



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 10:06 pm BST

PDB ID : 3QKW  
Title : Structure of Streptococcus parasanguinis Gtf3 glycosyltransferase  
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Wu, H.  
Deposited on : 2011-02-01  
Resolution : 2.29 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.

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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

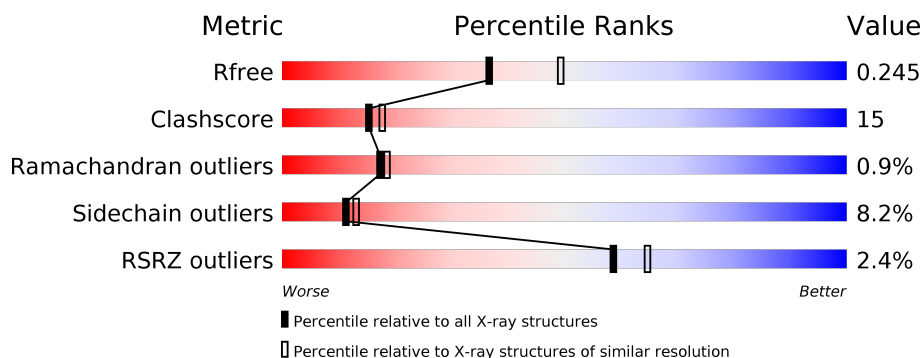
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.29 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	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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	332	<div> <div>2%</div> <div> <div></div> <div>69%</div> <div>25%</div> <div>5%</div> </div> </div>
1	B	332	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>20%</div> <div></div> </div> </div>
1	C	332	<div> <div>2%</div> <div> <div></div> <div>70%</div> <div>23%</div> <div>5%</div> </div> </div>
1	D	332	<div> <div>4%</div> <div> <div></div> <div>68%</div> <div>27%</div> <div></div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10871 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.

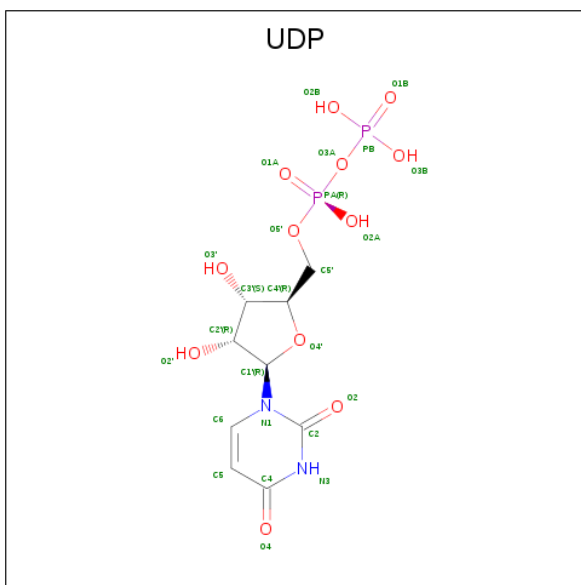
- Molecule 1 is a protein called Nucleotide sugar synthetase-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	0	0
			2657	1708	443	490	16			
1	B	324	Total	C	N	O	S	0	0	0
			2618	1687	437	477	17			
1	C	329	Total	C	N	O	S	0	0	0
			2669	1717	443	492	17			
1	D	326	Total	C	N	O	S	0	0	0
			2639	1697	441	485	16			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ALA	-	EXPRESSION TAG	UNP B5A7L9
A	0	CYS	-	EXPRESSION TAG	UNP B5A7L9
B	-1	ALA	-	EXPRESSION TAG	UNP B5A7L9
B	0	CYS	-	EXPRESSION TAG	UNP B5A7L9
C	-1	ALA	-	EXPRESSION TAG	UNP B5A7L9
C	0	CYS	-	EXPRESSION TAG	UNP B5A7L9
D	-1	ALA	-	EXPRESSION TAG	UNP B5A7L9
D	0	CYS	-	EXPRESSION TAG	UNP B5A7L9

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

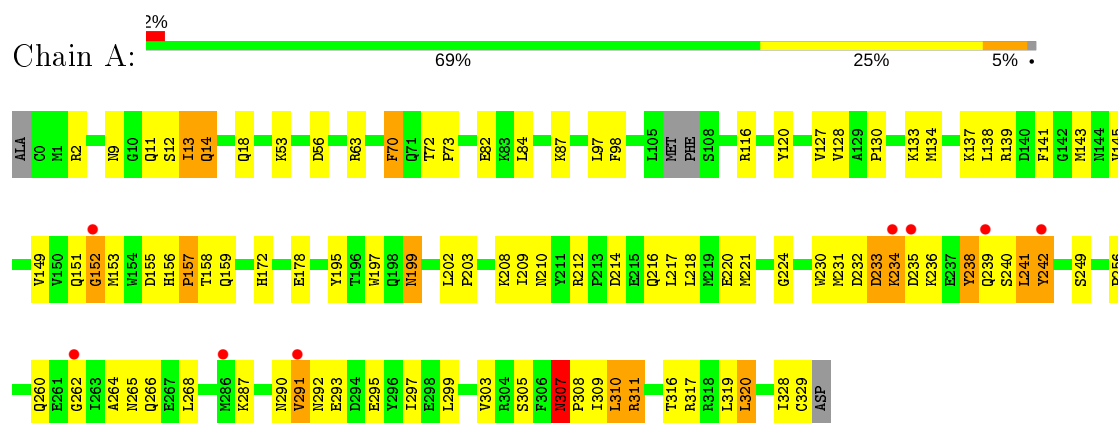
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	52	Total O 52 52	0	0
3	B	57	Total O 57 57	0	0
3	C	47	Total O 47 47	0	0
3	D	32	Total O 32 32	0	0

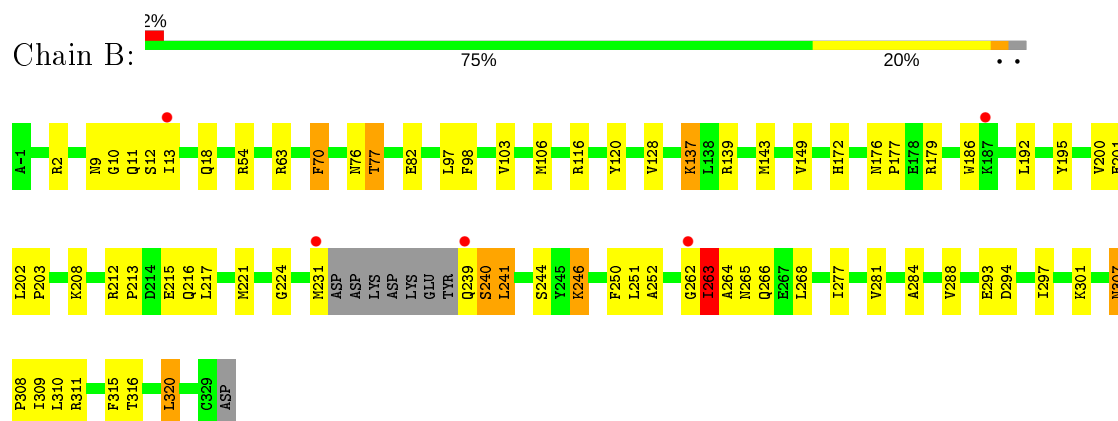
### 3 Residue-property plots

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

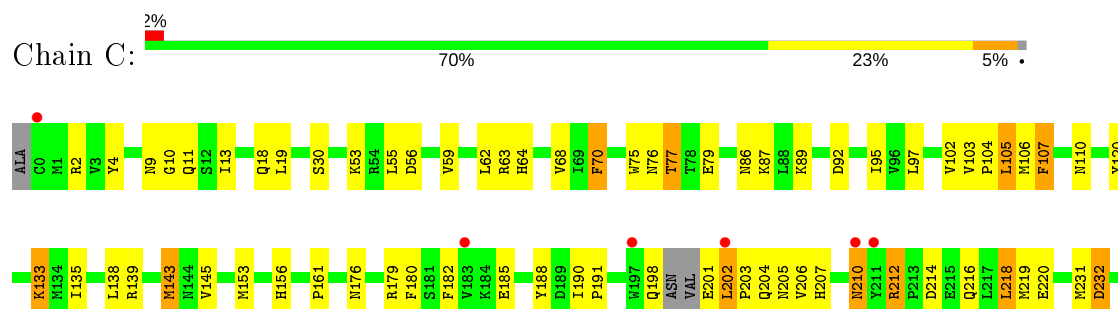
- Molecule 1: Nucleotide sugar synthetase-like protein



