C:112:GLN:HE21	1.91	0.82
1:D:282:LYS:HG2	1:D:283:LYS:H	1.43	0.80
1:D:17:PRO:O	1:D:82:SER:HB2	1.81	0.80
1:A:186:LEU:HD21	1:A:363:LEU:HD21	1.63	0.80
1:D:186:LEU:HB3	1:D:187:GLU:HG2	1.64	0.79
1:A:379:ASP:OD2	1:C:166:ASN:ND2	2.14	0.78
1:D:274:ILE:HD11	1:D:320:ILE:HD13	1.64	0.77
1:C:219:ARG:NH1	1:C:340:GLU:OE2	2.18	0.77
1:A:80:VAL:HG11	1:A:86:LEU:HD21	1.68	0.76
1:B:38:THR:HG23	1:B:62:HIS:HA	1.68	0.75
1:D:84:GLN:NE2	3:D:1002:HOH:O	2.19	0.75
1:D:312:ARG:HH12	1:D:314:TRP:C	1.89	0.75

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:GLU:HB3	1:A:302:GLU:HB2	1.68	0.74
1:C:17:PRO:HG2	1:C:98:LEU:HD11	1.70	0.73
1:D:46:LYS:HB2	1:D:81:GLU:HG2	1.70	0.73
1:A:353:MET:HG2	1:A:378:ARG:HA	1.71	0.73
1:B:373:ARG:HH21	1:B:376:GLU:HA	1.54	0.73
1:D:269:SER:HB2	1:D:388:ALA:HB1	1.71	0.72
1:A:254:LEU:HB3	1:A:258:GLN:HG3	1.72	0.72
1:B:244:VAL:HG22	1:B:328:SER:HB2	1.70	0.71
1:D:361:THR:O	1:D:365:VAL:HG22	1.90	0.71
1:D:226:PRO:HB2	1:D:229:LYS:HD3	1.72	0.70
1:D:255:THR:HG23	1:D:258:GLN:H	1.55	0.70
1:C:112:GLN:HG3	1:C:113:PRO:CD	2.18	0.70
1:B:369:ARG:HH12	1:C:369:ARG:HD2	1.56	0.70
1:C:331:THR:HG22	1:C:333:LEU:H	1.56	0.70
1:D:203:ARG:NH2	1:D:428:GLU:OE2	2.24	0.70
1:B:388:ALA:HA	1:B:391:LEU:HD12	1.74	0.69
1:C:417:LYS:HG2	1:C:423:TYR:HD2	1.58	0.68
1:D:195:LYS:HB2	1:D:204:ILE:HG23	1.75	0.68
1:C:55:ARG:HH21	1:C:65:THR:HG21	1.57	0.68
1:D:186:LEU:HB2	1:D:368:LEU:HD21	1.75	0.68
1:B:409:ARG:O	1:B:413:MET:HG2	1.93	0.68
1:D:341:GLY:HA3	1:D:348:LEU:HD21	1.75	0.68
1:D:353:MET:HG2	1:D:378:ARG:HA	1.74	0.68
1:B:358:PHE:CE2	1:D:161:ASP:HB2	2.29	0.68
1:B:161:ASP:HB2	1:D:358:PHE:CE2	2.29	0.68
1:A:206:THR:O	1:A:425:ASN:ND2	2.22	0.67
1:D:312:ARG:NH1	1:D:315:ALA:CA	2.57	0.67
1:A:219:ARG:NH2	1:A:340:GLU:OE2	2.23	0.67
1:D:256:ALA:O	1:D:259:THR:HG22	1.95	0.67
1:C:303:ARG:H	1:C:305:LYS:HG3	1.60	0.66
1:C:151:HIS:N	1:C:187:GLU:OE1	2.26	0.66
1:B:383:ASP:HB3	1:B:386:LYS:HG3	1.75	0.66
1:B:369:ARG:NH1	1:C:369:ARG:HD2	2.11	0.66
1:C:92:VAL:HG22	1:C:164:PHE:CE2	2.30	0.66
1:D:282:LYS:HG2	1:D:283:LYS:N	2.11	0.66
1:A:152:SER:O	1:A:156:MET:HG3	1.97	0.65
1:A:82:SER:HB3	1:A:84:GLN:OE1	1.97	0.65
1:B:239:GLU:HB3	1:B:240:ASP:HB3	1.77	0.65
1:B:353:MET:O	1:B:378:ARG:HB2	1.97	0.65
1:B:321:LEU:HD21	1:B:341:GLY:HA2	1.79	0.65
1:A:18:GLN:HG3	1:A:83:LEU:HD23	1.77	0.64

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:46:LYS:HZ2	1:D:81:GLU:HA	1.62	0.64
1:C:146:LEU:HB2	1:C:180:ILE:HG22	1.79	0.64
1:A:163:SER:HA	1:A:166:ASN:H	1.62	0.63
1:A:310:VAL:HG13	1:A:312:ARG:HH12	1.62	0.63
1:B:307:LYS:HA	1:B:309:LEU:HB3	1.80	0.63
1:C:173:THR:HG22	1:C:198:PHE:O	1.98	0.63
1:A:14:ILE:HG12	1:A:42:LEU:HB3	1.80	0.63
1:D:365:VAL:HG12	1:D:371:ALA:HB3	1.81	0.63
1:C:30:HIS:HB2	1:C:54:LEU:HD21	1.82	0.62
1:C:374:VAL:HG12	1:C:390:ILE:HD13	1.81	0.62
1:D:72:SER:O	1:D:100:ARG:NH2	2.33	0.62
1:B:54:LEU:HD13	1:B:213:PHE:CE2	2.34	0.62
1:C:353:MET:HG2	1:C:378:ARG:HA	1.82	0.62
1:B:401:GLU:OE2	1:C:369:ARG:NE	2.25	0.62
1:C:112:GLN:HG2	1:C:116:ASP:HB2	1.82	0.62
1:C:356:ASP:HB2	1:C:360:ASN:ND2	2.14	0.61
1:D:186:LEU:HA	1:D:188:PRO:HD3	1.81	0.61
1:B:372:VAL:HG23	1:B:402:ARG:HH21	1.65	0.60
1:D:162:ARG:HD2	1:D:169:GLU:OE1	2.00	0.60
1:C:10:HIS:HB3	1:C:118:PRO:HA	1.84	0.60
1:C:33:LEU:HD13	1:C:58:HIS:CD2	2.36	0.60
1:B:311:ILE:HG22	1:B:312:ARG:H	1.66	0.60
1:B:80:VAL:HG11	1:B:86:LEU:HD21	1.84	0.60
1:B:212:PRO:HG3	1:B:219:ARG:NE	2.16	0.59
1:C:164:PHE:HD1	1:C:164:PHE:H	1.50	0.59
1:B:276:ALA:HA	1:B:312:ARG:HG3	1.85	0.59
1:C:307:LYS:HA	1:C:309:LEU:HD13	1.84	0.59
1:B:230:VAL:HG21	1:B:312:ARG:HH12	1.68	0.59
1:C:223:SER:OG	3:C:1001:HOH:O	2.02	0.59
1:B:163:SER:OG	1:B:164:PHE:HA	2.04	0.58
1:C:400:PRO:O	1:C:403:VAL:HG12	2.03	0.58
1:D:365:VAL:HG23	1:D:366:ASP:OD1	2.04	0.58
1:A:378:ARG:NH1	1:C:163:SER:OG	2.37	0.57
1:A:18:GLN:NE2	1:A:251:GLN:OE1	2.33	0.57
1:D:408:LEU:HD23	1:D:411:LYS:HD3	1.86	0.57
1:C:186:LEU:HD21	1:C:335:TRP:HH2	1.68	0.57
1:D:28:LEU:HD21	1:D:122:LEU:HD21	1.87	0.57
1:D:163:SER:OG	1:D:164:PHE:HA	2.04	0.57
1:D:312:ARG:CZ	1:D:315:ALA:HA	2.35	0.57
1:A:389:ARG:HG3	1:A:390:ILE:N	2.20	0.56
1:A:356:ASP:O	1:A:360:ASN:ND2	2.39	0.56

