



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2022 – 04:06 PM EDT

PDB ID : 5UW4
Title : Activated yeast Glycogen Synthase in complex with UDP glucosamine
Authors : Mahalingan, K.K.; Hurley, T.D.
Deposited on : 2017-02-20
Resolution : 2.98 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.27
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) : 2.27

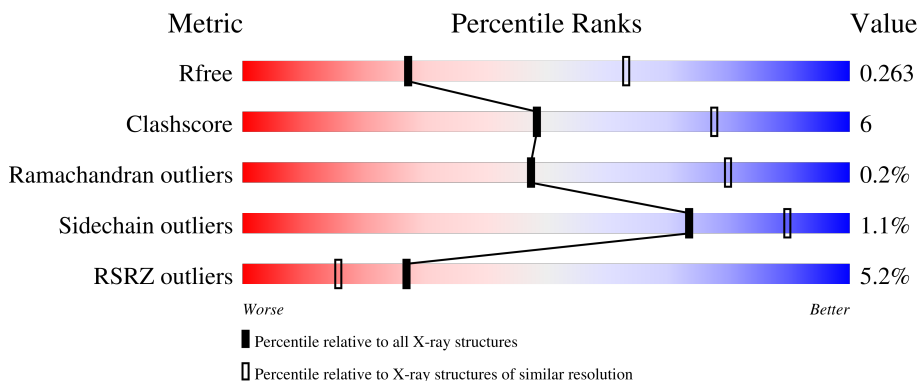
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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	2754 (3.00-2.96)
Clashscore	141614	3103 (3.00-2.96)
Ramachandran outliers	138981	2993 (3.00-2.96)
Sidechain outliers	138945	2996 (3.00-2.96)
RSRZ outliers	127900	2644 (3.00-2.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	720	 4% 77% 11% • 11%
1	B	720	 3% 79% 9% • 12%
1	C	720	 6% 77% 11% • 11%
1	D	720	 6% 75% 13% • 12%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 20374 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 Glycogen [starch] synthase isoform 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	C	638	5127	3272	894	942	19	0	0	0
1	B	636	5084	3246	886	933	19	0	0	0
1	A	638	4980	3185	853	924	18	0	0	0
1	D	636	5005	3192	868	927	18	0	0	0

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-19	MET	-	initiating methionine	UNP P27472
C	-18	GLY	-	expression tag	UNP P27472
C	-17	SER	-	expression tag	UNP P27472
C	-16	SER	-	expression tag	UNP P27472
C	-15	HIS	-	expression tag	UNP P27472
C	-14	HIS	-	expression tag	UNP P27472
C	-13	HIS	-	expression tag	UNP P27472
C	-12	HIS	-	expression tag	UNP P27472
C	-11	HIS	-	expression tag	UNP P27472
C	-10	HIS	-	expression tag	UNP P27472
C	-9	SER	-	expression tag	UNP P27472
C	-8	SER	-	expression tag	UNP P27472
C	-7	GLY	-	expression tag	UNP P27472
C	-6	LEU	-	expression tag	UNP P27472
C	-5	VAL	-	expression tag	UNP P27472
C	-4	PRO	-	expression tag	UNP P27472
C	-3	ARG	-	expression tag	UNP P27472
C	-2	GLY	-	expression tag	UNP P27472
C	-1	SER	-	expression tag	UNP P27472
C	0	HIS	-	expression tag	UNP P27472
C	169	GLN	GLU	engineered mutation	UNP P27472

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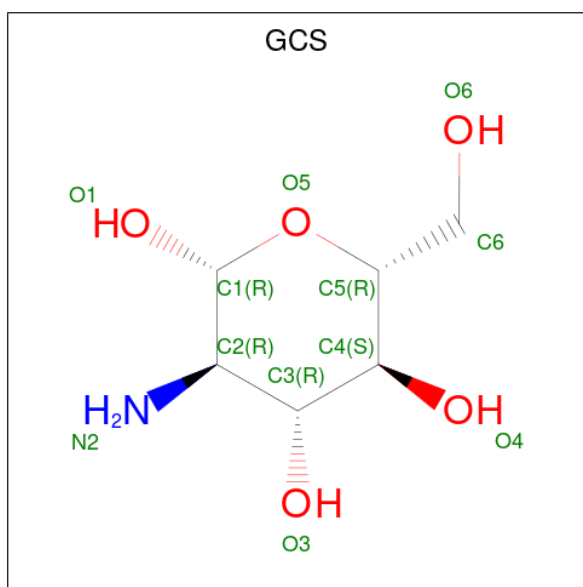
Chain	Residue	Modelled	Actual	Comment	Reference
C	535	SER	ALA	conflict	UNP P27472
B	-19	MET	-	initiating methionine	UNP P27472
B	-18	GLY	-	expression tag	UNP P27472
B	-17	SER	-	expression tag	UNP P27472
B	-16	SER	-	expression tag	UNP P27472
B	-15	HIS	-	expression tag	UNP P27472
B	-14	HIS	-	expression tag	UNP P27472
B	-13	HIS	-	expression tag	UNP P27472
B	-12	HIS	-	expression tag	UNP P27472
B	-11	HIS	-	expression tag	UNP P27472
B	-10	HIS	-	expression tag	UNP P27472
B	-9	SER	-	expression tag	UNP P27472
B	-8	SER	-	expression tag	UNP P27472
B	-7	GLY	-	expression tag	UNP P27472
B	-6	LEU	-	expression tag	UNP P27472
B	-5	VAL	-	expression tag	UNP P27472
B	-4	PRO	-	expression tag	UNP P27472
B	-3	ARG	-	expression tag	UNP P27472
B	-2	GLY	-	expression tag	UNP P27472
B	-1	SER	-	expression tag	UNP P27472
B	0	HIS	-	expression tag	UNP P27472
B	169	GLN	GLU	engineered mutation	UNP P27472
B	535	SER	ALA	conflict	UNP P27472
A	-19	MET	-	initiating methionine	UNP P27472
A	-18	GLY	-	expression tag	UNP P27472
A	-17	SER	-	expression tag	UNP P27472
A	-16	SER	-	expression tag	UNP P27472
A	-15	HIS	-	expression tag	UNP P27472
A	-14	HIS	-	expression tag	UNP P27472
A	-13	HIS	-	expression tag	UNP P27472
A	-12	HIS	-	expression tag	UNP P27472
A	-11	HIS	-	expression tag	UNP P27472
A	-10	HIS	-	expression tag	UNP P27472
A	-9	SER	-	expression tag	UNP P27472
A	-8	SER	-	expression tag	UNP P27472
A	-7	GLY	-	expression tag	UNP P27472
A	-6	LEU	-	expression tag	UNP P27472
A	-5	VAL	-	expression tag	UNP P27472
A	-4	PRO	-	expression tag	UNP P27472
A	-3	ARG	-	expression tag	UNP P27472
A	-2	GLY	-	expression tag	UNP P27472
A	-1	SER	-	expression tag	UNP P27472