- Molecule 1: Nucleotide sugar synthetase-like protein

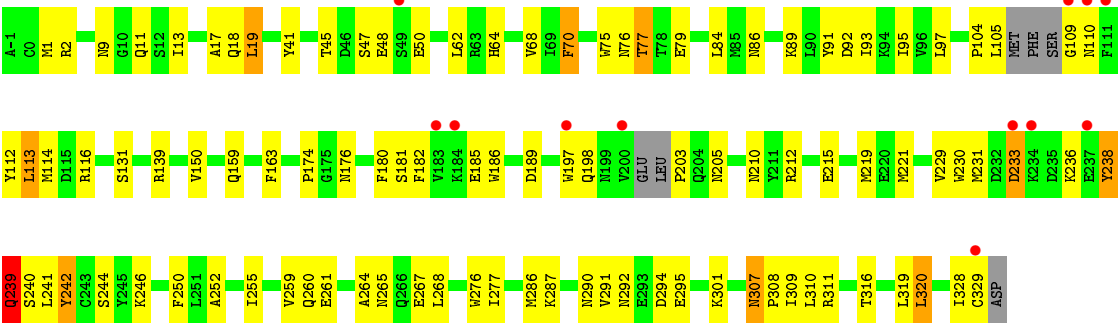


- Molecule 1: Nucleotide sugar synthetase-like protein





● Molecule 1: Nucleotide sugar synthetase-like protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.22Å 99.24Å 188.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.51 – 2.29 42.51 – 2.29	Depositor EDS
% Data completeness (in resolution range)	95.7 (42.51-2.29) 95.7 (42.51-2.29)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.78 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_351)	Depositor
R, $R_{free}$	0.187 , 0.250 0.189 , 0.245	Depositor DCC
$R_{free}$ test set	3218 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.9	Xtriage
Anisotropy	0.591	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 45.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	10871	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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

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	0.42	0/2716	0.59	2/3677 (0.1%)
1	B	0.49	0/2677	0.58	0/3625
1	C	0.40	0/2729	0.55	0/3693
1	D	0.38	0/2697	0.58	1/3649 (0.0%)
All	All	0.42	0/10819	0.58	3/14644 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	233	ASP	N-CA-C	5.68	126.34	111.00
1	A	311	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	A	152	GLY	N-CA-C	5.23	126.18	113.10

There are no chirality outliers.

There are no planarity outliers.

### 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	2657	0	2647	93	0
1	B	2618	0	2617	75	0
1	C	2669	0	2654	76	0
1	D	2639	0	2630	85	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	25	0	11	0	0
2	B	25	0	11	0	0
2	C	25	0	11	0	0
2	D	25	0	11	0	0
3	A	52	0	0	3	0
3	B	57	0	0	1	0
3	C	47	0	0	0	0
3	D	32	0	0	2	0
All	All	10871	0	10592	317	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 15.