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:347:MET:CB	1:B:402:ARG:HH11	2.19	0.56
1:D:149:ASN:HB3	1:D:335:TRP:CE2	2.41	0.56
1:C:269:SER:HB2	1:C:392:ALA:HB2	1.87	0.56
1:C:329:TYR:CZ	1:C:331:THR:HG23	2.40	0.56
1:D:363:LEU:HD22	1:D:364:ILE:HD13	1.87	0.56
1:A:304:VAL:HA	1:A:306:GLU:OE1	2.06	0.55
1:C:170:THR:O	1:C:173:THR:OG1	2.16	0.55
1:C:331:THR:CG2	1:C:333:LEU:H	2.19	0.55
1:D:342:MET:HA	1:D:405:LEU:HD11	1.87	0.55
1:A:306:GLU:HB2	1:A:309:LEU:HG	1.88	0.55
1:B:223:SER:HA	1:B:314:TRP:HD1	1.70	0.55
1:A:340:GLU:OE1	2:A:900:UDP:O2'	2.21	0.55
1:C:372:VAL:HG11	1:C:390:ILE:HD11	1.89	0.55
1:A:161:ASP:HB2	1:C:358:PHE:CD2	2.41	0.55
1:B:299:GLY:O	3:B:1001:HOH:O	2.18	0.55
1:D:358:PHE:O	1:D:361:THR:HG22	2.07	0.55
1:D:83:LEU:HB2	1:D:94:MET:HG3	1.89	0.55
1:B:343:VAL:HA	1:B:409:ARG:HG3	1.88	0.54
1:C:407:LYS:HE2	1:C:411:LYS:HE3	1.89	0.54
1:B:168:LEU:O	1:B:172:THR:OG1	2.13	0.54
1:B:217:VAL:HG13	1:B:218:ASP:H	1.72	0.54
1:D:165:PHE:O	1:D:169:GLU:HG3	2.07	0.54
1:A:383:ASP:OD2	1:A:386:LYS:HG3	2.08	0.54
1:C:365:VAL:HG13	1:C:371:ALA:O	2.07	0.54
1:C:409:ARG:O	1:C:413:MET:HG2	2.07	0.54
1:A:396:ARG:HB3	1:A:399:LEU:HD13	1.90	0.54
1:B:235:ASP:OD2	1:B:323:HIS:NE2	2.39	0.54
1:B:92:VAL:HG22	1:B:164:PHE:CZ	2.43	0.54
1:D:279:ASP:OD1	1:D:280:ALA:N	2.32	0.54
1:A:16:PHE:O	1:A:18:GLN:N	2.37	0.54
1:C:185:ASP:HA	1:C:188:PRO:HG3	1.90	0.53
1:D:245:TYR:OH	1:D:317:GLN:HG2	2.07	0.53
1:A:250:SER:HB2	2:A:900:UDP:O3B	2.08	0.53
1:D:126:PHE:CD1	1:D:168:LEU:HD11	2.44	0.53
1:B:234:LEU:HD21	1:B:274:ILE:HD11	1.91	0.53
1:C:33:LEU:HD11	1:C:63:PHE:HD1	1.74	0.53
1:B:402:ARG:O	1:B:405:LEU:HB3	2.08	0.53
1:C:190:PHE:O	1:C:194:VAL:HG12	2.08	0.53
1:D:205:TRP:HH2	1:D:432:GLU:HG3	1.73	0.53
1:A:234:LEU:HB3	1:A:323:HIS:NE2	2.23	0.53
1:D:409:ARG:HH11	1:D:413:MET:HG3	1.74	0.53

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:416:ILE:HA	1:C:422:SER:HB2	1.91	0.53
1:A:66:LEU:HD11	1:A:68:LEU:HD21	1.91	0.52
1:D:121:ILE:HD12	1:D:135:ALA:HB2	1.91	0.52
1:D:412:ALA:O	1:D:416:ILE:HG12	2.09	0.52
1:B:377:ASN:N	1:B:380:SER:OG	2.30	0.52
1:D:164:PHE:H	1:D:164:PHE:HD1	1.56	0.52
1:A:163:SER:HA	1:A:165:PHE:N	2.24	0.52
1:B:195:LYS:HE3	1:B:204:ILE:HG22	1.91	0.52
1:C:303:ARG:C	1:C:305:LYS:H	2.12	0.52
1:D:265:ALA:HA	1:D:384:SER:OG	2.09	0.52
1:D:141:LYS:HD3	1:D:176:TYR:CZ	2.44	0.52
1:B:152:SER:O	1:B:156:MET:HG3	2.10	0.52
1:D:55:ARG:HG3	1:D:63:PHE:CE2	2.45	0.52
1:D:255:THR:HG22	1:D:258:GLN:HB2	1.92	0.52
1:A:173:THR:HG23	1:A:198:PHE:O	2.11	0.51
1:B:266:LEU:HD22	1:B:273:PHE:CG	2.45	0.51
1:C:279:ASP:OD1	1:C:280:ALA:N	2.40	0.51
1:B:372:VAL:CG2	1:B:402:ARG:HH21	2.23	0.51
1:C:7:LYS:HD3	1:C:35:ARG:O	2.11	0.51
1:B:102:HIS:ND1	1:B:103:ASP:OD1	2.39	0.51
1:B:268:LYS:NZ	1:B:384:SER:HB2	2.26	0.51
1:D:379:ASP:O	1:D:381:VAL:HG23	2.10	0.51
1:A:358:PHE:CE2	1:C:161:ASP:HB2	2.46	0.51
1:D:189:GLU:O	1:D:193:THR:OG1	2.24	0.51
1:D:195:LYS:O	1:D:202:HIS:ND1	2.44	0.51
1:B:332:HIS:HD2	1:B:357:HIS:NE2	2.10	0.50
1:C:163:SER:OG	1:C:164:PHE:HA	2.12	0.50
1:C:43:VAL:HG21	1:C:48:SER:HA	1.93	0.50
1:D:55:ARG:HG3	1:D:63:PHE:HE2	1.76	0.50
1:B:155:VAL:O	1:B:159:GLN:HG3	2.12	0.50
1:C:302:GLU:HB2	1:C:303:ARG:HG2	1.94	0.50
1:D:312:ARG:NH1	1:D:314:TRP:C	2.63	0.50
1:C:416:ILE:O	1:C:423:TYR:HB2	2.12	0.50
1:D:151:HIS:HB3	1:D:363:LEU:HD12	1.94	0.50
1:D:163:SER:HA	1:D:166:ASN:H	1.77	0.49
1:D:249:GLY:O	1:D:278:ARG:HD3	2.12	0.49
1:D:86:LEU:HB2	1:D:87:PRO:CD	2.42	0.49
1:B:153:ILE:HG23	1:B:157:TRP:CE3	2.48	0.49
1:A:377:ASN:H	1:A:380:SER:HB2	1.77	0.49
1:C:10:HIS:HD2	1:C:38:THR:HG22	1.77	0.49
1:D:170:THR:O	1:D:174:GLU:HG2	2.13	0.49