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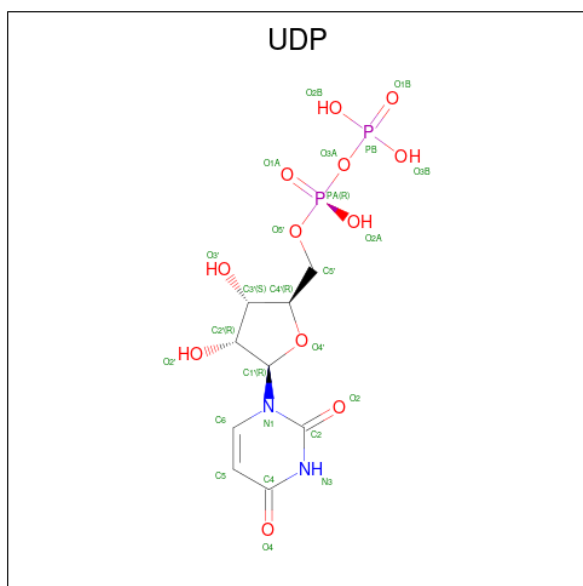
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	expression tag	UNP P27472
A	169	GLN	GLU	engineered mutation	UNP P27472
A	535	SER	ALA	conflict	UNP P27472
D	-19	MET	-	initiating methionine	UNP P27472
D	-18	GLY	-	expression tag	UNP P27472
D	-17	SER	-	expression tag	UNP P27472
D	-16	SER	-	expression tag	UNP P27472
D	-15	HIS	-	expression tag	UNP P27472
D	-14	HIS	-	expression tag	UNP P27472
D	-13	HIS	-	expression tag	UNP P27472
D	-12	HIS	-	expression tag	UNP P27472
D	-11	HIS	-	expression tag	UNP P27472
D	-10	HIS	-	expression tag	UNP P27472
D	-9	SER	-	expression tag	UNP P27472
D	-8	SER	-	expression tag	UNP P27472
D	-7	GLY	-	expression tag	UNP P27472
D	-6	LEU	-	expression tag	UNP P27472
D	-5	VAL	-	expression tag	UNP P27472
D	-4	PRO	-	expression tag	UNP P27472
D	-3	ARG	-	expression tag	UNP P27472
D	-2	GLY	-	expression tag	UNP P27472
D	-1	SER	-	expression tag	UNP P27472
D	0	HIS	-	expression tag	UNP P27472
D	169	GLN	GLU	engineered mutation	UNP P27472
D	535	SER	ALA	conflict	UNP P27472

- Molecule 2 is 2-amino-2-deoxy-beta-D-glucopyranose (three-letter code: GCS) (formula: $C_6H_{13}NO_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	C	1	12	6	1	5	0	0

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



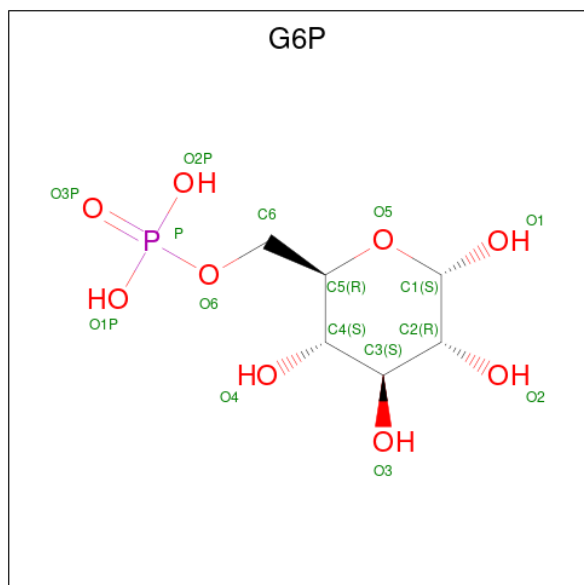
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	C	1	25	9	2	12	2	0	0
3	B	1	25	9	2	12	2	0	0
3	A	1	25	9	2	12	2	0	0

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

- Molecule 4 is 6-O-phosphono-alpha-D-glucopyranose (three-letter code: G6P) (formula: $C_6H_{13}O_9P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
4	C	1	16	6	9	1	0	0
4	B	1	16	6	9	1	0	0
4	A	1	16	6	9	1	0	0
4	D	1	16	6	9	1	0	0

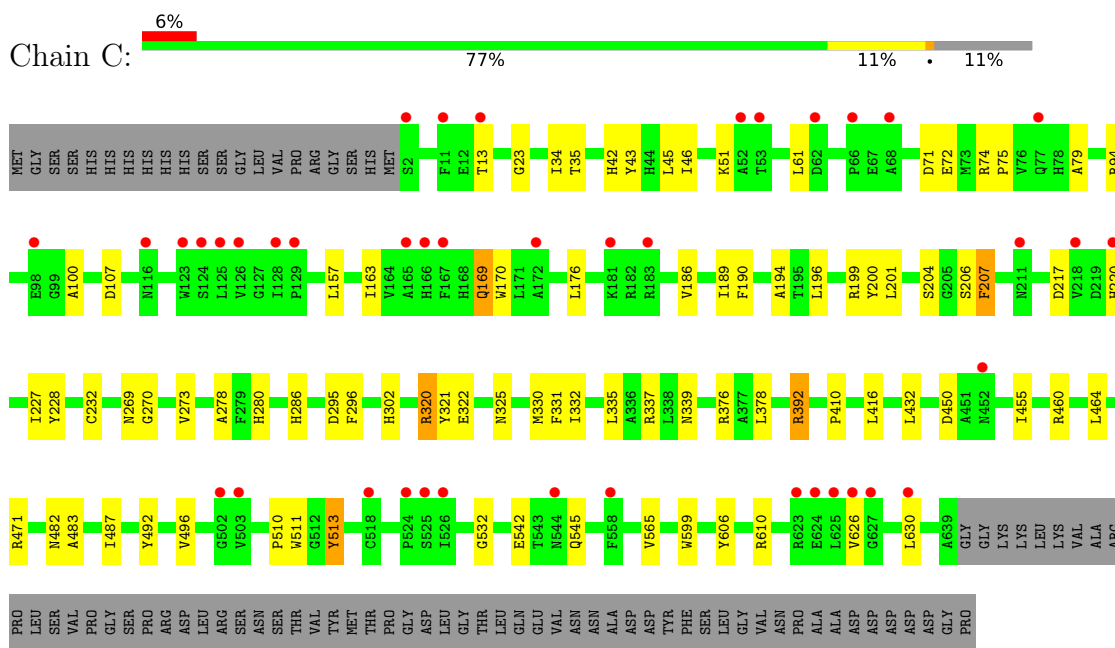
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	C	1	1	1	0	0
5	B	1	1	1	0	0

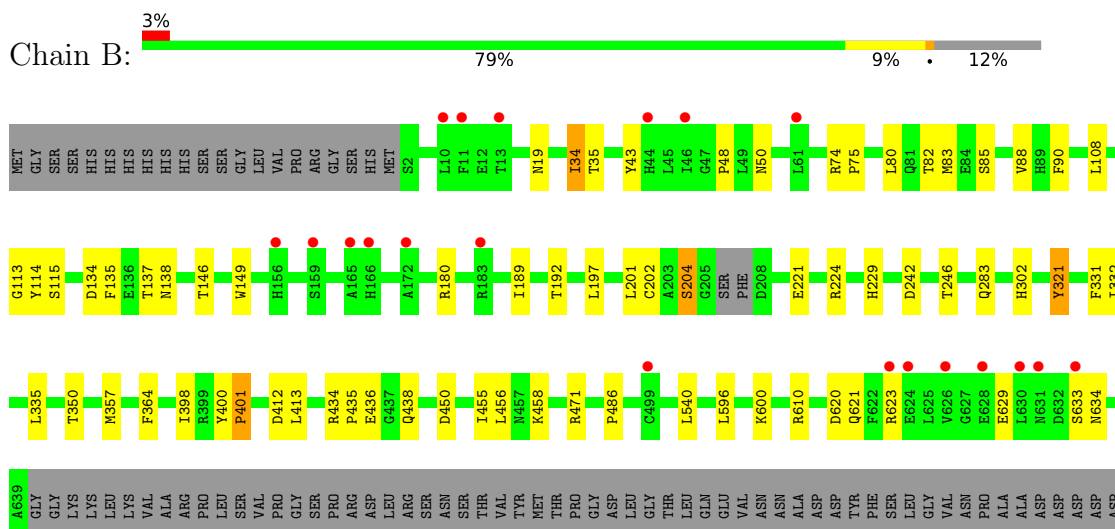
3 Residue-property plots [i](#)

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: Glycogen [starch] synthase isoform 2



- Molecule 1: Glycogen [starch] synthase isoform 2



4 Data and refinement statistics i

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	193.04Å 204.70Å 205.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	145.16 – 2.98 29.95 – 2.98	Depositor EDS
% Data completeness (in resolution range)	99.6 (145.16-2.98) 99.7 (29.95-2.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 3.00Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.188 , 0.257 0.197 , 0.263	Depositor DCC
R_{free} test set	4474 reflections (5.42%)	wwPDB-VP
Wilson B-factor (Å ²)	74.6	Xtrriage
Anisotropy	0.088	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 59.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.003 for -h,-l,-k	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	20374	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GCS, UDP, G6P