All (317) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:ASN:OD1	1:A:295:GLU:HG3	1.24	1.27
1:B:263:ILE:O	1:B:263:ILE:HD13	1.40	1.19
1:C:212:ARG:HB2	1:C:212:ARG:HH11	1.05	1.15
1:A:292:ASN:CG	1:A:295:GLU:HG3	1.75	1.06
1:D:64:HIS:HD2	1:D:92:ASP:H	1.10	0.99
1:D:231:MET:HG3	1:D:239:GLN:HB3	1.47	0.97
1:D:77:THR:HG22	1:D:79:GLU:H	1.26	0.97
1:C:64:HIS:HD2	1:C:92:ASP:H	1.13	0.96
1:C:212:ARG:HB2	1:C:212:ARG:NH1	1.81	0.93
1:B:263:ILE:C	1:B:263:ILE:HD13	1.83	0.91
1:B:11:GLN:H	1:B:18:GLN:HE22	1.16	0.90
1:A:240:SER:O	1:A:264:ALA:HB2	1.73	0.89
1:A:292:ASN:OD1	1:A:295:GLU:CG	2.17	0.87
1:A:317:ARG:HD2	3:A:371:HOH:O	1.75	0.87
1:C:77:THR:HG22	1:C:79:GLU:H	1.40	0.86
1:D:105:LEU:HB3	1:D:238:TYR:HE1	1.41	0.85
1:A:152:GLY:O	1:A:153:MET:HB2	1.79	0.82
1:D:291:VAL:HG23	1:D:295:GLU:HG3	1.61	0.81
1:A:11:GLN:H	1:A:18:GLN:HE22	1.26	0.80
1:B:251:LEU:O	1:B:307:ASN:ND2	2.13	0.79
1:B:297:ILE:O	1:B:301:LYS:HG2	1.83	0.79
1:D:64:HIS:CD2	1:D:92:ASP:H	1.98	0.78
1:D:316:THR:HG22	1:D:320:LEU:HD22	1.65	0.77
1:D:113:LEU:HD12	3:D:356:HOH:O	1.83	0.77
1:D:11:GLN:H	1:D:18:GLN:HE22	1.32	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:GLN:H	1:C:18:GLN:HE22	1.32	0.76
1:A:234:LYS:HB3	1:A:234:LYS:HZ2	1.49	0.75
1:B:262:GLY:H	1:B:266:GLN:HE22	1.32	0.75
1:B:263:ILE:O	1:B:263:ILE:CD1	2.30	0.74
1:B:179:ARG:HH21	1:B:231:MET:HE1	1.53	0.74
1:B:316:THR:HG22	1:B:320:LEU:HD22	1.69	0.74
1:A:234:LYS:H	1:A:234:LYS:HD2	1.51	0.74
1:C:202:LEU:HB2	1:C:203:PRO:HD2	1.70	0.74
1:B:263:ILE:HD11	1:B:266:GLN:HB3	1.69	0.73
1:A:155:ASP:HB3	1:A:157:PRO:HD3	1.69	0.72
1:C:75:TRP:HH2	1:C:106:MET:HE3	1.52	0.72
1:B:263:ILE:C	1:B:263:ILE:CD1	2.54	0.71
1:A:316:THR:HG22	1:A:320:LEU:HD22	1.72	0.71
1:A:56:ASP:OD1	1:A:87:LYS:HE3	1.89	0.71
1:A:230:TRP:CZ3	1:A:260:GLN:HG2	2.25	0.71
1:D:105:LEU:HB3	1:D:238:TYR:CE1	2.25	0.71
1:C:212:ARG:CB	1:C:212:ARG:HH11	1.95	0.71
1:B:202:LEU:HD12	1:B:208:LYS:HB2	1.73	0.70
1:A:156:HIS:HE1	1:A:214:ASP:OD2	1.74	0.70
1:A:216:GLN:O	1:A:220:GLU:HG3	1.91	0.70
1:C:103:VAL:HG23	1:C:106:MET:HE2	1.75	0.69
1:C:103:VAL:HG12	1:C:104:PRO:HD3	1.75	0.68
1:B:179:ARG:NH2	1:B:231:MET:CE	2.56	0.68
1:B:315:PHE:O	3:B:336:HOH:O	2.11	0.68
1:A:231:MET:HG2	1:A:239:GLN:HB2	1.77	0.67
1:C:13:ILE:H	1:C:13:ILE:HD12	1.60	0.67
1:B:76:ASN:O	1:B:77:THR:HB	1.95	0.67
1:C:202:LEU:CB	1:C:203:PRO:HD2	2.26	0.66
1:A:158:THR:HG21	1:A:218:LEU:HD23	1.77	0.66
1:A:236:LYS:O	1:A:239:GLN:HB3	1.96	0.66
1:A:238:TYR:HD2	1:A:238:TYR:O	1.78	0.66
1:D:238:TYR:O	1:D:241:LEU:N	2.24	0.66
1:A:152:GLY:O	1:A:268:LEU:CD1	2.44	0.66
1:A:172:HIS:HD2	1:A:195:TYR:OH	1.77	0.66
1:B:240:SER:CB	1:B:263:ILE:HA	2.26	0.66
1:B:179:ARG:NH2	1:B:231:MET:HE1	2.10	0.66
1:C:307:ASN:C	1:C:307:ASN:HD22	1.99	0.66
1:C:106:MET:HB3	1:C:238:TYR:CE1	2.31	0.65
1:B:240:SER:HB3	1:B:263:ILE:HA	1.79	0.65
1:C:307:ASN:HD22	1:C:308:PRO:HD3	1.61	0.65
1:C:316:THR:HG22	1:C:320:LEU:HD22	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:ILE:CD1	1:B:266:GLN:HB3	2.27	0.64
1:D:307:ASN:HD22	1:D:308:PRO:HD3	1.62	0.64
1:D:70:PHE:CZ	1:D:84:LEU:HD23	2.33	0.64
1:D:17:ALA:HB1	1:D:75:TRP:CZ3	2.32	0.64
1:D:76:ASN:O	1:D:77:THR:HB	1.96	0.64
1:A:230:TRP:CH2	1:A:260:GLN:HG2	2.32	0.64
1:A:11:GLN:HB2	1:A:18:GLN:NE2	2.13	0.64
1:A:290:ASN:O	1:A:291:VAL:C	2.37	0.63
1:B:172:HIS:HD2	1:B:195:TYR:OH	1.80	0.63
1:C:212:ARG:HG2	1:C:216:GLN:CD	2.19	0.63
1:C:210:ASN:O	1:C:212:ARG:HD3	1.98	0.63
1:B:239:GLN:O	1:B:241:LEU:N	2.28	0.62
1:B:263:ILE:N	1:B:263:ILE:HD12	2.14	0.62
1:C:156:HIS:HE1	1:C:214:ASP:OD2	1.81	0.62
1:A:199:ASN:H	1:A:199:ASN:HD22	1.48	0.62
1:A:158:THR:HG21	1:A:218:LEU:CD2	2.30	0.62
1:B:137:LYS:NZ	1:B:137:LYS:HB2	2.15	0.61
1:C:76:ASN:O	1:C:77:THR:HB	2.00	0.61
1:B:307:ASN:CB	1:B:308:PRO:HD3	2.30	0.61
1:C:102:VAL:HG11	1:C:105:LEU:HD12	1.81	0.61
1:A:292:ASN:ND2	1:A:295:GLU:HG3	2.15	0.61
1:C:64:HIS:CD2	1:C:92:ASP:H	2.05	0.61
1:B:76:ASN:O	1:B:77:THR:CB	2.48	0.61
1:C:135:ILE:O	1:C:139:ARG:HG2	2.00	0.61
1:A:156:HIS:N	1:A:157:PRO:HD3	2.16	0.61
1:C:179:ARG:HH21	1:C:231:MET:CE	2.14	0.60
1:C:307:ASN:HD22	1:C:308:PRO:CD	2.14	0.60
1:A:70:PHE:CZ	1:A:84:LEU:HD23	2.37	0.60
1:B:239:GLN:OE1	1:B:239:GLN:HA	2.01	0.60
1:A:307:ASN:HD22	1:A:307:ASN:C	2.04	0.60
1:D:104:PRO:O	1:D:105:LEU:HG	2.01	0.60
1:D:11:GLN:HB2	1:D:18:GLN:NE2	2.16	0.60
1:D:11:GLN:HB2	1:D:18:GLN:HE21	1.66	0.60
1:D:233:ASP:HA	1:D:236:LYS:HG3	1.83	0.59
1:A:307:ASN:HD22	1:A:308:PRO:HD3	1.68	0.59
1:D:197:TRP:CD1	1:D:198:GLN:HG3	2.37	0.59
1:D:186:TRP:HB3	1:D:203:PRO:HG3	1.84	0.58
1:D:292:ASN:OD1	1:D:295:GLU:HG2	2.03	0.58
1:D:221:MET:HE1	1:D:255:ILE:HD11	1.85	0.58
1:A:159:GLN:HB2	3:A:358:HOH:O	2.03	0.58
1:D:221:MET:CE	1:D:255:ILE:HD11	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:236:LYS:HA	1:D:239:GLN:HG2	1.85	0.57
1:D:104:PRO:HG3	1:D:114:MET:HA	1.87	0.57
1:D:64:HIS:HD2	1:D:92:ASP:N	1.91	0.57
1:D:105:LEU:CB	1:D:238:TYR:HE1	2.15	0.57
1:A:133:LYS:HG2	1:A:241:LEU:HD12	1.86	0.57
1:A:262:GLY:H	1:A:266:GLN:HE22	1.53	0.57
1:D:131:SER:HA	1:D:150:VAL:HG13	1.86	0.57
1:B:11:GLN:HB2	1:B:18:GLN:HE21	1.69	0.56
1:C:179:ARG:HH21	1:C:231:MET:HE1	1.71	0.56
1:A:157:PRO:HA	1:A:311:ARG:O	2.05	0.56
1:B:11:GLN:HB2	1:B:18:GLN:NE2	2.21	0.56
1:C:76:ASN:O	1:C:77:THR:CB	2.54	0.56
1:D:89:LYS:HD3	1:D:95:ILE:HD12	1.88	0.56
1:C:176:ASN:HB3	1:C:198:GLN:HE21	1.71	0.56
1:D:229:VAL:HB	1:D:259:VAL:HG12	1.87	0.56
1:D:186:TRP:HB3	1:D:203:PRO:CG	2.36	0.55
1:C:143:MET:HE3	1:C:145:VAL:H	1.71	0.55
1:D:238:TYR:O	1:D:240:SER:N	2.39	0.55
1:A:307:ASN:HD22	1:A:308:PRO:CD	2.20	0.55
1:B:179:ARG:NH2	1:B:231:MET:HE2	2.22	0.55
1:D:197:TRP:NE1	1:D:198:GLN:HG3	2.22	0.55
1:D:316:THR:O	1:D:320:LEU:HB2	2.07	0.55
1:A:265:ASN:ND2	3:A:352:HOH:O	2.39	0.55
1:D:11:GLN:H	1:D:18:GLN:NE2	2.04	0.54
1:D:261:GLU:HA	1:D:277:ILE:CG2	2.37	0.54
1:D:307:ASN:HD22	1:D:308:PRO:CD	2.21	0.54
1:D:231:MET:CG	1:D:239:GLN:HB3	2.31	0.54
1:C:307:ASN:HD22	1:C:308:PRO:N	2.06	0.54
1:A:11:GLN:H	1:A:18:GLN:NE2	2.00	0.54
1:B:11:GLN:H	1:B:18:GLN:NE2	1.97	0.54
1:A:172:HIS:HE1	1:A:224:GLY:O	1.91	0.54
1:D:112:TYR:N	3:D:356:HOH:O	2.41	0.54
1:D:286:MET:HG2	1:D:290:ASN:HD21	1.73	0.54
1:D:77:THR:HG22	1:D:79:GLU:N	2.09	0.53
1:B:294:ASP:HA	1:B:297:ILE:HD12	1.90	0.53
1:A:156:HIS:N	1:A:157:PRO:CD	2.69	0.53
1:B:11:GLN:N	1:B:18:GLN:HE22	1.95	0.53
1:B:200:VAL:HG12	1:B:201:GLU:O	2.08	0.53
1:B:103:VAL:O	1:B:106:MET:HG3	2.08	0.53
1:C:176:ASN:HB3	1:C:198:GLN:NE2	2.24	0.53
1:A:127:VAL:HG23	1:A:145:VAL:HG11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LYS:NZ	1:A:234:LYS:HB3	2.22	0.52
1:B:9:ASN:ND2	1:B:10:GLY:H	2.06	0.52
1:D:238:TYR:O	1:D:239:GLN:C	2.48	0.52
1:A:139:ARG:HG3	1:A:139:ARG:HH11	1.74	0.52
1:D:91:TYR:HB3	1:D:93:ILE:HG23	1.90	0.52
1:A:234:LYS:H	1:A:234:LYS:CD	2.20	0.52
1:B:263:ILE:CD1	1:B:263:ILE:N	2.70	0.52
1:A:137:LYS:HG3	1:A:141:PHE:HE2	1.75	0.52
1:B:293:GLU:O	1:B:297:ILE:HD12	2.09	0.52
1:A:9:ASN:ND2	1:B:63:ARG:HE	2.08	0.51
1:B:284:ALA:O	1:B:288:VAL:HG23	2.11	0.51
1:D:221:MET:HE3	1:D:250:PHE:CE1	2.46	0.51
1:C:107:PHE:HB3	1:C:110:ASN:HD22	1.76	0.51