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:ASP:HB2	1:D:358:PHE:HE2	1.75	0.49
1:C:112:GLN:NE2	1:C:112:GLN:CA	2.67	0.49
1:D:356:ASP:OD2	1:D:360:ASN:ND2	2.46	0.49
1:A:311:ILE:O	1:A:312:ARG:HD3	2.13	0.49
1:C:212:PRO:HG3	1:C:219:ARG:CZ	2.42	0.49
1:D:107:ASP:OD1	1:D:111:ARG:NH1	2.45	0.49
1:D:416:ILE:O	1:D:423:TYR:HB2	2.12	0.49
1:D:9:PRO:HD2	1:D:37:ALA:HA	1.94	0.49
1:A:255:THR:HG22	1:A:256:ALA:N	2.28	0.49
1:C:234:LEU:HD11	1:C:320:ILE:HG23	1.94	0.49
1:C:335:TRP:NE1	1:C:339:LEU:HD11	2.28	0.49
1:D:186:LEU:HB3	1:D:187:GLU:CG	2.40	0.49
1:D:312:ARG:NH2	1:D:316:PRO:HD3	2.28	0.49
1:D:386:LYS:O	1:D:390:ILE:HG13	2.12	0.49
1:A:132:ASN:O	1:A:136:ASP:HB2	2.13	0.48
1:D:43:VAL:HG23	1:D:67:ILE:HG12	1.96	0.48
1:A:209:PRO:HG3	1:A:339:LEU:HD22	1.93	0.48
1:D:254:LEU:HB3	1:D:258:GLN:HB3	1.93	0.48
1:D:282:LYS:O	1:D:283:LYS:HB2	2.13	0.48
1:D:86:LEU:HB2	1:D:87:PRO:HD2	1.94	0.48
1:B:165:PHE:HA	1:B:168:LEU:HD13	1.94	0.48
1:B:52:ASP:OD1	1:B:55:ARG:NH1	2.46	0.48
1:A:20:GLY:HA3	2:A:900:UDP:O1B	2.13	0.48
1:C:46:LYS:HE3	1:C:79:GLY:O	2.13	0.48
1:B:306:GLU:O	1:B:309:LEU:HD13	2.14	0.48
1:D:255:THR:CG2	1:D:258:GLN:HB2	2.44	0.48
1:A:30:HIS:HE1	1:A:34:LEU:HD21	1.79	0.48
1:D:88:LEU:HD22	1:D:354:GLN:OE1	2.14	0.47
1:A:254:LEU:HB2	1:A:259:THR:HG23	1.95	0.47
1:C:274:ILE:HG23	1:C:312:ARG:HH12	1.79	0.47
1:D:426:LEU:HA	1:D:426:LEU:HD23	1.64	0.47
1:A:75:CYS:CB	1:C:75:CYS:HG	2.27	0.47
1:A:164:PHE:H	1:A:164:PHE:HD1	1.63	0.47
1:B:150:ALA:N	1:B:187:GLU:OE2	2.43	0.47
1:A:274:ILE:HG12	1:A:310:VAL:HG12	1.96	0.47
1:B:184:TYR:OH	1:B:192:GLU:OE2	2.21	0.47
1:D:203:ARG:HD2	1:D:203:ARG:HA	1.69	0.47
1:B:274:ILE:HG23	1:B:310:VAL:HB	1.96	0.47
1:B:347:MET:HA	1:B:402:ARG:HD2	1.97	0.47
1:A:239:GLU:O	1:A:242:SER:OG	2.20	0.47
1:A:333:LEU:HD13	1:A:350:ALA:HB2	1.95	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:250:SER:OG	2:B:900:UDP:O3B	2.23	0.47
1:A:212:PRO:HB3	1:A:219:ARG:CD	2.45	0.47
1:A:234:LEU:HB3	1:A:323:HIS:CD2	2.50	0.47
1:B:163:SER:HB3	1:D:378:ARG:NH1	2.29	0.47
1:A:245:TYR:OH	1:A:317:GLN:HG2	2.15	0.47
1:B:316:PRO:HD2	1:B:320:ILE:HD11	1.97	0.47
1:B:331:THR:HG22	1:B:333:LEU:H	1.79	0.47
1:D:264:ALA:HA	1:D:267:GLU:CD	2.35	0.47
1:B:212:PRO:HG3	1:B:219:ARG:HE	1.80	0.46
1:A:410:GLU:O	1:A:414:GLU:HG2	2.15	0.46
1:C:340:GLU:OE2	2:C:900:UDP:O3'	2.30	0.46
1:D:65:THR:HG22	1:D:67:ILE:HG13	1.96	0.46
1:D:11:VAL:HG21	1:D:32:ILE:HD12	1.96	0.46
1:D:258:GLN:HE22	1:D:379:ASP:HA	1.81	0.46
1:A:19:SER:OG	1:A:84:GLN:HG2	2.16	0.46
1:D:304:VAL:HG13	1:D:305:LYS:H	1.80	0.46
1:A:241:ASN:HD21	1:A:402:ARG:HH22	1.64	0.46
1:A:266:LEU:HD22	1:A:273:PHE:CG	2.50	0.46
1:B:18:GLN:NE2	1:B:251:GLN:HG2	2.31	0.46
1:C:16:PHE:O	1:C:18:GLN:N	2.46	0.46
1:D:84:GLN:HG2	1:D:84:GLN:O	2.16	0.46
1:A:274:ILE:HG23	1:A:310:VAL:HG13	1.96	0.46
1:A:383:ASP:HB3	1:A:386:LYS:HB2	1.98	0.46
1:C:240:ASP:O	1:C:325:ALA:HA	2.16	0.46
1:D:409:ARG:NH1	1:D:413:MET:HG3	2.30	0.46
1:B:424:LYS:HD2	1:B:424:LYS:HA	1.64	0.45
1:C:117:LEU:HA	1:C:117:LEU:HD23	1.82	0.45
1:C:186:LEU:HD21	1:C:335:TRP:CH2	2.50	0.45
1:D:311:ILE:O	1:D:312:ARG:HG2	2.15	0.45
1:A:150:ALA:O	1:A:154:SER:OG	2.33	0.45
1:B:255:THR:HG22	1:B:256:ALA:N	2.32	0.45
1:C:255:THR:HG23	1:C:258:GLN:OE1	2.16	0.45
1:C:184:TYR:HD2	1:C:185:ASP:OD1	1.99	0.45
1:C:317:GLN:OE1	2:C:900:UDP:H2'	2.16	0.45
1:D:153:ILE:HG23	1:D:157:TRP:CE3	2.51	0.45
1:D:303:ARG:HA	1:D:304:VAL:HA	1.43	0.45
1:C:312:ARG:HB3	1:C:313:GLY:HA3	1.99	0.45
1:B:183:PHE:CZ	1:B:186:LEU:HB2	2.51	0.45
1:B:190:PHE:O	1:B:194:VAL:HG23	2.16	0.45
1:B:353:MET:HG2	1:B:378:ARG:HA	1.98	0.45
1:C:162:ARG:HD2	1:C:169:GLU:OE1	2.17	0.45