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	0.56	0/5101	0.76	3/6943 (0.0%)
1	B	0.59	0/5206	0.83	1/7061 (0.0%)
1	C	0.57	0/5251	0.80	6/7119 (0.1%)
1	D	0.56	0/5123	0.78	3/6962 (0.0%)
All	All	0.57	0/20681	0.80	13/28085 (0.0%)

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	D	0	1

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	199	ARG	NE-CZ-NH2	-6.70	116.95	120.30
1	B	34	ILE	N-CA-CB	6.51	125.78	110.80
1	C	320	ARG	NE-CZ-NH1	6.48	123.54	120.30
1	A	131	PRO	N-CA-C	-6.29	95.73	112.10
1	C	199	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	D	125	LEU	N-CA-C	6.03	127.28	111.00
1	A	124	SER	O-C-N	5.94	132.20	122.70
1	C	392	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	C	471	ARG	CG-CD-NE	5.59	123.55	111.80
1	D	306	ASP	CB-CG-OD1	5.37	123.13	118.30
1	D	124	SER	N-CA-C	5.28	125.26	111.00
1	A	376	ARG	NE-CZ-NH1	5.28	122.94	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	337	ARG	NE-CZ-NH1	5.19	122.89	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	18	ALA	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	4980	0	4753	61	0
1	B	5084	0	4955	49	0
1	C	5127	0	5015	64	0
1	D	5005	0	4827	70	0
2	C	12	0	13	1	0
3	A	25	0	11	1	0
3	B	25	0	11	0	0
3	C	25	0	11	1	0
3	D	25	0	11	2	0
4	A	16	0	11	1	0
4	B	16	0	11	1	0
4	C	16	0	11	1	0
4	D	16	0	11	2	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
All	All	20374	0	19651	239	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 6.

All (239) 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:217:ASP:CB	1:C:220:HIS:HD2	1.36	1.39