1:A:137:LYS:HE2	1:A:141:PHE:CZ	2.46	0.51
1:A:133:LYS:HG2	1:A:241:LEU:CD1	2.41	0.50
1:A:172:HIS:HB3	1:A:221:MET:HE1	1.93	0.50
1:C:307:ASN:ND2	1:C:308:PRO:HD3	2.25	0.50
1:D:181:SER:O	1:D:185:GLU:HG2	2.11	0.50
1:D:76:ASN:O	1:D:77:THR:CB	2.58	0.50
1:B:213:PRO:HD2	1:B:216:GLN:HE21	1.75	0.50
1:B:70:PHE:CD1	1:B:70:PHE:C	2.85	0.50
1:C:307:ASN:ND2	1:C:307:ASN:C	2.63	0.50
1:A:217:LEU:HD11	1:A:221:MET:CE	2.42	0.50
1:A:290:ASN:O	1:A:291:VAL:O	2.30	0.50
1:A:293:GLU:O	1:A:297:ILE:HG12	2.12	0.50
1:C:89:LYS:HD3	1:C:95:ILE:HD12	1.94	0.50
1:A:13:ILE:O	1:A:13:ILE:HD13	2.11	0.50
1:A:156:HIS:O	1:A:157:PRO:C	2.51	0.49
1:A:291:VAL:O	1:A:291:VAL:HG13	2.12	0.49
1:A:11:GLN:HB2	1:A:18:GLN:HE21	1.77	0.49
1:B:217:LEU:O	1:B:221:MET:HG3	2.12	0.49
1:A:53:LYS:HE3	1:B:54:ARG:NH2	2.27	0.49
1:C:107:PHE:HB3	1:C:110:ASN:ND2	2.28	0.49
1:A:233:ASP:O	1:A:236:LYS:HB2	2.12	0.49
1:A:238:TYR:O	1:A:238:TYR:CD2	2.62	0.49
1:A:307:ASN:ND2	1:A:307:ASN:C	2.66	0.49
1:B:9:ASN:ND2	1:B:10:GLY:N	2.61	0.49
1:C:75:TRP:CH2	1:C:106:MET:HE3	2.40	0.49
1:C:204:GLN:HG3	1:C:205:ASN:N	2.28	0.49
1:C:53:LYS:HD2	1:D:50:GLU:HB3	1.95	0.49
1:A:53:LYS:HE3	1:B:54:ARG:HH21	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ARG:HD3	1:A:143:MET:O	2.12	0.48
1:D:180:PHE:HB3	1:D:182:PHE:CE1	2.48	0.48
1:A:82:GLU:OE2	1:A:116:ARG:HB3	2.13	0.48
1:C:240:SER:OG	1:C:241:LEU:HD13	2.14	0.48
1:D:252:ALA:O	1:D:311:ARG:NH1	2.41	0.48
1:D:319:LEU:HD12	1:D:319:LEU:C	2.33	0.48
1:A:292:ASN:CG	1:A:295:GLU:CG	2.66	0.48
1:B:213:PRO:HD2	1:B:216:GLN:NE2	2.28	0.48
1:B:186:TRP:CE3	1:B:192:LEU:HD22	2.49	0.48
1:D:264:ALA:O	1:D:265:ASN:HB2	2.13	0.48
1:C:212:ARG:NH2	1:C:220:GLU:OE1	2.47	0.48
1:A:152:GLY:O	1:A:268:LEU:HD11	2.14	0.47
1:C:214:ASP:OD2	1:C:218:LEU:HD22	2.15	0.47
1:A:202:LEU:HB3	1:A:203:PRO:HD2	1.96	0.47
1:C:70:PHE:C	1:C:70:PHE:CD1	2.87	0.47
1:B:246:LYS:HG2	1:B:250:PHE:CE2	2.50	0.47
1:C:9:ASN:ND2	1:C:10:GLY:H	2.12	0.47
1:C:133:LYS:HD3	1:C:133:LYS:N	2.28	0.47
1:D:176:ASN:HB3	1:D:198:GLN:OE1	2.15	0.47
1:D:242:TYR:CD2	1:D:242:TYR:C	2.87	0.47
1:D:47:SER:OG	1:D:50:GLU:HG3	2.15	0.47
1:A:72:THR:OG1	1:A:73:PRO:HA	2.15	0.46
1:D:1:MET:SD	1:D:328:ILE:HD11	2.55	0.46
1:A:307:ASN:HD22	1:A:308:PRO:N	2.13	0.46
1:A:307:ASN:N	1:A:308:PRO:CD	2.77	0.46
1:C:4:TYR:HB2	1:C:68:VAL:HG22	1.96	0.46
1:B:179:ARG:HH21	1:B:231:MET:CE	2.21	0.46
1:C:216:GLN:O	1:C:220:GLU:HG3	2.15	0.46
1:B:212:ARG:HB2	1:B:216:GLN:NE2	2.31	0.46
1:D:109:GLY:HA2	1:D:110:ASN:HA	1.78	0.46
1:B:262:GLY:O	1:B:263:ILE:O	2.34	0.46
1:A:156:HIS:CD2	1:A:249:SER:OG	2.69	0.46
1:C:207:HIS:N	1:C:207:HIS:CD2	2.83	0.46
1:C:242:TYR:CD1	1:C:242:TYR:C	2.89	0.46
1:C:283:GLU:O	1:C:287:LYS:HG3	2.16	0.46
1:C:232:ASP:N	1:C:232:ASP:OD1	2.46	0.45
1:B:307:ASN:CB	1:B:308:PRO:CD	2.94	0.45
1:C:212:ARG:HG2	1:C:216:GLN:NE2	2.32	0.45
1:D:113:LEU:HD23	1:D:116:ARG:HH11	1.81	0.45
1:B:13:ILE:N	1:B:13:ILE:HD13	2.31	0.45
1:C:56:ASP:OD1	1:C:87:LYS:NZ	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:PRO:HA	1:A:151:GLN:HE21	1.82	0.45
1:A:328:ILE:HG21	1:D:309:ILE:HG13	1.98	0.45
1:C:237:GLU:OE1	1:C:237:GLU:HA	2.16	0.45
1:D:189:ASP:HA	1:D:205:ASN:ND2	2.32	0.45
1:C:70:PHE:C	1:C:70:PHE:HD1	2.20	0.45
1:A:11:GLN:N	1:A:18:GLN:HE22	2.04	0.45
1:D:89:LYS:HA	1:D:89:LYS:HD2	1.76	0.45
1:B:139:ARG:HG3	1:B:139:ARG:HH11	1.81	0.45
1:D:112:TYR:CD2	1:D:113:LEU:HG	2.52	0.45
1:A:234:LYS:N	1:A:234:LYS:HZ3	2.13	0.45
1:A:242:TYR:C	1:A:242:TYR:CD1	2.88	0.45
1:A:195:TYR:HA	1:A:209:ILE:O	2.17	0.44
1:A:217:LEU:HD11	1:A:221:MET:HE3	1.99	0.44
1:B:82:GLU:OE2	1:B:116:ARG:HB3	2.16	0.44
1:B:262:GLY:H	1:B:266:GLN:NE2	2.06	0.44
1:C:307:ASN:N	1:C:308:PRO:CD	2.80	0.44
1:D:307:ASN:HD22	1:D:307:ASN:C	2.19	0.44
1:D:319:LEU:HD12	1:D:320:LEU:N	2.32	0.44
1:A:292:ASN:ND2	1:A:295:GLU:CG	2.80	0.44
1:B:307:ASN:HB2	1:B:308:PRO:HD3	1.99	0.44
1:A:12:SER:C	1:A:14:GLN:H	2.21	0.44
1:A:236:LYS:NZ	1:A:260:GLN:HE22	2.16	0.44
1:B:139:ARG:HD3	1:B:143:MET:O	2.17	0.44
1:C:103:VAL:CG1	1:C:104:PRO:HD3	2.45	0.44
1:C:161:PRO:HD2	1:C:219:MET:SD	2.58	0.44
1:D:19:LEU:HD22	1:D:19:LEU:HA	1.65	0.44
1:B:216:GLN:HG3	1:D:219:MET:HE1	1.99	0.44
1:D:45:THR:O	1:D:45:THR:HG22	2.16	0.44
1:D:113:LEU:HD23	1:D:116:ARG:NH1	2.33	0.44
1:A:63:ARG:HE	1:B:9:ASN:ND2	2.16	0.44
1:C:143:MET:HE3	1:C:145:VAL:N	2.32	0.44
1:C:202:LEU:CB	1:C:203:PRO:CD	2.95	0.44
1:D:174:PRO:HD2	1:D:246:LYS:HD3	2.00	0.44
1:A:299:LEU:O	1:A:303:VAL:HG23	2.18	0.43
1:A:9:ASN:HD21	1:B:63:ARG:HE	1.66	0.43
1:C:62:LEU:HD23	1:C:62:LEU:HA	1.88	0.43
1:A:256:PRO:HB3	1:A:299:LEU:HB3	2.00	0.43
1:C:294:ASP:O	1:C:297:ILE:HB	2.19	0.43
1:C:153:MET:CE	1:C:268:LEU:HD23	2.49	0.43
1:C:55:LEU:O	1:C:59:VAL:HG22	2.18	0.43
1:C:89:LYS:HA	1:C:89:LYS:HD2	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:218:LEU:HD12	1:C:253:ALA:HB2	2.00	0.43
1:D:13:ILE:H	1:D:13:ILE:HG13	1.50	0.43
1:B:262:GLY:O	1:B:263:ILE:C	2.57	0.43
1:A:287:LYS:HE3	1:A:287:LYS:HB3	1.86	0.43
1:A:199:ASN:HD22	1:A:199:ASN:N	2.15	0.43
1:C:300:VAL:O	1:C:304:ARG:HG3	2.19	0.42
1:C:77:THR:HG22	1:C:79:GLU:N	2.20	0.42
1:B:172:HIS:CD2	1:B:195:TYR:OH	2.67	0.42
1:B:264:ALA:O	1:B:265:ASN:HB2	2.18	0.42
1:B:137:LYS:HZ3	1:B:137:LYS:HB2	1.84	0.42
1:B:215:GLU:OE1	1:D:215:GLU:OE1	2.37	0.42
1:D:139:ARG:HH11	1:D:139:ARG:HG2	1.84	0.42
1:C:262:GLY:H	1:C:266:GLN:HE22	1.66	0.42
1:A:319:LEU:C	1:A:319:LEU:HD12	2.39	0.42
1:A:156:HIS:CE1	1:A:214:ASP:OD2	2.63	0.42
1:D:230:TRP:CH2	1:D:260:GLN:HG3	2.55	0.41
1:A:128:VAL:HG22	1:A:149:VAL:HB	2.02	0.41
1:A:309:ILE:HG22	1:A:310:LEU:N	2.35	0.41
1:B:70:PHE:HD1	1:B:70:PHE:C	2.22	0.41
1:A:12:SER:C	1:A:14:GLN:N	2.74	0.41
1:D:104:PRO:O	1:D:105:LEU:CG	2.68	0.41
1:B:172:HIS:HE1	1:B:224:GLY:O	2.03	0.41
1:C:11:GLN:H	1:C:18:GLN:NE2	2.09	0.41
1:B:309:ILE:HD11	1:C:329:CYS:HB2	2.03	0.41
1:D:104:PRO:HA	1:D:110:ASN:O	2.21	0.41
1:B:128:VAL:HG22	1:B:149:VAL:HB	2.02	0.41
1:B:277:ILE:HD12	1:B:277:ILE:N	2.35	0.41
1:D:62:LEU:HD21	1:D:68:VAL:HG21	2.01	0.41
1:C:63:ARG:HE	1:D:9:ASN:ND2	2.18	0.41
1:A:14:GLN:HG2	1:A:197:TRP:CZ3	2.56	0.41
1:A:230:TRP:CE3	1:A:260:GLN:HG2	2.55	0.41
1:D:197:TRP:CD1	1:D:198:GLN:N	2.89	0.41
1:D:215:GLU:O	1:D:219:MET:HG3	2.21	0.41
1:D:221:MET:HE1	1:D:250:PHE:HD1	1.85	0.41
1:B:176:ASN:HA	1:B:177:PRO:HD3	1.77	0.41
1:C:180:PHE:HB3	1:C:182:PHE:CE1	2.56	0.41
1:D:276:TRP:CD2	1:D:287:LYS:HD3	2.56	0.41
1:D:291:VAL:CG2	1:D:295:GLU:HG3	2.43	0.41
1:C:190:ILE:HA	1:C:191:PRO:HD3	1.90	0.40
1:D:221:MET:CE	1:D:250:PHE:CD1	3.04	0.40
1:B:9:ASN:HD22	1:B:10:GLY:H	1.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:206:VAL:C	1:C:207:HIS:HD2	2.24	0.40
1:D:238:TYR:C	1:D:240:SER:N	2.74	0.40
1:C:86:ASN:HA	1:C:86:ASN:HD22	1.64	0.40
1:B:252:ALA:O	1:B:311:ARG:NH1	2.47	0.40
1:B:216:GLN:NE2	1:D:163:PHE:HZ	2.20	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	324/332 (98%)	308 (95%)	12 (4%)	4 (1%)	13	12
1	B	320/332 (96%)	308 (96%)	8 (2%)	4 (1%)	12	11
1	C	325/332 (98%)	316 (97%)	7 (2%)	2 (1%)	25	29
1	D	320/332 (96%)	304 (95%)	14 (4%)	2 (1%)	25	29
All	All	1289/1328 (97%)	1236 (96%)	41 (3%)	12 (1%)	17	18