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:185:ASP:HB2	1:D:186:LEU:HD23	1.99	0.45
1:A:273:PHE:O	1:A:309:LEU:HA	2.17	0.44
1:A:332:HIS:CE1	1:A:334:GLY:HA3	2.52	0.44
1:B:245:TYR:OH	1:B:317:GLN:HG2	2.18	0.44
1:A:212:PRO:HB3	1:A:219:ARG:HD3	1.99	0.44
1:A:48:SER:O	1:A:51:LEU:HB2	2.17	0.44
1:B:173:THR:HG23	1:B:198:PHE:O	2.17	0.44
1:C:7:LYS:HA	1:C:8:LYS:HA	1.56	0.44
1:A:311:ILE:HG22	1:A:312:ARG:H	1.83	0.44
1:A:173:THR:HG23	1:A:198:PHE:C	2.38	0.44
1:B:266:LEU:HB3	1:B:273:PHE:CE1	2.52	0.44
1:B:18:GLN:HE21	1:B:251:GLN:HG2	1.83	0.44
1:C:417:LYS:HG2	1:C:423:TYR:CD2	2.46	0.44
1:C:113:PRO:O	1:C:116:ASP:N	2.49	0.44
1:D:7:LYS:HB3	1:D:35:ARG:O	2.18	0.44
1:C:140:ILE:HG22	1:C:141:LYS:O	2.18	0.44
1:C:223:SER:HA	1:C:314:TRP:O	2.18	0.44
1:C:242:SER:O	1:C:242:SER:OG	2.35	0.44
1:C:307:LYS:HA	1:C:309:LEU:HB2	2.00	0.44
1:D:261:ALA:O	1:D:264:ALA:HB3	2.18	0.44
1:A:52:ASP:HA	1:A:55:ARG:NH1	2.33	0.44
1:D:88:LEU:HG	1:D:88:LEU:O	2.17	0.44
1:C:20:GLY:HA2	2:C:900:UDP:H5'1	1.98	0.44
1:A:367:LYS:HE2	1:A:367:LYS:HB3	1.87	0.43
1:B:358:PHE:CD2	1:D:161:ASP:HB2	2.52	0.43
1:D:387:LEU:HD12	1:D:390:ILE:HD12	2.00	0.43
1:B:108:PHE:O	1:B:112:GLN:HG2	2.18	0.43
1:B:310:VAL:HG12	1:B:311:ILE:H	1.83	0.43
1:B:70:PHE:CD2	1:B:81:GLU:HB3	2.54	0.43
1:A:283:LYS:HB3	1:A:283:LYS:NZ	2.33	0.43
1:A:273:PHE:CZ	1:A:309:LEU:HD23	2.54	0.43
1:C:219:ARG:NH1	1:C:318:THR:HG21	2.21	0.43
1:D:127:LEU:O	1:D:131:ILE:HG13	2.18	0.43
1:D:128:SER:HA	1:D:131:ILE:HB	2.00	0.43
1:D:21:HIS:CE1	1:D:126:PHE:CE2	3.06	0.43
1:A:16:PHE:CD1	1:A:17:PRO:HD2	2.53	0.43
1:A:387:LEU:O	1:A:391:LEU:HG	2.18	0.43
1:B:113:PRO:O	1:B:116:ASP:N	2.51	0.43
1:C:169:GLU:O	1:C:173:THR:HG23	2.19	0.43
1:C:352:PRO:HA	1:C:357:HIS:HB3	2.01	0.43
1:C:61:GLU:O	1:C:64:LYS:NZ	2.47	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:413:MET:SD	1:D:416:ILE:HD11	2.59	0.43
1:B:239:GLU:HA	1:B:240:ASP:HA	1.56	0.43
1:C:224:SER:OG	1:C:313:GLY:HA3	2.18	0.43
1:D:340:GLU:OE2	2:D:900:UDP:O3'	2.26	0.43
1:B:212:PRO:HG3	1:B:219:ARG:CZ	2.49	0.43
1:B:246:VAL:HG13	1:B:330:LEU:HD23	2.00	0.43
1:A:270:SER:OG	1:A:270:SER:O	2.35	0.42
1:A:332:HIS:CD2	1:A:332:HIS:H	2.36	0.42
1:B:304:VAL:HG11	1:B:309:LEU:HD22	2.01	0.42
1:A:146:LEU:HD13	1:A:148:ILE:HD11	2.01	0.42
1:B:321:LEU:CD2	1:B:341:GLY:HA2	2.49	0.42
1:B:426:LEU:HA	1:B:426:LEU:HD23	1.89	0.42
1:C:235:ASP:OD1	1:C:323:HIS:NE2	2.47	0.42
1:A:31:GLN:OE1	1:A:35:ARG:NH2	2.48	0.42
1:D:45:PRO:HD3	1:D:68:LEU:O	2.19	0.42
1:B:307:LYS:HG3	1:B:307:LYS:O	2.19	0.42
1:C:332:HIS:HD2	1:C:357:HIS:NE2	2.18	0.42
1:C:55:ARG:NH2	1:C:65:THR:HG21	2.30	0.42
1:D:227:PRO:HB3	1:D:319:MET:HE1	2.02	0.42
1:B:341:GLY:O	1:B:346:VAL:HG22	2.20	0.42
1:C:240:ASP:O	1:C:242:SER:N	2.49	0.42
1:D:117:LEU:HA	1:D:117:LEU:HD23	1.91	0.42
1:A:75:CYS:SG	1:C:75:CYS:SG	3.16	0.42
1:B:373:ARG:HH21	1:B:376:GLU:CA	2.29	0.42
1:C:354:GLN:O	1:C:357:HIS:CG	2.73	0.42
1:A:303:ARG:HB2	1:A:305:LYS:HG2	2.01	0.42
1:D:77:PRO:O	1:D:80:VAL:HG12	2.20	0.42
1:A:234:LEU:O	1:A:237:CYS:N	2.42	0.42
1:A:304:VAL:O	1:A:304:VAL:HG23	2.20	0.42
1:B:372:VAL:HG23	1:B:402:ARG:NH2	2.32	0.42
1:B:162:ARG:HD2	1:B:169:GLU:OE1	2.20	0.41
1:B:329:TYR:CZ	1:B:331:THR:HG23	2.55	0.41
1:C:209:PRO:HD2	1:C:416:ILE:HG23	2.01	0.41
1:C:98:LEU:HD23	1:C:98:LEU:HA	1.73	0.41
1:B:13:VAL:HG12	1:B:15:PRO:HD3	2.03	0.41
1:C:266:LEU:HG	1:C:391:LEU:HD21	2.02	0.41
1:D:250:SER:O	1:D:278:ARG:NE	2.53	0.41
1:B:178:LEU:HG	1:B:180:ILE:HG23	2.01	0.41
1:A:193:THR:HG22	1:D:189:GLU:OE1	2.20	0.41
1:B:225:ILE:HG23	1:B:229:LYS:NZ	2.36	0.41
1:B:365:VAL:HG12	1:B:366:ASP:OD1	2.20	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:372:VAL:HG23	1:D:374:VAL:HG13	2.03	0.41
1:D:393:GLU:OE2	1:D:399:LEU:HD21	2.20	0.41
1:D:46:LYS:HD2	1:D:81:GLU:HA	2.02	0.41
1:C:68:LEU:HD13	1:C:104:PRO:HB2	2.03	0.41
1:C:263:ALA:HB2	1:C:275:TRP:CZ3	2.55	0.41
1:D:185:ASP:HB3	1:D:188:PRO:HG3	2.01	0.41
1:A:44:THR:HB	1:A:45:PRO:CD	2.50	0.41
1:C:70:PHE:CD2	1:C:81:GLU:HB3	2.56	0.41
1:D:140:ILE:HD13	1:D:140:ILE:HG21	1.75	0.41
1:D:41:VAL:HG12	1:D:43:VAL:HG13	2.02	0.41
1:A:163:SER:HB3	1:A:166:ASN:HB2	2.02	0.41
1:A:269:SER:HB3	1:A:388:ALA:O	2.21	0.41
1:C:157:TRP:CZ2	1:C:169:GLU:HA	2.56	0.41
1:D:354:GLN:O	1:D:357:HIS:CG	2.74	0.41
1:D:386:LYS:O	1:D:389:ARG:HG2	2.21	0.41
1:A:189:GLU:O	1:A:193:THR:HG23	2.20	0.41
1:A:231:SER:OG	1:A:319:MET:SD	2.79	0.41
1:A:235:ASP:HA	1:A:236:SER:HA	1.55	0.41
1:A:238:PRO:HB2	1:A:239:GLU:HG3	2.02	0.41
1:C:302:GLU:HA	1:C:303:ARG:HA	1.46	0.41
1:C:54:LEU:HD23	1:C:54:LEU:HA	1.80	0.41
1:D:263:ALA:HA	1:D:266:LEU:HD13	2.02	0.41
1:A:429:LEU:O	1:A:433:MET:HG3	2.20	0.41
1:D:331:THR:HG23	1:D:333:LEU:H	1.85	0.41
1:B:255:THR:HG22	1:B:256:ALA:H	1.86	0.41
1:D:224:SER:HB3	1:D:314:TRP:O	2.21	0.41
1:D:46:LYS:H	1:D:46:LYS:HG2	1.40	0.41
1:A:16:PHE:CE1	1:A:95:PHE:HE1	2.39	0.40
1:C:361:THR:HG21	1:C:373:ARG:HD2	2.03	0.40
1:D:178:LEU:O	1:D:205:TRP:HD1	2.04	0.40
1:D:183:PHE:CD2	1:D:185:ASP:O	2.74	0.40
1:A:360:ASN:HD22	1:A:360:ASN:N	2.20	0.40
1:A:377:ASN:H	1:A:380:SER:CB	2.33	0.40
1:B:235:ASP:HA	1:B:236:SER:HA	1.68	0.40
1:C:271:VAL:CG1	1:C:392:ALA:HA	2.50	0.40
1:D:186:LEU:HA	1:D:187:GLU:HA	1.80	0.40
1:D:195:LYS:HB2	1:D:204:ILE:CG2	2.48	0.40
1:C:356:ASP:O	1:C:359:PHE:N	2.54	0.40
1:C:55:ARG:HD2	1:C:55:ARG:HH11	1.74	0.40
1:D:151:HIS:O	1:D:154:SER:OG	2.38	0.40
1:D:351:TRP:C	1:D:351:TRP:CD1	2.93	0.40