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:217:ASP:CB	1:C:220:HIS:CD2	2.28	1.16
1:C:217:ASP:HB3	1:C:220:HIS:CD2	1.80	1.16
1:C:217:ASP:CG	1:C:220:HIS:CD2	2.20	1.14
1:C:217:ASP:HB3	1:C:220:HIS:HD2	0.97	1.10
1:B:620:ASP:OD1	1:B:621:GLN:N	1.92	1.02
1:A:121:ASP:OD1	1:A:125:LEU:HD23	1.59	1.02
1:B:620:ASP:OD1	1:B:621:GLN:HG3	1.62	0.99
1:D:70:SER:OG	1:D:73:MET:HB2	1.62	0.99
1:A:123:TRP:O	1:A:127:GLY:N	1.97	0.97
1:C:13:THR:OG1	1:C:46:ILE:HB	1.70	0.90
1:A:121:ASP:OD1	1:A:125:LEU:CD2	2.18	0.90
1:D:49:LEU:HD12	1:D:50:ASN:N	1.88	0.88
1:D:41:ASP:OD1	1:D:73:MET:HG3	1.75	0.85
1:C:217:ASP:CG	1:C:220:HIS:HD2	1.66	0.82
1:C:169:GLN:HE21	1:C:170:TRP:H	1.29	0.81
1:A:96:LEU:HG	1:A:96:LEU:O	1.81	0.78
1:A:28:LEU:HD22	1:A:45:LEU:HD21	1.69	0.74
1:D:49:LEU:HB2	1:D:105:LEU:HB3	1.69	0.74
1:B:620:ASP:OD1	1:B:621:GLN:CG	2.36	0.73
1:C:450:ASP:OD1	1:C:460:ARG:NH2	2.22	0.72
1:A:626:VAL:HG11	1:A:630:LEU:HD11	1.70	0.72
1:D:128:ILE:HD12	1:D:232:CYS:HB3	1.72	0.72
1:D:39:TYR:HB3	1:D:42:HIS:HB2	1.73	0.71
1:D:214:GLU:HB3	1:D:257:HIS:ND1	2.07	0.70
1:A:170:TRP:CZ3	1:A:171:LEU:HD23	2.26	0.70
1:C:169:GLN:NE2	1:C:169:GLN:HA	2.05	0.70
1:A:73:MET:O	1:A:76:VAL:HG12	1.91	0.69
1:C:61:LEU:HD21	1:C:94:ARG:HH22	1.59	0.68
1:D:128:ILE:HG23	1:D:232:CYS:HB3	1.76	0.68
1:C:330:MET:HG3	1:C:565:VAL:HG22	1.78	0.66
1:B:357:MET:CE	1:B:455:ILE:HD11	2.25	0.66
1:A:170:TRP:O	1:A:173:GLY:N	2.30	0.65
1:C:23:GLY:N	3:C:802:UDP:O3B	2.31	0.64
1:D:14:ALA:HB2	1:D:168:HIS:HB2	1.80	0.63
1:D:213:LEU:HD12	1:D:258:LEU:HD21	1.81	0.63
1:C:626:VAL:HG11	1:C:630:LEU:HD11	1.82	0.62
1:C:61:LEU:HD21	1:C:94:ARG:NH2	2.14	0.62
1:C:169:GLN:HE21	1:C:170:TRP:N	1.98	0.61
1:C:79:ALA:HB2	1:C:157:LEU:HD12	1.83	0.61
1:A:280:HIS:CE1	1:D:283:GLN:HG2	2.37	0.60
1:B:80:LEU:HD22	1:B:90:PHE:CE2	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:398:ILE:CD1	1:A:298:ARG:HD3	2.31	0.60
1:D:49:LEU:HD12	1:D:49:LEU:C	2.19	0.59
1:C:280:HIS:CE1	1:B:283:GLN:HG2	2.38	0.59
1:D:350:THR:OG1	1:D:471:ARG:NH1	2.36	0.59
1:D:372:GLN:HG3	1:D:487:ILE:HG22	1.85	0.59
1:B:540:LEU:HD23	1:B:596:LEU:HD13	1.86	0.58
1:B:197:LEU:O	1:B:201:LEU:HB2	2.03	0.58
1:A:549:TYR:CD2	1:A:593:LEU:HD13	2.39	0.58
1:B:114:TYR:O	1:B:115:SER:C	2.42	0.58
1:B:82:THR:O	1:B:85:SER:HB3	2.04	0.58
1:A:131:PRO:HG2	1:A:229:HIS:CE1	2.39	0.58
1:D:187:VAL:HG12	1:D:613:ALA:HB1	1.84	0.58
1:A:381:THR:HG21	1:A:428:ARG:HE	1.68	0.57
1:D:91:VAL:HG13	1:D:105:LEU:HB2	1.84	0.57
1:D:292:LYS:HD2	1:D:490:LEU:HD21	1.87	0.57
1:D:49:LEU:HD11	1:D:54:TYR:CD1	2.40	0.56
1:A:15:THR:HG23	1:A:16:GLU:HG3	1.86	0.56
1:C:295:ASP:OD2	1:C:376:ARG:NH2	2.39	0.56
1:A:126:VAL:HG12	1:A:126:VAL:O	2.04	0.56
1:B:202:CYS:C	1:B:204:SER:H	2.09	0.56
1:A:170:TRP:CH2	1:A:171:LEU:CD2	2.89	0.55
1:C:492:TYR:O	1:C:496:VAL:HG23	2.05	0.55
1:C:278:ALA:HB1	1:C:280:HIS:CE1	2.42	0.55
1:B:202:CYS:O	1:B:204:SER:N	2.41	0.54
1:D:205:GLY:O	1:D:208:ASP:N	2.41	0.54
1:A:57:GLU:O	1:A:96:LEU:HB3	2.07	0.54
1:A:97:ILE:HG13	1:A:98:GLU:N	2.23	0.54
1:A:170:TRP:O	1:A:171:LEU:C	2.46	0.54
1:C:228:TYR:CE2	1:C:232:CYS:SG	3.01	0.54
1:B:400:TYR:N	1:B:401:PRO:CD	2.70	0.54
1:D:331:PHE:CZ	1:D:335:LEU:HD11	2.42	0.54
1:C:286:HIS:NE2	4:C:803:G6P:O2P	2.35	0.53
1:A:19:ASN:ND2	1:A:21:VAL:HG23	2.22	0.53
1:C:217:ASP:OD1	1:C:220:HIS:CD2	2.59	0.53
1:B:398:ILE:HD13	1:A:298:ARG:HD3	1.89	0.53
1:D:210:TYR:CZ	1:D:530:VAL:HG13	2.43	0.53
1:C:201:LEU:HB3	1:C:207:PHE:HE2	1.73	0.53
1:A:126:VAL:O	1:A:126:VAL:CG1	2.57	0.53
1:B:180:ARG:NH2	1:B:242:ASP:OD1	2.41	0.52
1:A:283:GLN:HG3	4:A:802:G6P:O1	2.09	0.52
1:C:43:TYR:OH	1:C:45:LEU:HD21	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:214:GLU:HB3	1:D:257:HIS:CE1	2.45	0.52
1:D:74:ARG:NH1	1:D:77:GLN:OE1	2.43	0.52
1:A:95:TRP:CE3	1:A:97:ILE:HG22	2.45	0.52
1:D:41:ASP:CG	1:D:73:MET:HG3	2.29	0.52
1:B:321:TYR:CZ	1:B:455:ILE:HD12	2.45	0.52
1:D:111:VAL:HG11	1:D:118:TRP:CH2	2.45	0.52
1:D:49:LEU:HA	1:D:105:LEU:HD23	1.92	0.52
1:A:271:LEU:HD13	1:A:520:VAL:HG21	1.92	0.51
1:A:538:GLU:HG3	1:A:543:THR:HG21	1.93	0.51
1:C:204:SER:O	1:C:206:SER:N	2.43	0.51
1:A:16:GLU:HG2	1:A:21:VAL:HB	1.92	0.51
1:A:18:ALA:HB2	1:A:105:LEU:CD2	2.40	0.51
1:A:95:TRP:HB3	1:A:101:PRO:HD2	1.91	0.51
1:C:169:GLN:NE2	1:C:169:GLN:CA	2.73	0.51
1:A:322:GLU:HB3	1:A:325:ASN:HB2	1.93	0.51
1:B:350:THR:OG1	1:B:471:ARG:NH1	2.44	0.51
1:B:283:GLN:HG3	4:B:802:G6P:O1	2.11	0.51
1:B:435:PRO:O	1:B:436:GLU:C	2.47	0.51
1:D:293:ILE:O	1:D:297:VAL:HG23	2.10	0.51
1:D:507:TYR:HB2	1:D:556:ARG:NH1	2.25	0.50
1:B:35:THR:HG21	1:B:43:TYR:CE1	2.46	0.50
1:A:131:PRO:O	1:A:133:ASN:N	2.45	0.50
1:A:296:PHE:HE1	1:A:487:ILE:HD12	1.76	0.50
1:B:202:CYS:C	1:B:204:SER:N	2.65	0.50
1:B:357:MET:HE3	1:B:455:ILE:HD11	1.93	0.50
1:D:34:ILE:HG21	1:D:599:TRP:HB3	1.94	0.50
1:B:189:ILE:HD11	1:B:610:ARG:HA	1.94	0.49
1:D:286:HIS:NE2	4:D:802:G6P:O3P	2.32	0.49
1:D:74:ARG:N	1:D:75:PRO:CD	2.76	0.49
1:A:357:MET:O	1:A:478:PRO:HA	2.12	0.49
1:B:633:SER:O	1:B:634:ASN:ND2	2.45	0.49
1:A:97:ILE:HG13	1:A:98:GLU:H	1.77	0.49
1:C:510:PRO:O	1:C:532:GLY:HA3	2.13	0.49
1:B:134:ASP:O	1:B:138:ASN:ND2	2.46	0.48
1:A:170:TRP:CH2	1:A:171:LEU:HD23	2.48	0.48
1:A:510:PRO:O	1:A:532:GLY:HA3	2.13	0.48
1:D:388:SER:HB3	1:D:392:ARG:NH2	2.29	0.48
1:D:128:ILE:HG23	1:D:232:CYS:CB	2.42	0.48
1:D:192:THR:HG23	1:D:246:THR:HG22	1.95	0.48
1:C:378:LEU:HD22	1:C:432:LEU:HD11	1.95	0.47
1:A:151:LEU:HA	1:A:154:VAL:HG12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:TYR:CE2	1:C:227:ILE:HD11	2.49	0.47
1:A:176:LEU:HB2	1:A:177:PRO:HD3	1.97	0.47
1:B:331:PHE:CZ	1:B:335:LEU:HD11	2.49	0.47
1:B:364:PHE:CE2	1:B:486:PRO:HD2	2.50	0.47
1:A:191:THR:HA	1:A:245:THR:O	2.15	0.47
1:D:176:LEU:HD22	1:D:241:ALA:HB2	1.97	0.47
1:A:74:ARG:O	1:A:75:PRO:C	2.53	0.47
1:D:138:ASN:O	1:D:141:ILE:HG22	2.15	0.47
1:B:19:ASN:HD22	1:B:50:ASN:HD22	1.63	0.46
1:C:273:VAL:HG13	1:C:273:VAL:O	2.15	0.46
1:D:304:CYS:SG	1:D:434:ARG:HD3	2.55	0.46
1:D:580:ARG:NH1	4:D:802:G6P:O1P	2.41	0.46
1:D:325:ASN:ND2	1:D:508:TYR:HB3	2.31	0.45
1:B:623:ARG:HG3	1:B:629:GLU:HB2	1.98	0.45
1:B:434:ARG:HH21	1:B:438:GLN:NE2	2.14	0.45
1:D:47:GLY:O	1:D:105:LEU:HA	2.15	0.45
1:D:126:VAL:HG23	1:D:128:ILE:HG12	1.99	0.45
1:D:278:ALA:HB1	1:D:280:HIS:CE1	2.52	0.45
1:C:270:GLY:O	1:C:513:TYR:OH	2.20	0.45
1:A:231:TYR:CD1	1:A:231:TYR:C	2.89	0.45
1:B:135:PHE:HA	1:B:138:ASN:HD22	1.82	0.45
1:D:144:GLY:HA3	1:D:174:VAL:CG2	2.47	0.45
1:C:207:PHE:HD2	1:C:207:PHE:N	2.15	0.45
1:C:322:GLU:HB3	1:C:325:ASN:HB2	1.99	0.45
1:B:302:HIS:O	1:B:434:ARG:HD2	2.16	0.45
1:C:71:ASP:OD1	1:C:74:ARG:NH1	2.50	0.45
1:C:228:TYR:CD2	1:C:232:CYS:SG	3.10	0.45
1:C:320:ARG:NH2	1:C:322:GLU:OE1	2.50	0.45
1:B:302:HIS:O	1:B:434:ARG:NH1	2.44	0.45
1:A:218:VAL:HG11	1:A:231:TYR:OH	2.17	0.45
1:D:144:GLY:HA3	1:D:174:VAL:HG23	1.99	0.45
1:C:207:PHE:N	1:C:207:PHE:CD2	2.84	0.44
1:C:201:LEU:HB3	1:C:207:PHE:CE2	2.51	0.44
1:B:192:THR:CG2	1:B:246:THR:HG22	2.47	0.44
1:A:170:TRP:CD2	1:A:171:LEU:N	2.86	0.44
1:A:74:ARG:O	1:A:76:VAL:N	2.51	0.44
1:D:434:ARG:HB2	1:D:435:PRO:HD2	1.99	0.44
1:C:74:ARG:N	1:C:75:PRO:CD	2.81	0.44
1:C:339:ASN:HD22	1:C:464:LEU:CD2	2.31	0.44
1:A:15:THR:HG21	1:A:171:LEU:HD13	1.98	0.44
1:A:16:GLU:OE1	1:A:24:ILE:HD12	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:42:HIS:NE2	1:C:72:GLU:OE2	2.51	0.43
1:C:606:TYR:HB3	1:C:610:ARG:NH2	2.33	0.43
1:B:34:ILE:HD12	1:B:600:LYS:HA	1.99	0.43
1:A:97:ILE:CG1	1:A:98:GLU:H	2.32	0.43
1:A:217:ASP:O	1:A:221:GLU:HG2	2.18	0.43
1:C:35:THR:HG21	1:C:43:TYR:CZ	2.53	0.43
1:B:412:ASP:C	1:B:413:LEU:O	2.54	0.43
1:D:96:LEU:O	1:D:97:ILE:HD12	2.19	0.43
1:D:197:LEU:HD13	1:D:258:LEU:HB3	2.00	0.43
1:D:331:PHE:CE1	1:D:335:LEU:HD11	2.54	0.43
1:C:296:PHE:HE1	1:C:487:ILE:HD12	1.84	0.43
1:C:482:ASN:OD1	1:C:483:ALA:N	2.52	0.43
1:B:74:ARG:N	1:B:75:PRO:CD	2.82	0.43
1:D:356:VAL:HG12	1:D:356:VAL:O	2.19	0.43
1:D:166:HIS:CD2	1:D:189:ILE:HB	2.54	0.43
1:B:83:MET:HE2	1:B:88:VAL:HG11	2.01	0.42
1:C:194:ALA:HB2	1:C:510:PRO:HG2	2.01	0.42
1:B:221:GLU:OE2	1:B:224:ARG:NH1	2.52	0.42
1:B:332:ILE:HG21	1:B:458:LYS:HG3	2.01	0.42
1:D:163:ILE:HB	1:D:186:VAL:HG12	2.00	0.42
1:B:113:GLY:O	1:B:115:SER:N	2.52	0.42
1:D:228:TYR:CE2	1:D:232:CYS:SG	3.13	0.42
1:D:548:ASP:O	1:D:589:ARG:HD2	2.18	0.42
1:C:189:ILE:HD11	1:C:610:ARG:HA	2.01	0.42
1:A:278:ALA:HB1	1:A:280:HIS:CE1	2.55	0.42
1:A:492:TYR:CD1	3:A:801:UDP:C5	3.07	0.42
1:B:400:TYR:N	1:B:401:PRO:HD3	2.33	0.42
1:A:29:LYS:HA	1:A:95:TRP:HZ3	1.84	0.42
1:A:36:VAL:O	1:A:40:LYS:N	2.48	0.42
1:B:137:THR:HG21	1:B:229:HIS:HD2	1.85	0.42
1:D:428:ARG:HA	1:D:428:ARG:HD3	1.85	0.42
1:C:163:ILE:HB	1:C:186:VAL:HG12	2.02	0.42
1:C:269:ASN:HB2	1:C:511:TRP:CD1	2.55	0.42
1:D:42:HIS:ND1	1:D:73:MET:SD	2.86	0.42
1:C:331:PHE:CZ	1:C:335:LEU:HD11	2.54	0.42
1:A:526:ILE:HG21	1:A:568:LEU:HD13	2.01	0.42
1:D:183:ARG:HD2	1:D:183:ARG:HA	1.60	0.42
1:C:94:ARG:CG	1:C:100:ALA:HB1	2.49	0.42
1:C:332:ILE:HD11	1:C:455:ILE:HG23	2.01	0.41
1:C:34:ILE:HD13	1:C:599:TRP:O	2.19	0.41
1:B:48:PRO:HB3	1:B:108:LEU:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:LEU:C	1:D:49:LEU:CD1	2.85	0.41
1:C:410:PRO:HG2	1:C:416:LEU:HD21	2.02	0.41
1:D:321:TYR:C	1:D:321:TYR:CD2	2.94	0.41
1:D:192:THR:CG2	1:D:246:THR:HG22	2.49	0.41
1:D:227:ILE:HG22	1:D:230:ARG:CB	2.50	0.41
1:C:170:TRP:NE1	1:C:196:LEU:HD21	2.35	0.41
1:C:176:LEU:HD11	1:C:190:PHE:HB2	2.03	0.41
1:C:392:ARG:NH1	1:C:416:LEU:O	2.53	0.41
1:B:450:ASP:OD1	1:B:456:LEU:HD13	2.20	0.41
1:A:584:ILE:HD13	1:D:278:ALA:HB2	2.02	0.41
1:D:282:PHE:CE1	1:D:591:GLU:HG3	2.56	0.41
1:C:51:LYS:NZ	1:C:107:ASP:OD1	2.42	0.41
1:B:146:THR:O	1:B:149:TRP:HB3	2.20	0.41
1:A:19:ASN:OD1	1:A:19:ASN:N	2.54	0.41
1:C:227:ILE:O	1:C:228:TYR:C	2.59	0.41
1:A:170:TRP:CH2	1:A:171:LEU:HD21	2.55	0.41
1:D:187:VAL:CG1	1:D:613:ALA:HB1	2.50	0.41
1:D:292:LYS:CD	1:D:490:LEU:HD21	2.48	0.41
1:A:170:TRP:O	1:A:172:ALA:N	2.53	0.41
1:C:511:TRP:HB3	2:C:801:GCS:O4	2.21	0.41
1:D:178:LEU:O	1:D:182:ARG:CB	2.69	0.41
1:A:626:VAL:HG12	1:A:626:VAL:O	2.21	0.41
1:A:332:ILE:HD11	1:A:455:ILE:HG23	2.01	0.40
1:D:95:TRP:CD1	1:D:97:ILE:HD13	2.56	0.40
1:D:514:THR:HG23	3:D:801:UDP:H5'2	2.03	0.40
1:C:302:HIS:O	1:C:302:HIS:CG	2.74	0.40
1:B:192:THR:HG22	1:B:246:THR:HG22	2.04	0.40
1:D:396:HIS:ND1	1:D:415:GLU:OE1	2.54	0.40
1:D:514:THR:HG23	3:D:801:UDP:O1B	2.22	0.40
1:B:113:GLY:O	1:B:114:TYR:C	2.60	0.40
1:C:542:GLU:OE1	1:C:545:GLN:N	2.47	0.40
1:D:296:PHE:HA	1:D:372:GLN:HE22	1.86	0.40
1:A:170:TRP:CE3	1:A:171:LEU:N	2.89	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	636/720 (88%)	575 (90%)	58 (9%)	3 (0%)	29	66
1	B	632/720 (88%)	585 (93%)	46 (7%)	1 (0%)	47	80
1	C	636/720 (88%)	586 (92%)	50 (8%)	0	100	100
1	D	632/720 (88%)	579 (92%)	52 (8%)	1 (0%)	47	80
All	All	2536/2880 (88%)	2325 (92%)	206 (8%)	5 (0%)	47	80