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	291	VAL
1	B	77	THR
1	A	232	ASP
1	B	240	SER
1	B	263	ILE
1	C	77	THR
1	D	77	THR
1	A	157	PRO
1	D	239	GLN
1	C	307	ASN

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Mol	Chain	Res	Type
1	B	203	PRO
1	A	307	ASN

### 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	295/298 (99%)	270 (92%)	25 (8%)	10	12
1	B	290/298 (97%)	274 (94%)	16 (6%)	21	27
1	C	296/298 (99%)	264 (89%)	32 (11%)	6	6
1	D	292/298 (98%)	269 (92%)	23 (8%)	12	14
All	All	1173/1192 (98%)	1077 (92%)	96 (8%)	11	13

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	13	ILE
1	A	14	GLN
1	A	70	PHE
1	A	97	LEU
1	A	98	PHE
1	A	120	TYR
1	A	134	MET
1	A	138	LEU
1	A	178	GLU
1	A	199	ASN
1	A	208	LYS
1	A	210	ASN
1	A	212	ARG
1	A	233	ASP
1	A	234	LYS
1	A	235	ASP
1	A	238	TYR
1	A	241	LEU

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Mol	Chain	Res	Type
1	A	242	TYR
1	A	305	SER
1	A	307	ASN
1	A	310	LEU
1	A	320	LEU
1	A	329	CYS
1	B	2	ARG
1	B	12	SER
1	B	70	PHE
1	B	97	LEU
1	B	98	PHE
1	B	120	TYR
1	B	137	LYS
1	B	241	LEU
1	B	244	SER
1	B	246	LYS
1	B	263	ILE
1	B	268	LEU
1	B	281	VAL
1	B	307	ASN
1	B	310	LEU
1	B	320	LEU
1	C	2	ARG
1	C	19	LEU
1	C	30	SER
1	C	70	PHE
1	C	97	LEU
1	C	105	LEU
1	C	107	PHE
1	C	120	TYR
1	C	133	LYS
1	C	138	LEU
1	C	143	MET
1	C	185	GLU
1	C	188	TYR
1	C	201	GLU
1	C	202	LEU
1	C	210	ASN
1	C	212	ARG
1	C	218	LEU
1	C	232	ASP
1	C	239	GLN

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Mol	Chain	Res	Type
1	C	241	LEU
1	C	242	TYR
1	C	244	SER
1	C	246	LYS
1	C	266	GLN
1	C	268	LEU
1	C	270	GLU
1	C	294	ASP
1	C	301	LYS
1	C	307	ASN
1	C	310	LEU
1	C	320	LEU
1	D	2	ARG
1	D	19	LEU
1	D	41	TYR
1	D	48	GLU
1	D	70	PHE
1	D	86	ASN
1	D	97	LEU
1	D	113	LEU
1	D	159	GLN
1	D	210	ASN
1	D	212	ARG
1	D	238	TYR
1	D	239	GLN
1	D	242	TYR
1	D	244	SER
1	D	267	GLU
1	D	268	LEU
1	D	294	ASP
1	D	301	LYS
1	D	307	ASN
1	D	310	LEU
1	D	320	LEU
1	D	329	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	18	GLN
1	A	22	ASN