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:46:LYS:NZ	1:D:81:GLU:HA	2.33	0.40
1:B:220:GLY:HA3	1:B:314:TRP:HZ2	1.86	0.40
1:C:165:PHE:O	1:C:169:GLU:HG3	2.22	0.40
1:D:234:LEU:HD23	1:D:323:HIS:CG	2.55	0.40
1:D:271:VAL:O	1:D:273:PHE:HD1	2.04	0.40
1:B:416:ILE:O	1:B:423:TYR:HB2	2.22	0.40
1:D:354:GLN:HB3	1:D:354:GLN:HE21	1.51	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	404/435 (93%)	382 (95%)	21 (5%)	1 (0%)	47 82
1	B	400/435 (92%)	378 (94%)	20 (5%)	2 (0%)	29 68
1	C	407/435 (94%)	380 (93%)	23 (6%)	4 (1%)	15 53
1	D	376/435 (86%)	355 (94%)	19 (5%)	2 (0%)	29 68
All	All	1587/1740 (91%)	1495 (94%)	83 (5%)	9 (1%)	25 64

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	304	VAL
1	B	305	LYS
1	C	240	ASP
1	C	238	PRO
1	C	305	LYS
1	D	313	GLY
1	B	223	SER
1	A	17	PRO

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	D	17	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	351/377 (93%)	340 (97%)	11 (3%)	40	75
1	B	348/377 (92%)	341 (98%)	7 (2%)	55	83
1	C	357/377 (95%)	353 (99%)	4 (1%)	73	90
1	D	336/377 (89%)	329 (98%)	7 (2%)	53	82
All	All	1392/1508 (92%)	1363 (98%)	29 (2%)	53	82

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	SER
1	A	103	ASP
1	A	136	ASP
1	A	141	LYS
1	A	197	ARG
1	A	241	ASN
1	A	253	ARG
1	A	282	LYS
1	A	300	PHE
1	A	303	ARG
1	A	398	ASP
1	B	72	SER
1	B	82	SER
1	B	185	ASP
1	B	219	ARG
1	B	282	LYS
1	B	300	PHE
1	B	312	ARG
1	C	142	SER
1	C	213	PHE

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	219	ARG
1	C	305	LYS
1	D	72	SER
1	D	86	LEU
1	D	269	SER
1	D	312	ARG
1	D	378	ARG
1	D	409	ARG
1	D	434	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	323	HIS
1	A	360	ASN
1	B	84	GLN
1	B	251	GLN
1	C	112	GLN
1	D	354	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
2	UDP	A	900	-	20,26,26	0.97	1 (5%)	25,40,40	1.03	1 (4%)
2	UDP	C	900	-	20,26,26	1.09	1 (5%)	25,40,40	1.17	2 (8%)
2	UDP	B	900	-	20,26,26	0.96	1 (5%)	25,40,40	1.07	2 (8%)
2	UDP	D	900	-	20,26,26	0.96	1 (5%)	25,40,40	1.19	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	UDP	A	900	-	-	3/14/32/32	0/2/2/2
2	UDP	C	900	-	-	6/14/32/32	0/2/2/2
2	UDP	B	900	-	-	3/14/32/32	0/2/2/2
2	UDP	D	900	-	-	1/14/32/32	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	900	UDP	C4-N3	3.16	1.38	1.33
2	A	900	UDP	C4-N3	3.00	1.38	1.33
2	B	900	UDP	C4-N3	2.91	1.38	1.33
2	D	900	UDP	C4-N3	2.82	1.37	1.33