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	401	PRO
1	A	544	ASN
1	A	183	ARG
1	A	363	SER
1	D	320	ARG

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	516/621 (83%)	504 (98%)	12 (2%)	50	79
1	B	539/621 (87%)	537 (100%)	2 (0%)	91	97
1	C	547/621 (88%)	543 (99%)	4 (1%)	84	93
1	D	523/621 (84%)	517 (99%)	6 (1%)	73	90
All	All	2125/2484 (86%)	2101 (99%)	24 (1%)	73	90

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

Mol	Chain	Res	Type
1	C	169	GLN
1	C	207	PHE
1	C	321	TYR
1	C	513	TYR
1	B	204	SER
1	B	321	TYR
1	A	2	SER
1	A	45	LEU
1	A	58	VAL
1	A	71	ASP
1	A	95	TRP
1	A	124	SER
1	A	125	LEU
1	A	326	LYS
1	A	363	SER
1	A	467	SER
1	A	469	SER
1	A	513	TYR
1	D	41	ASP
1	D	49	LEU
1	D	123	TRP
1	D	183	ARG
1	D	184	ILE
1	D	321	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	220	HIS
1	B	19	ASN
1	B	283	GLN
1	B	403	ASN
1	B	634	ASN
1	D	283	GLN
1	D	372	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	UDP	B	801	-	20,26,26	0.78	0	25,40,40	1.01	0
4	G6P	A	802	-	16,16,16	0.70	0	24,24,24	1.19	1 (4%)
3	UDP	D	801	-	20,26,26	0.80	0	25,40,40	1.15	1 (4%)
4	G6P	C	803	-	16,16,16	0.77	0	24,24,24	1.34	3 (12%)
3	UDP	C	802	-	20,26,26	0.96	1 (5%)	25,40,40	1.36	5 (20%)
4	G6P	B	802	-	16,16,16	0.59	0	24,24,24	1.27	2 (8%)
2	GCS	C	801	-	12,12,12	0.73	0	16,17,17	1.62	3 (18%)
3	UDP	A	801	-	20,26,26	0.79	0	25,40,40	1.37	4 (16%)
4	G6P	D	802	-	16,16,16	0.65	0	24,24,24	1.42	4 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	UDP	B	801	-	-	5/14/32/32	0/2/2/2
4	G6P	A	802	-	-	2/6/26/26	0/1/1/1
3	UDP	D	801	-	-	1/14/32/32	0/2/2/2
4	G6P	C	803	-	-	1/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	UDP	C	802	-	-	5/14/32/32	0/2/2/2
4	G6P	B	802	-	-	0/6/26/26	0/1/1/1
2	GCS	C	801	-	-	0/2/22/22	0/1/1/1
3	UDP	A	801	-	-	7/14/32/32	0/2/2/2
4	G6P	D	802	-	-	1/6/26/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	802	UDP	O4'-C1'	2.72	1.44	1.41