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Mol	Chain	Res	Type
1	A	86	ASN
1	A	100	HIS
1	A	151	GLN
1	A	156	HIS
1	A	172	HIS
1	A	199	ASN
1	A	207	HIS
1	A	210	ASN
1	A	216	GLN
1	A	260	GLN
1	A	266	GLN
1	A	307	ASN
1	A	326	GLN
1	B	9	ASN
1	B	18	GLN
1	B	86	ASN
1	B	110	ASN
1	B	151	GLN
1	B	172	HIS
1	B	210	ASN
1	B	216	GLN
1	B	266	GLN
1	B	326	GLN
1	C	9	ASN
1	C	18	GLN
1	C	64	HIS
1	C	86	ASN
1	C	110	ASN
1	C	151	GLN
1	C	156	HIS
1	C	207	HIS
1	C	210	ASN
1	C	216	GLN
1	C	223	GLN
1	C	239	GLN
1	C	266	GLN
1	C	307	ASN
1	C	326	GLN
1	D	9	ASN
1	D	18	GLN
1	D	64	HIS
1	D	86	ASN

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Mol	Chain	Res	Type
1	D	151	GLN
1	D	239	GLN
1	D	266	GLN
1	D	290	ASN
1	D	307	ASN
1	D	326	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	B	400	-	20,26,26	1.22	1 (5%)	25,40,40	0.89	1 (4%)
2	UDP	C	400	-	20,26,26	1.15	1 (5%)	25,40,40	1.08	3 (12%)
2	UDP	A	400	-	20,26,26	1.16	1 (5%)	25,40,40	0.98	2 (8%)
2	UDP	D	400	-	20,26,26	1.14	1 (5%)	25,40,40	1.10	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	B	400	-	-	1/14/32/32	0/2/2/2
2	UDP	C	400	-	-	3/14/32/32	0/2/2/2
2	UDP	A	400	-	-	0/14/32/32	0/2/2/2
2	UDP	D	400	-	-	2/14/32/32	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	400	UDP	C4-N3	3.39	1.38	1.33
2	D	400	UDP	C4-N3	3.25	1.38	1.33
2	C	400	UDP	C4-N3	3.17	1.38	1.33
2	A	400	UDP	C4-N3	2.72	1.37	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	400	UDP	PA-O3A-PB	-2.54	124.12	132.83
2	C	400	UDP	PA-O3A-PB	-2.49	124.29	132.83
2	A	400	UDP	C5-C4-N3	-2.34	118.17	123.31
2	B	400	UDP	O2B-PB-O3A	2.32	112.42	104.64
2	C	400	UDP	C3'-C2'-C1'	2.19	104.28	100.98
2	A	400	UDP	PA-O3A-PB	-2.14	125.48	132.83
2	C	400	UDP	O3B-PB-O3A	2.05	111.51	104.64
2	D	400	UDP	O3B-PB-O3A	2.01	111.36	104.64

There are no chirality outliers.

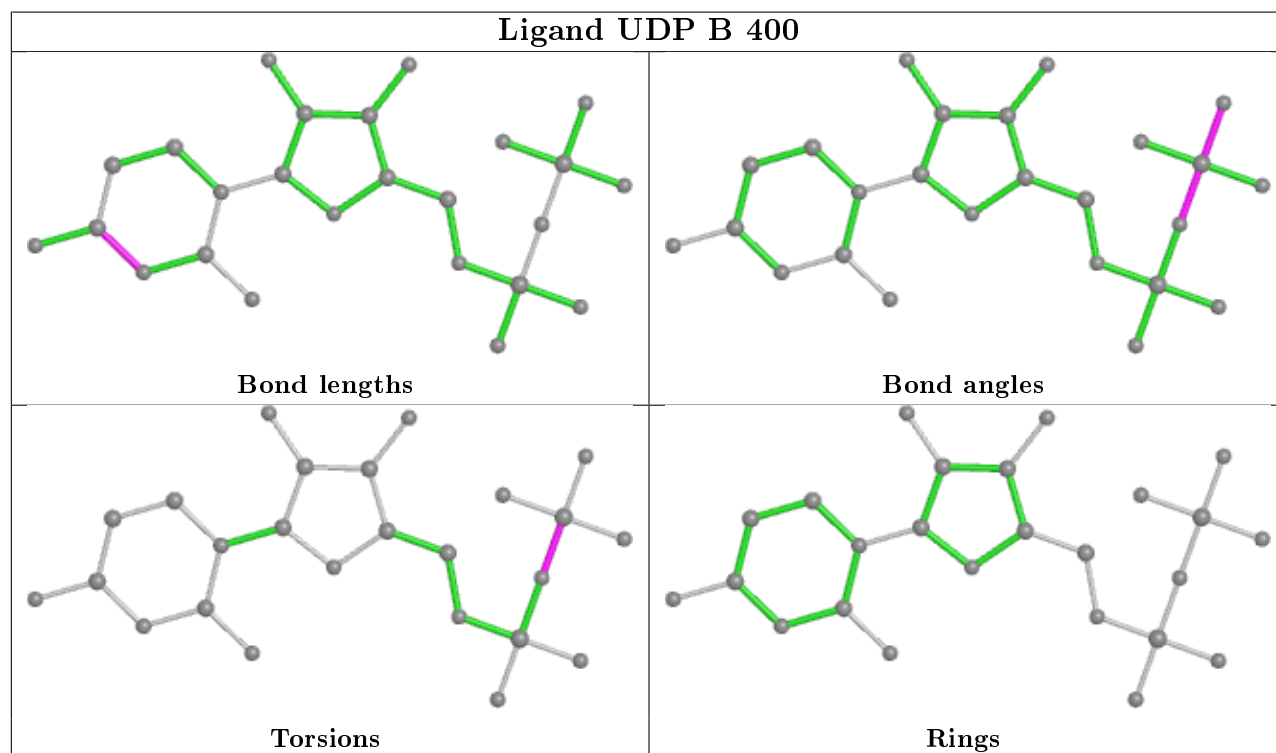
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	400	UDP	PA-O3A-PB-O2B
2	D	400	UDP	PA-O3A-PB-O3B
2	B	400	UDP	PA-O3A-PB-O1B
2	C	400	UDP	PA-O3A-PB-O3B
2	D	400	UDP	C5'-O5'-PA-O1A
2	C	400	UDP	PA-O3A-PB-O1B