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	900	UDP	C5-C4-N3	-3.93	114.66	123.31
2	C	900	UDP	C5-C4-N3	-3.93	114.67	123.31
2	D	900	UDP	C5-C4-N3	-3.85	114.83	123.31
2	B	900	UDP	C5-C4-N3	-3.82	114.91	123.31
2	B	900	UDP	PA-O3A-PB	-2.66	123.68	132.83
2	C	900	UDP	O3B-PB-O3A	2.30	112.36	104.64
2	D	900	UDP	O3B-PB-O3A	2.30	112.33	104.64

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	900	UDP	O4'-C4'-C5'-O5'
2	C	900	UDP	C5'-O5'-PA-O3A
2	C	900	UDP	O4'-C4'-C5'-O5'
2	B	900	UDP	C3'-C4'-C5'-O5'
2	B	900	UDP	O4'-C4'-C5'-O5'
2	A	900	UDP	C3'-C4'-C5'-O5'
2	C	900	UDP	C3'-C4'-C5'-O5'
2	A	900	UDP	PB-O3A-PA-O5'
2	B	900	UDP	PB-O3A-PA-O5'
2	C	900	UDP	C5'-O5'-PA-O1A
2	C	900	UDP	C5'-O5'-PA-O2A
2	C	900	UDP	PB-O3A-PA-O1A
2	D	900	UDP	O4'-C4'-C5'-O5'

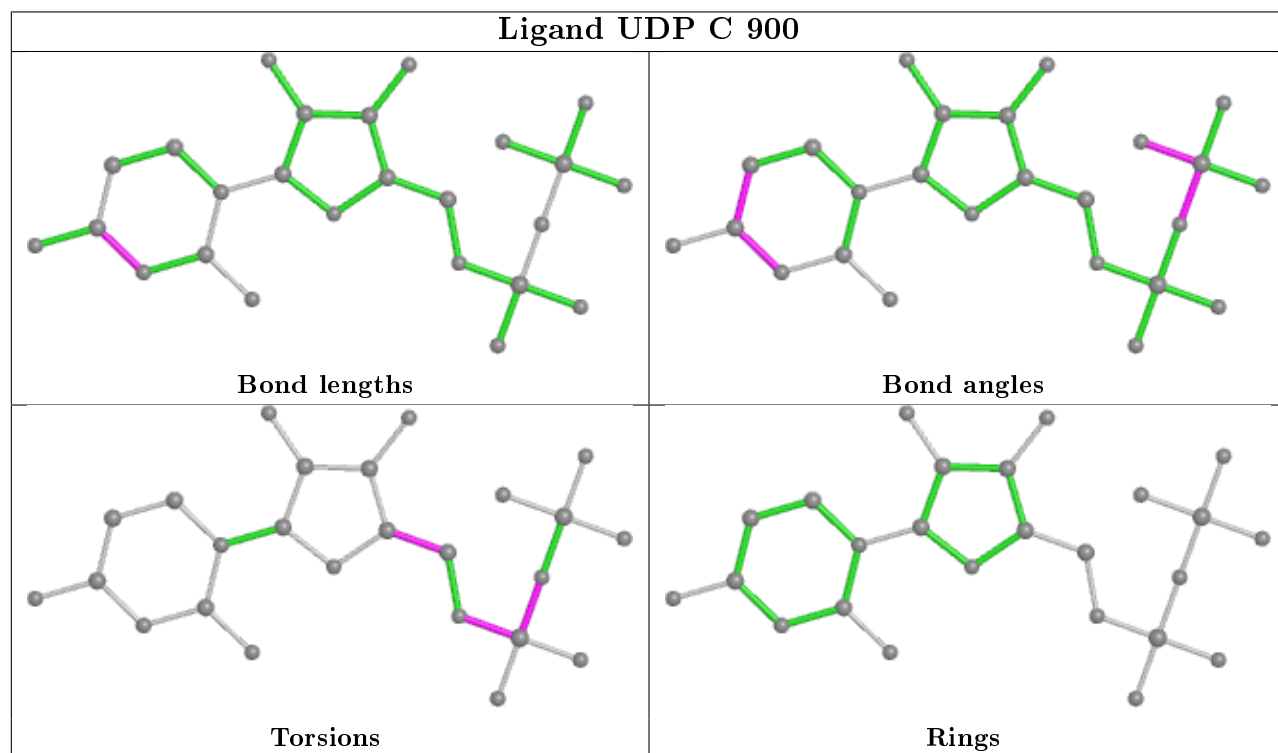
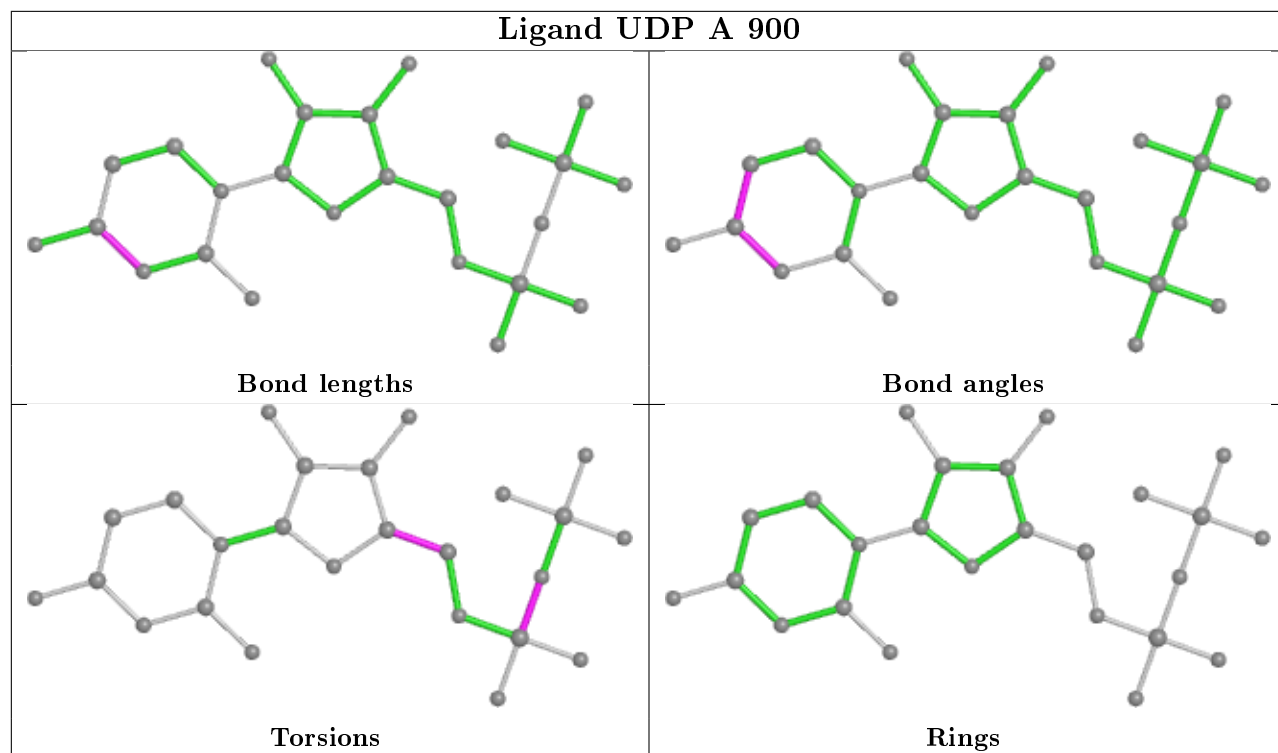
There are no ring outliers.