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	801	GCS	O5-C1-C2	4.03	114.21	109.51
4	D	802	G6P	O3-C3-C2	-3.55	102.15	110.35
3	A	801	UDP	PA-O3A-PB	-3.44	121.03	132.83
3	D	801	UDP	C3'-C2'-C1'	3.15	105.72	100.98
4	B	802	G6P	O2P-P-O1P	2.85	118.51	107.64
3	C	802	UDP	PA-O3A-PB	-2.79	123.25	132.83
3	A	801	UDP	O3B-PB-O2B	2.77	118.24	107.64
3	A	801	UDP	C3'-C2'-C1'	2.77	105.15	100.98
4	C	803	G6P	O2P-P-O1P	2.60	117.57	107.64
4	D	802	G6P	O3-C3-C4	-2.57	104.41	110.35
3	C	802	UDP	C3'-C2'-C1'	2.56	104.83	100.98
4	C	803	G6P	C4-C3-C2	2.46	115.11	110.82
2	C	801	GCS	C3-C2-N2	-2.42	106.10	111.05
4	D	802	G6P	C4-C3-C2	2.41	115.03	110.82
3	C	802	UDP	C6-N1-C2	-2.39	117.40	121.20
4	B	802	G6P	O3-C3-C4	-2.32	104.98	110.35
4	C	803	G6P	O3-C3-C2	-2.26	105.12	110.35
3	C	802	UDP	O3A-PB-O1B	-2.23	98.81	111.19
3	A	801	UDP	O2A-PA-O1A	2.06	122.42	112.24
3	C	802	UDP	O3B-PB-O2B	2.04	115.43	107.64
2	C	801	GCS	C1-C2-C3	-2.03	107.93	110.60
4	D	802	G6P	O2P-P-O1P	2.03	115.39	107.64
4	A	802	G6P	O6-C6-C5	2.02	115.96	108.99

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	802	UDP	C2'-C1'-N1-C6
3	C	802	UDP	O4'-C1'-N1-C6
3	B	801	UDP	C3'-C4'-C5'-O5'
3	B	801	UDP	PA-O3A-PB-O3B
3	A	801	UDP	C2'-C1'-N1-C6
3	A	801	UDP	O4'-C1'-N1-C6
3	A	801	UDP	C5'-O5'-PA-O1A
4	A	802	G6P	C4-C5-C6-O6
4	A	802	G6P	O5-C5-C6-O6
4	D	802	G6P	C4-C5-C6-O6
3	C	802	UDP	O4'-C4'-C5'-O5'
3	B	801	UDP	O4'-C4'-C5'-O5'
3	A	801	UDP	O4'-C4'-C5'-O5'
3	C	802	UDP	PB-O3A-PA-O5'
3	A	801	UDP	PB-O3A-PA-O5'
3	D	801	UDP	PB-O3A-PA-O2A
3	C	802	UDP	C3'-C4'-C5'-O5'
3	A	801	UDP	C3'-C4'-C5'-O5'
4	C	803	G6P	C4-C5-C6-O6
3	B	801	UDP	PA-O3A-PB-O1B
3	B	801	UDP	PA-O3A-PB-O2B
3	A	801	UDP	PA-O3A-PB-O2B

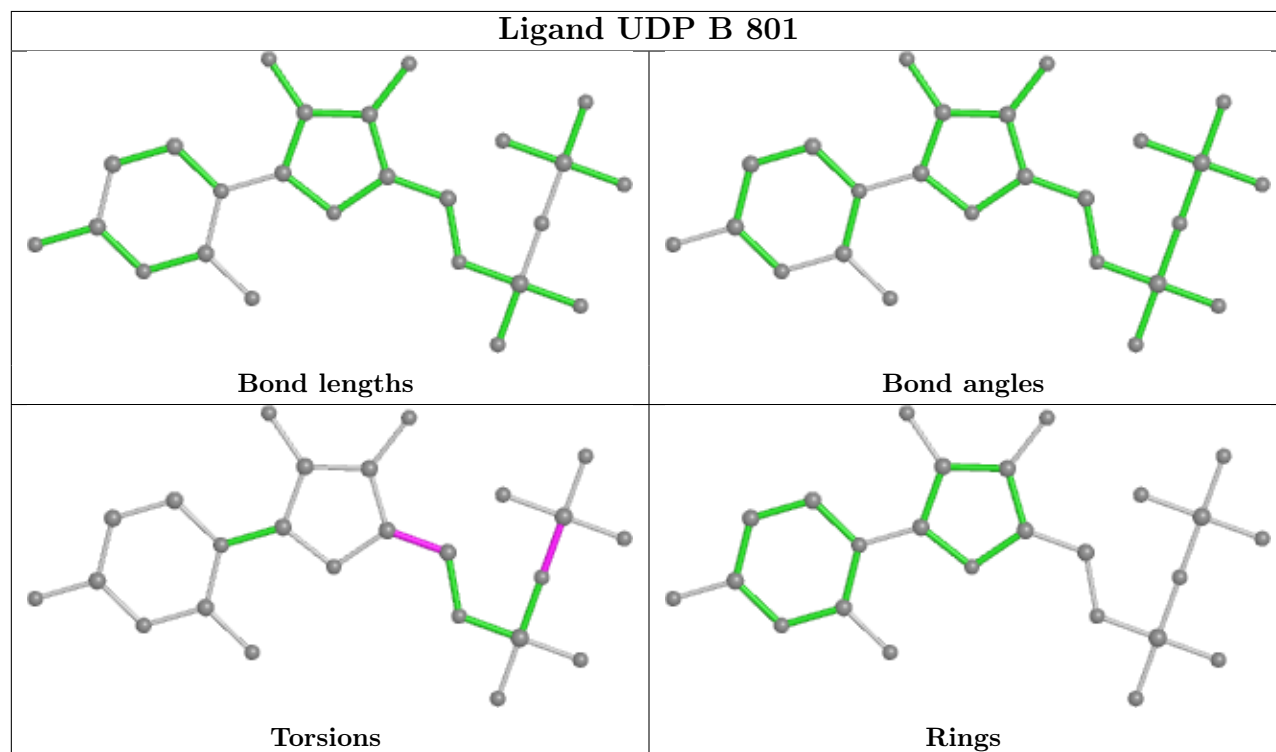
There are no ring outliers.