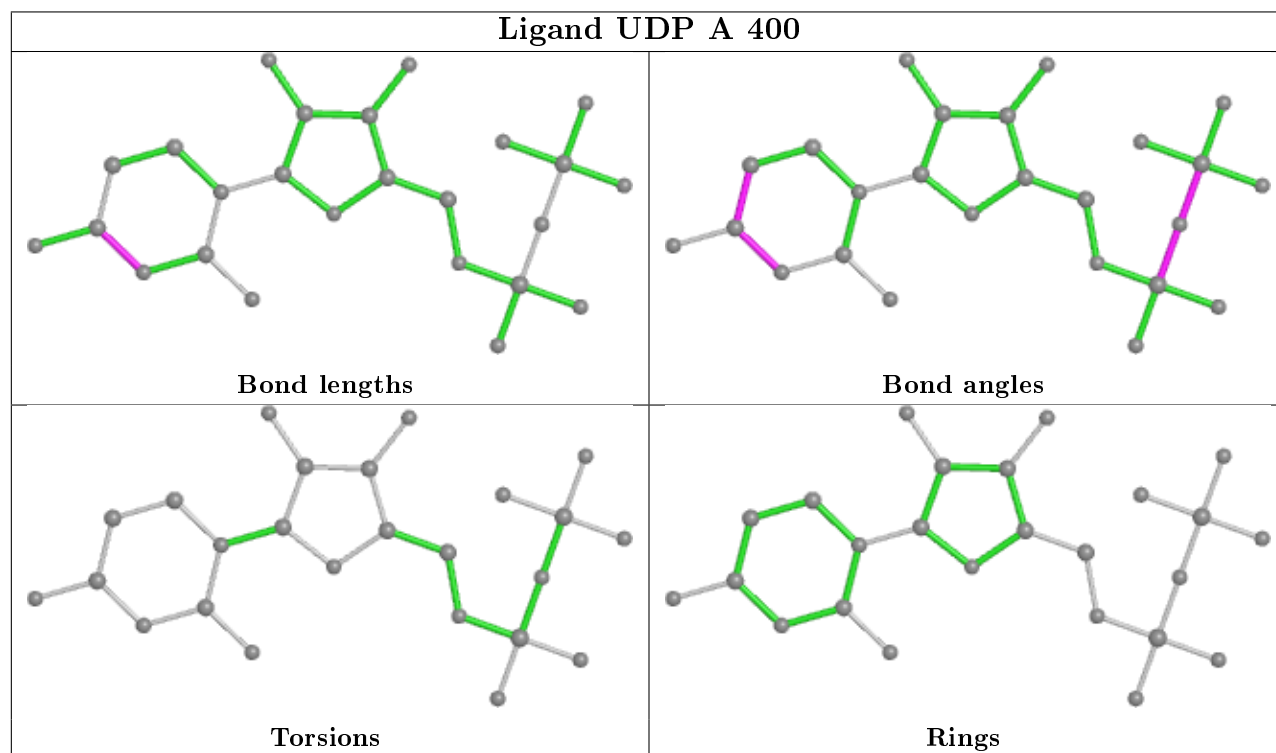
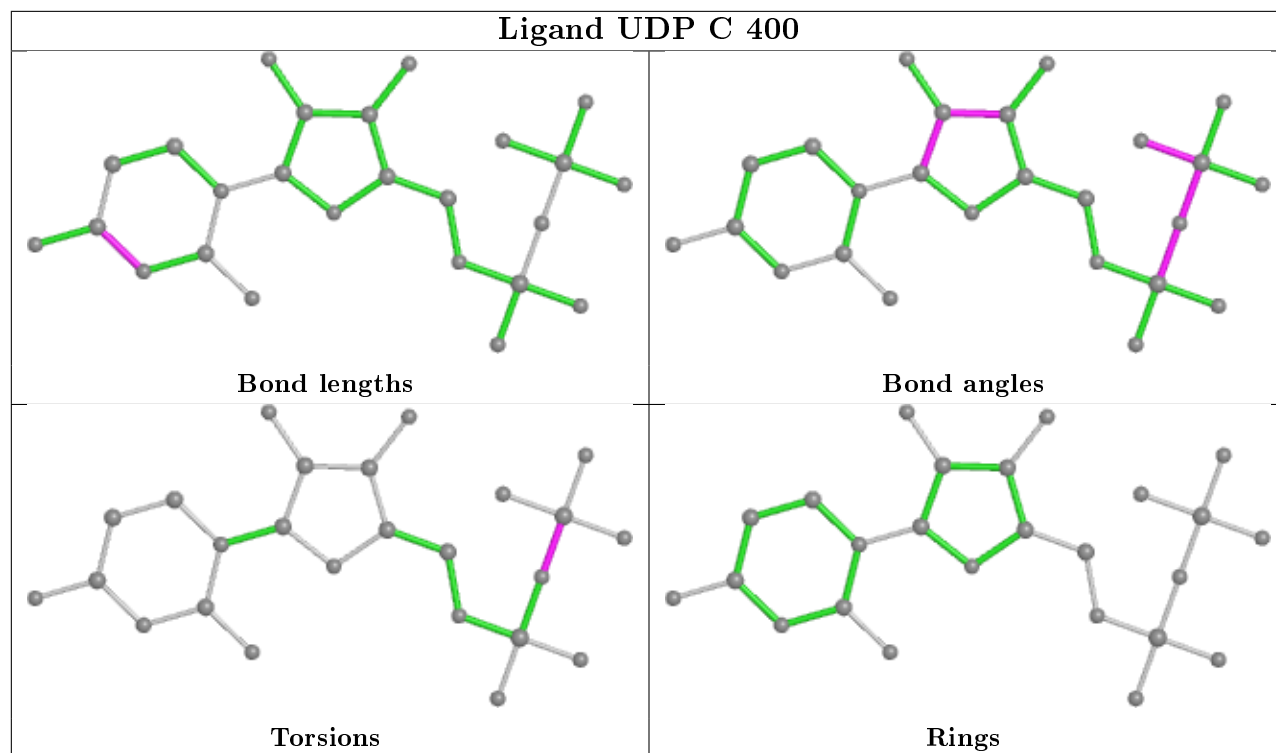
There are no ring outliers.

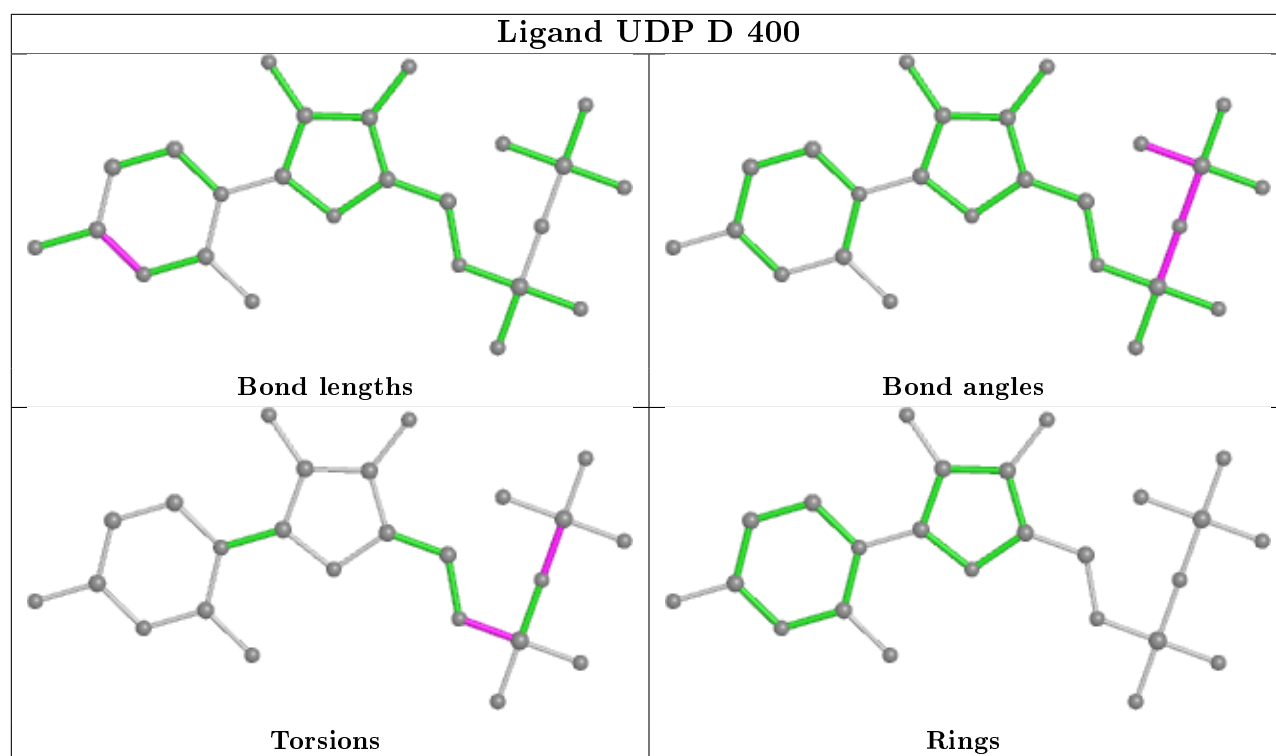
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	328/332 (98%)	-0.13	8 (2%) 59 65	27, 43, 82, 124	0
1	B	324/332 (97%)	-0.27	5 (1%) 73 78	20, 43, 73, 102	0
1	C	329/332 (99%)	-0.09	6 (1%) 68 74	28, 50, 91, 109	0
1	D	326/332 (98%)	0.05	12 (3%) 41 47	31, 53, 92, 122	0
All	All	1307/1328 (98%)	-0.11	31 (2%) 59 65	20, 47, 86, 124	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	200	VAL	6.0
1	A	239	GLN	4.2
1	C	197	TRP	3.3
1	D	183	VAL	3.2
1	A	262	GLY	3.2
1	D	329	CYS	3.1
1	A	234	LYS	3.0
1	C	202	LEU	2.9
1	D	234	LYS	2.8
1	A	152	GLY	2.7
1	B	231	MET	2.7
1	B	187	LYS	2.7
1	B	262	GLY	2.6
1	D	197	TRP	2.5
1	A	242	TYR	2.4
1	B	239	GLN	2.3
1	D	111	PHE	2.3
1	C	0	CYS	2.2
1	C	183	VAL	2.2
1	A	235	ASP	2.2
1	A	286	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	237	GLU	2.2
1	D	110	ASN	2.1
1	A	291	VAL	2.1
1	C	210	ASN	2.1
1	D	184	LYS	2.1
1	D	49	SER	2.1
1	D	109	GLY	2.0
1	C	211	TYR	2.0
1	B	13	ILE	2.0
1	D	233	ASP	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](#)