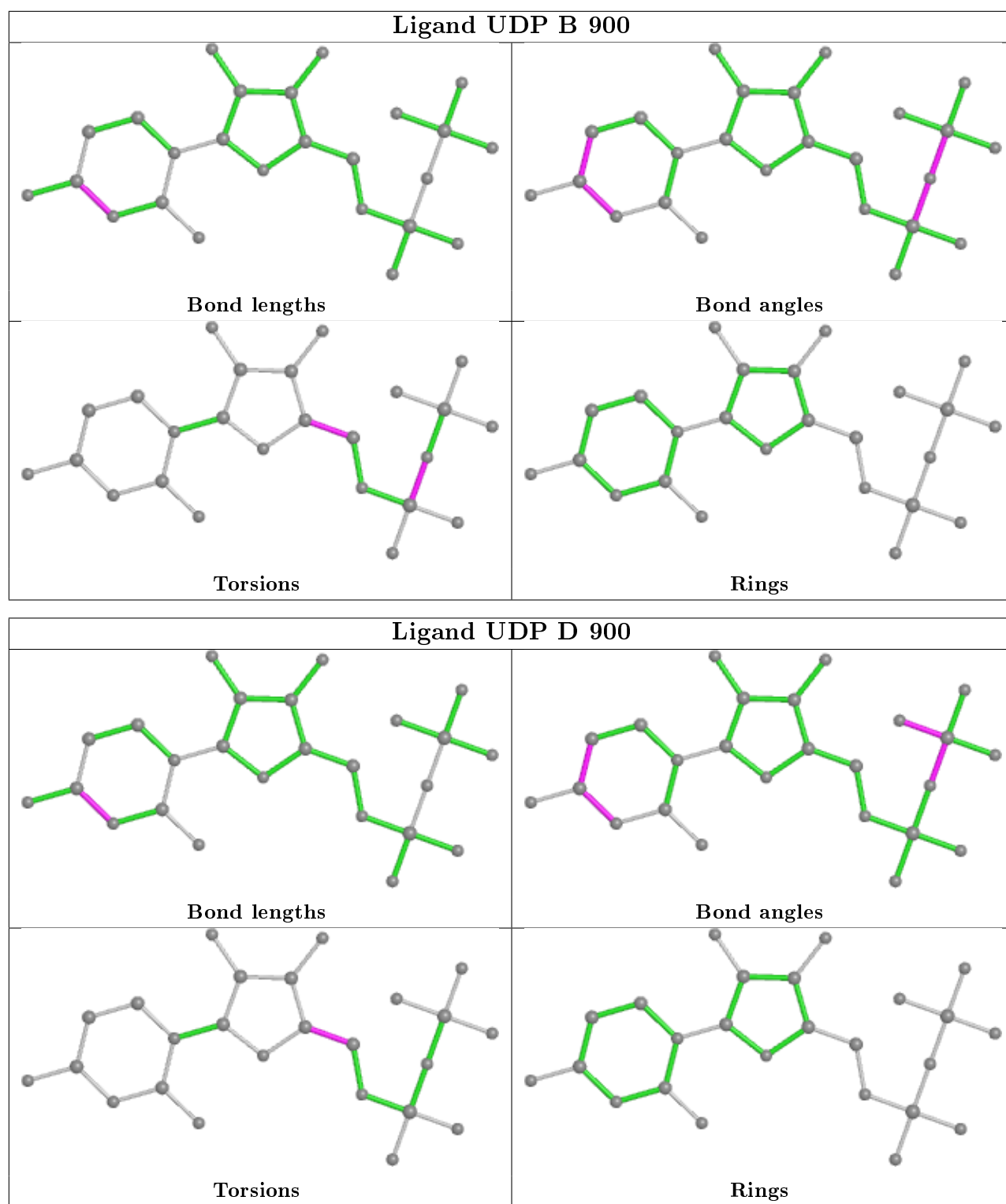
4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	900	UDP	3	0
2	C	900	UDP	3	0
2	B	900	UDP	1	0
2	D	900	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<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	410/435 (94%)	-0.41	3 (0%) 87 69	17, 30, 58, 70	0
1	B	408/435 (93%)	-0.28	2 (0%) 91 75	15, 39, 82, 96	0
1	C	413/435 (94%)	-0.37	4 (0%) 82 59	14, 33, 73, 90	0
1	D	388/435 (89%)	-0.18	7 (1%) 68 40	21, 42, 87, 102	0
All	All	1619/1740 (93%)	-0.31	16 (0%) 82 59	14, 36, 79, 102	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	224	SER	2.9
1	D	226	PRO	2.9
1	A	309	LEU	2.8
1	A	223	SER	2.6
1	C	303	ARG	2.5
1	D	228	ALA	2.5
1	D	259	THR	2.4
1	B	398	ASP	2.3
1	D	387	LEU	2.3
1	A	238	PRO	2.2
1	C	297	PRO	2.2
1	D	229	LYS	2.1
1	C	302	GLU	2.1
1	C	222	GLN	2.1
1	D	227	PRO	2.1
1	B	410	GLU	2.0

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

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

## 6.3 Carbohydrates [i](#)

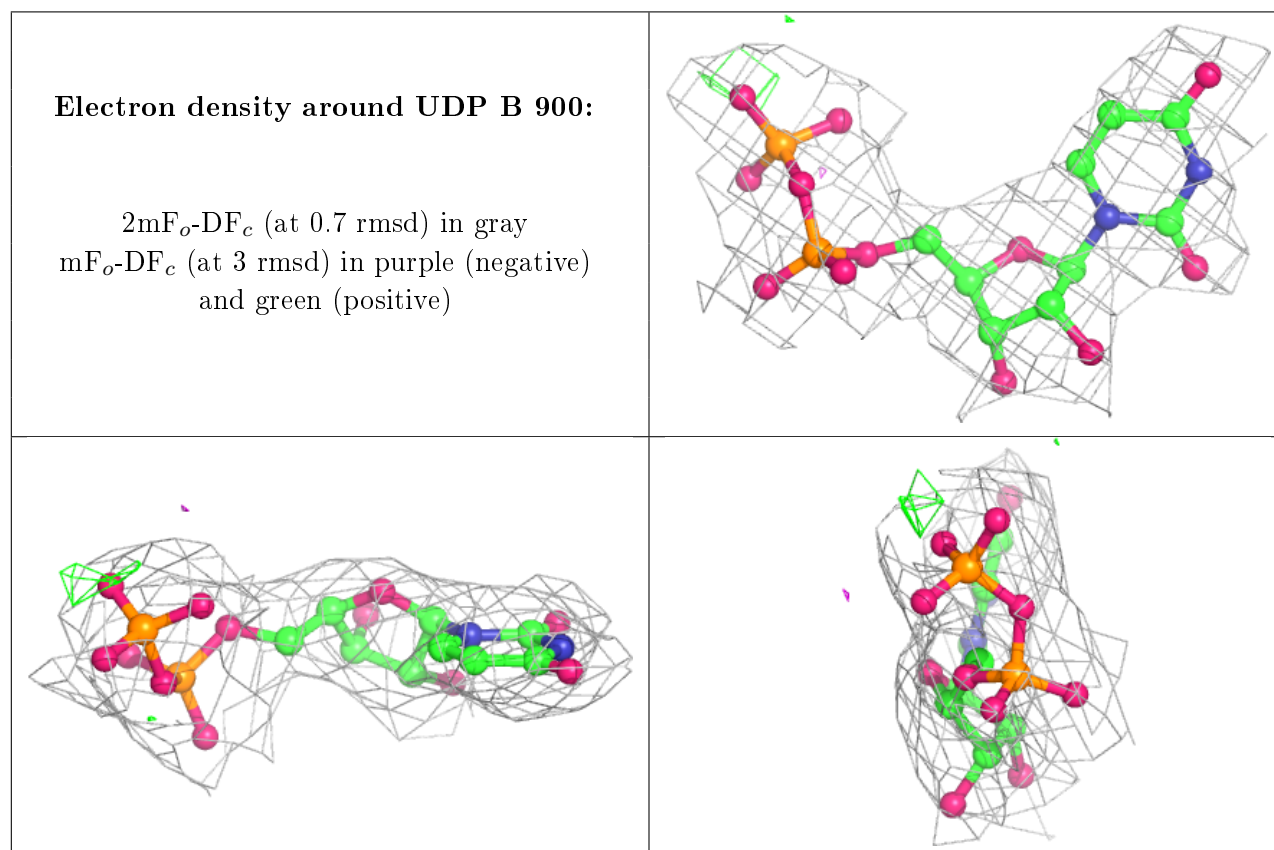
There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<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.

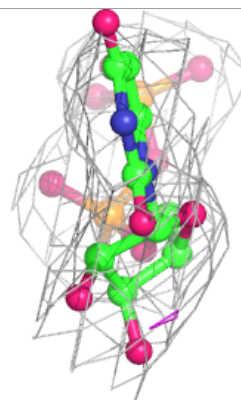
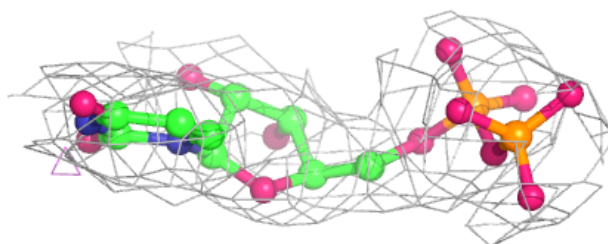
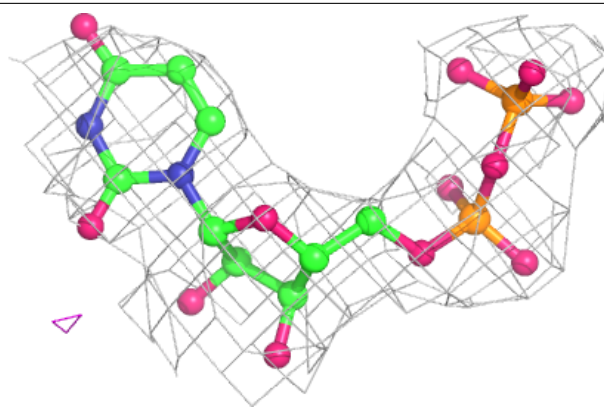
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	UDP	B	900	25/25	0.94	0.20	30,43,53,55	0
2	UDP	D	900	25/25	0.94	0.15	39,51,62,64	0
2	UDP	A	900	25/25	0.96	0.15	18,30,41,43	0
2	UDP	C	900	25/25	0.96	0.15	21,34,44,46	0

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.

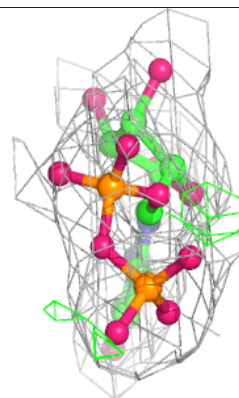
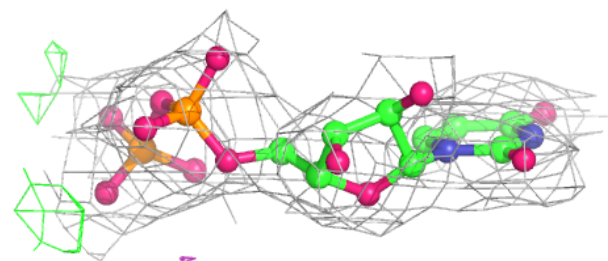
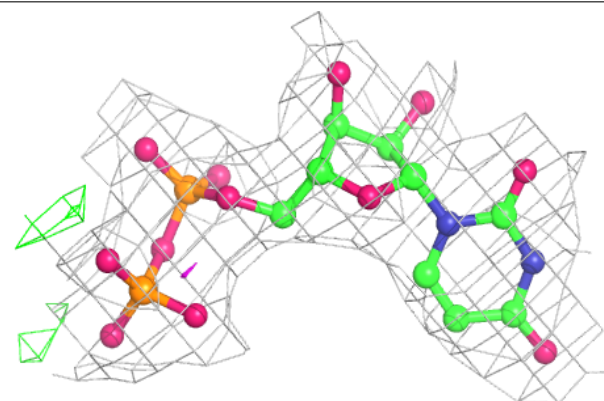


**Electron density around UDP D 900:**

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

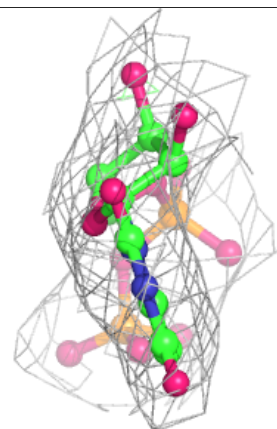
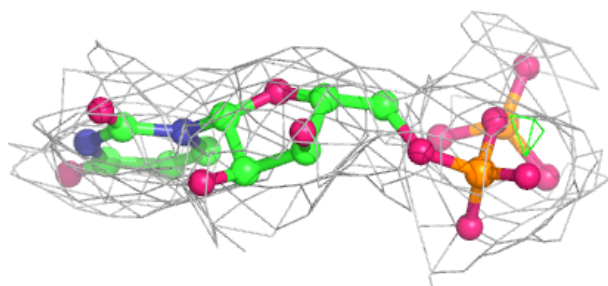
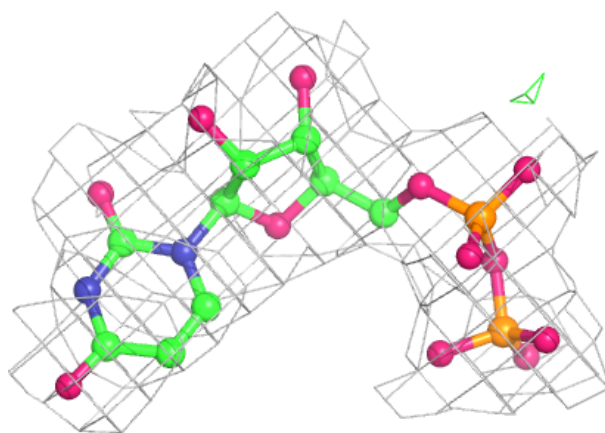
**Electron density around UDP A 900:**

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



**Electron density around UDP C 900:**

$2mF_o - DF_c$  (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.