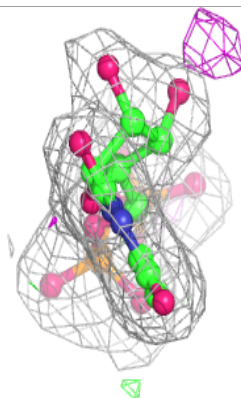
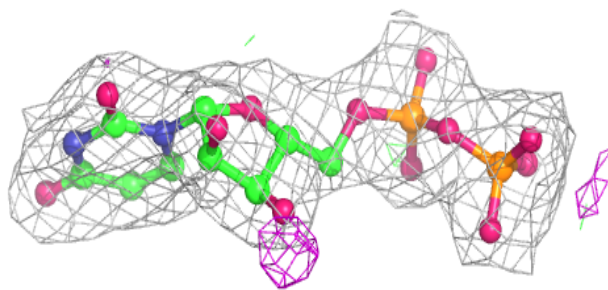
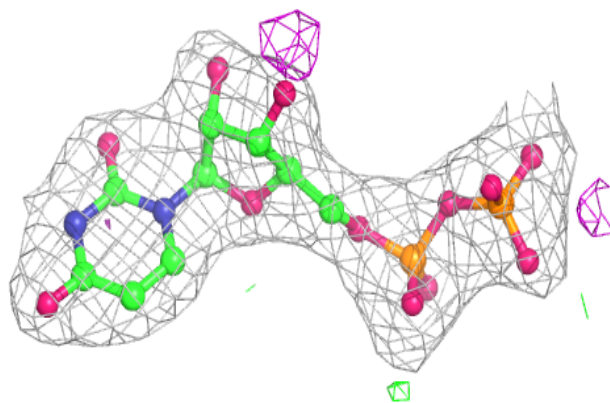
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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	400	25/25	0.98	0.10	25,38,49,51	0
2	UDP	C	400	25/25	0.98	0.11	39,51,63,66	0
2	UDP	A	400	25/25	0.98	0.11	29,38,51,56	0
2	UDP	D	400	25/25	0.98	0.11	43,50,57,60	0

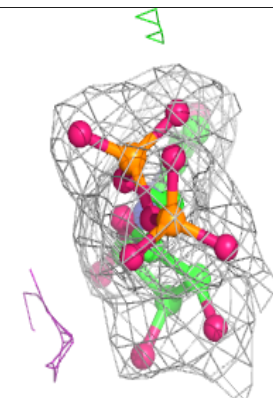
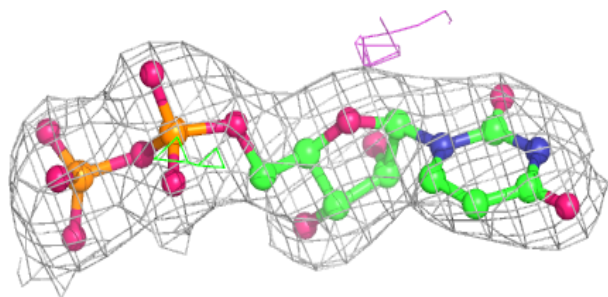
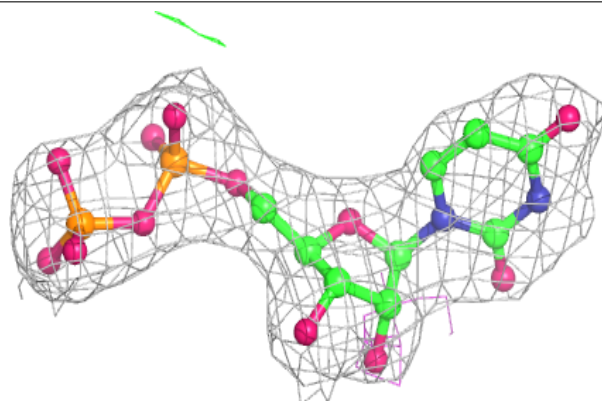
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around UDP B 400:**

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

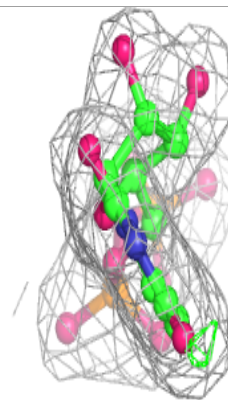
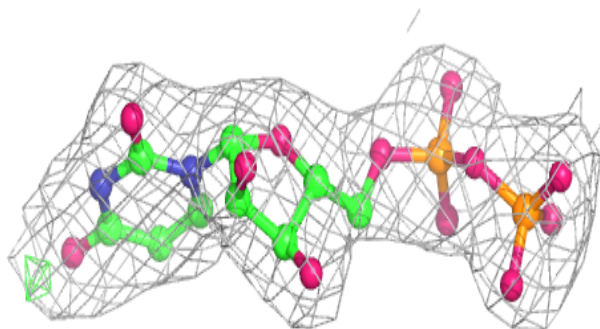
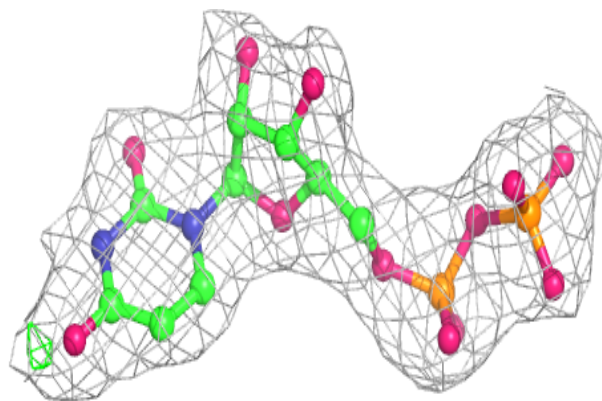
**Electron density around UDP C 400:**

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

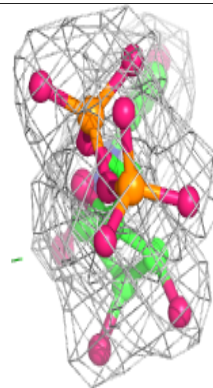
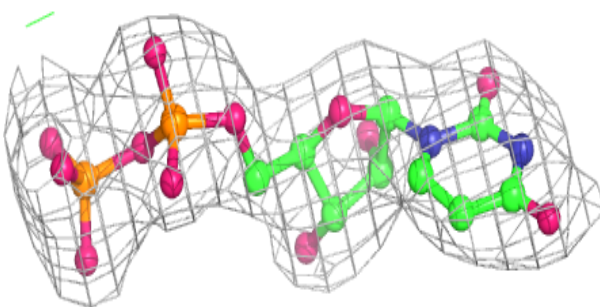
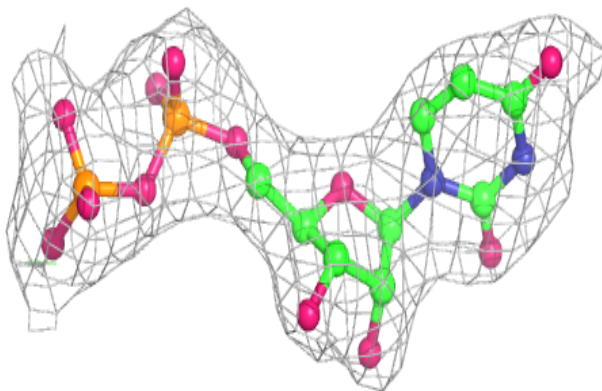


**Electron density around UDP A 400:**

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

**Electron density around UDP D 400:**

$2mF_o-DF_c$  (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.