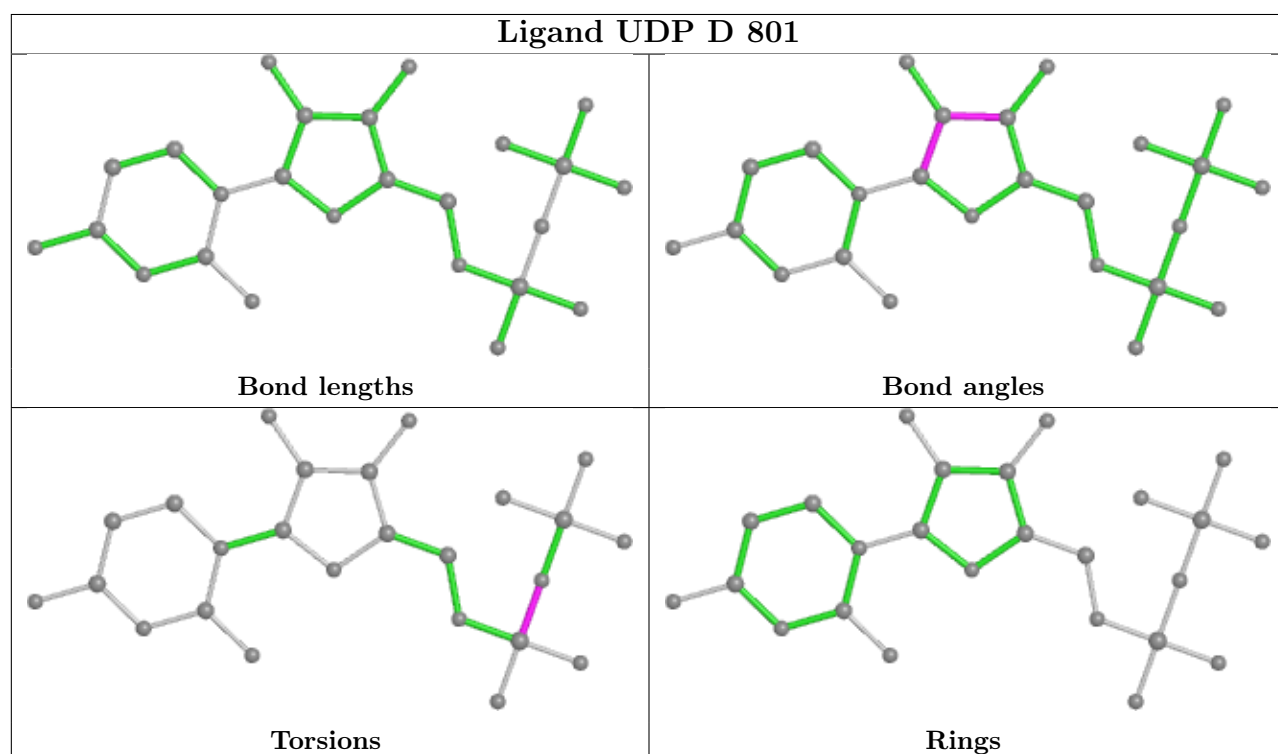
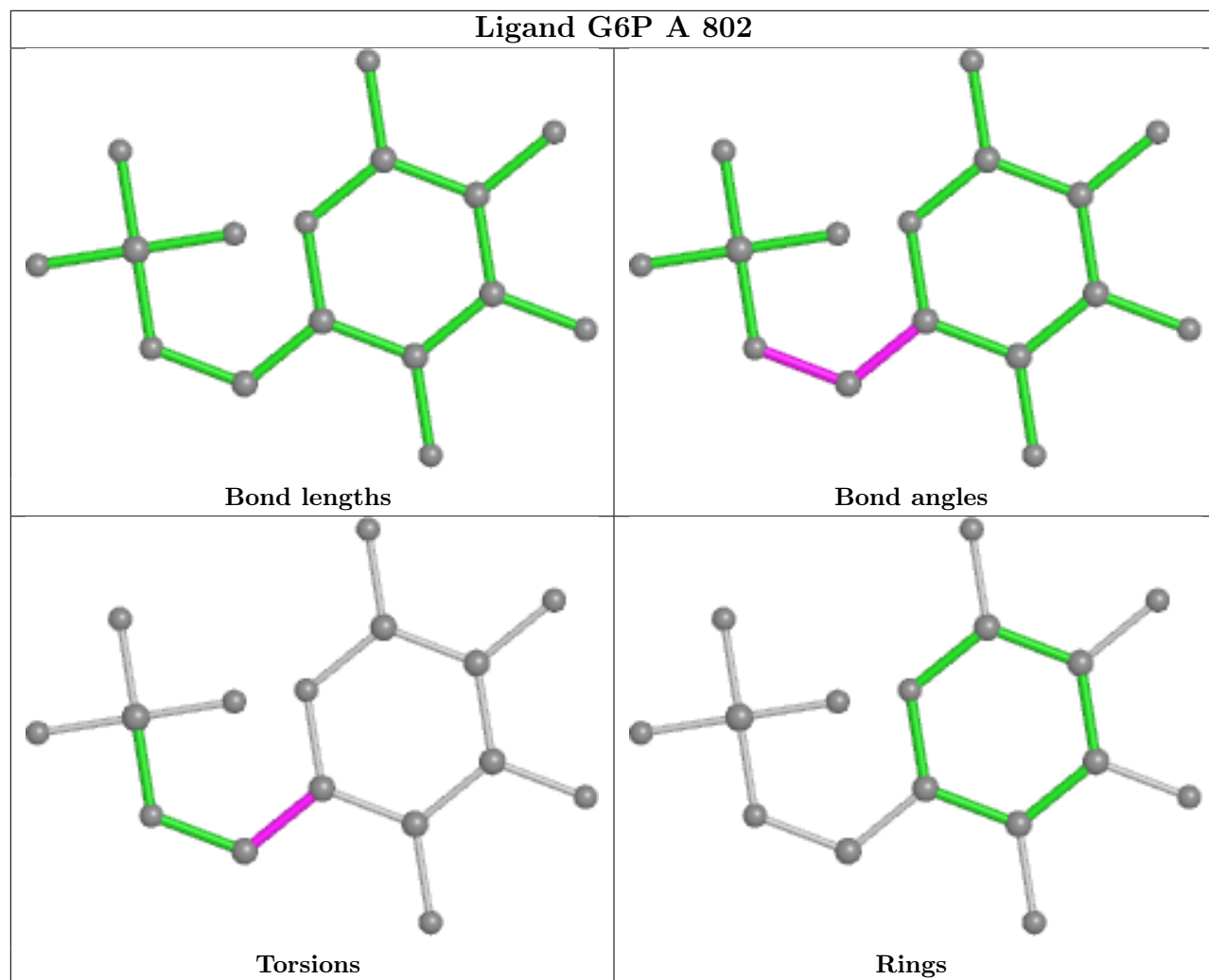
8 monomers are involved in 10 short contacts:

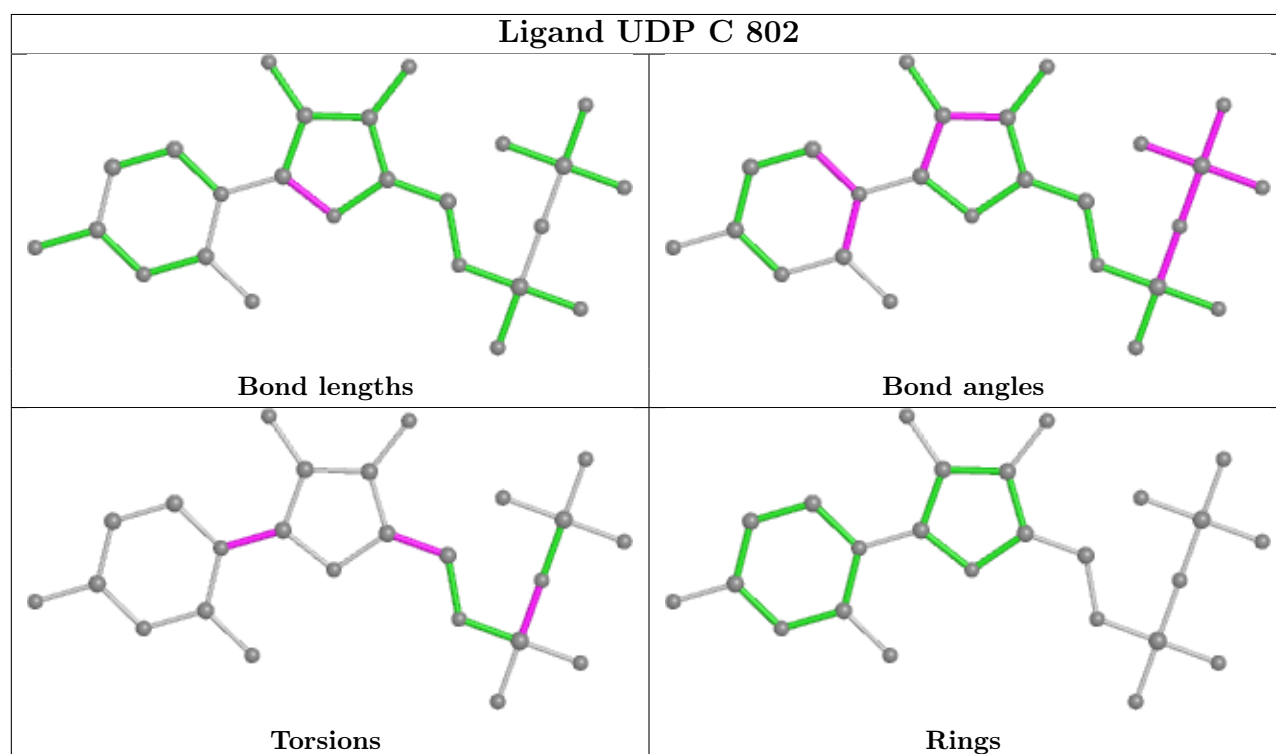
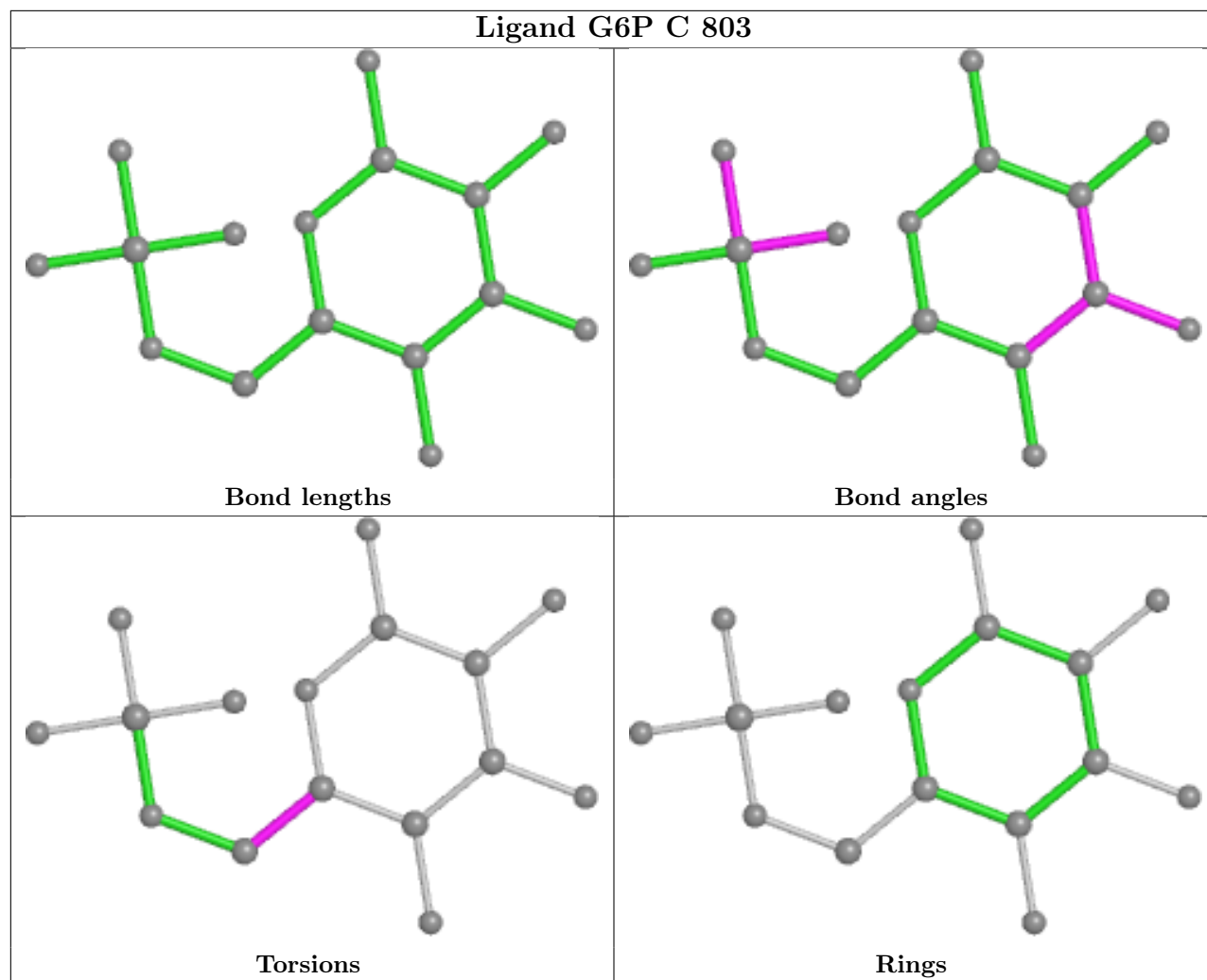
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	802	G6P	1	0
3	D	801	UDP	2	0
4	C	803	G6P	1	0
3	C	802	UDP	1	0
4	B	802	G6P	1	0
2	C	801	GCS	1	0
3	A	801	UDP	1	0
4	D	802	G6P	2	0

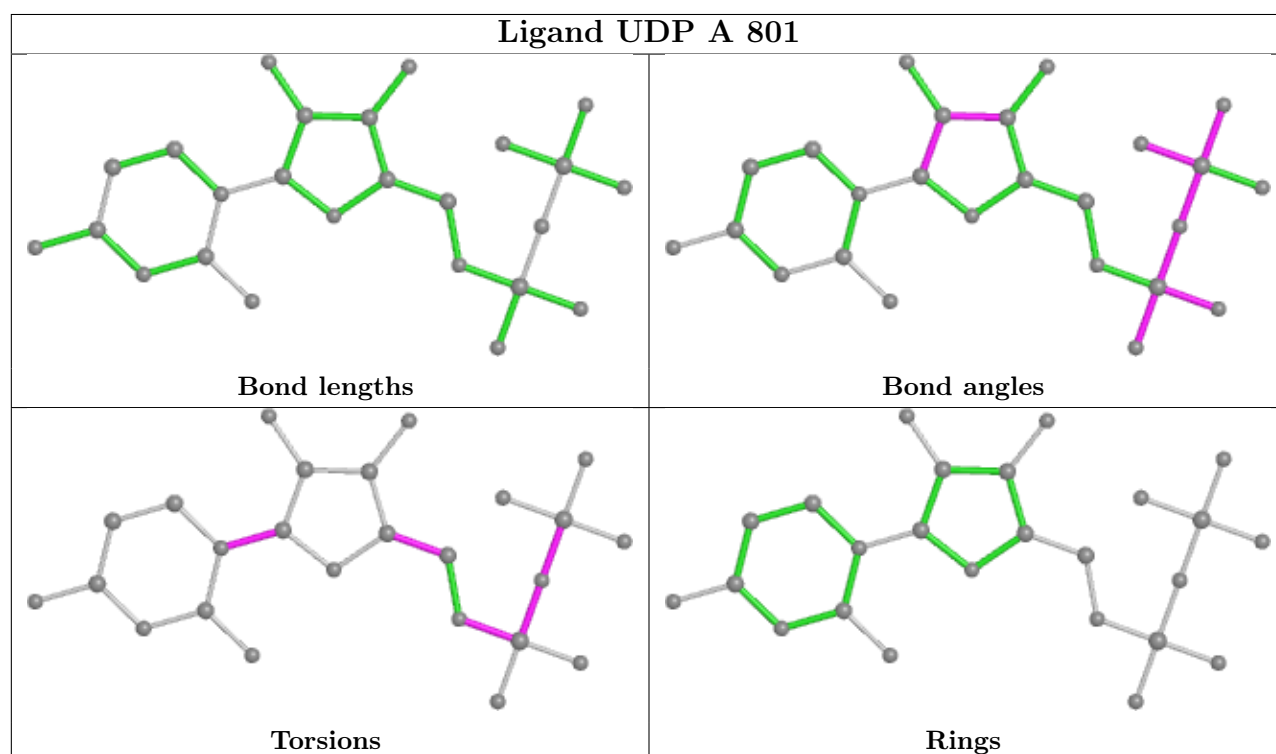
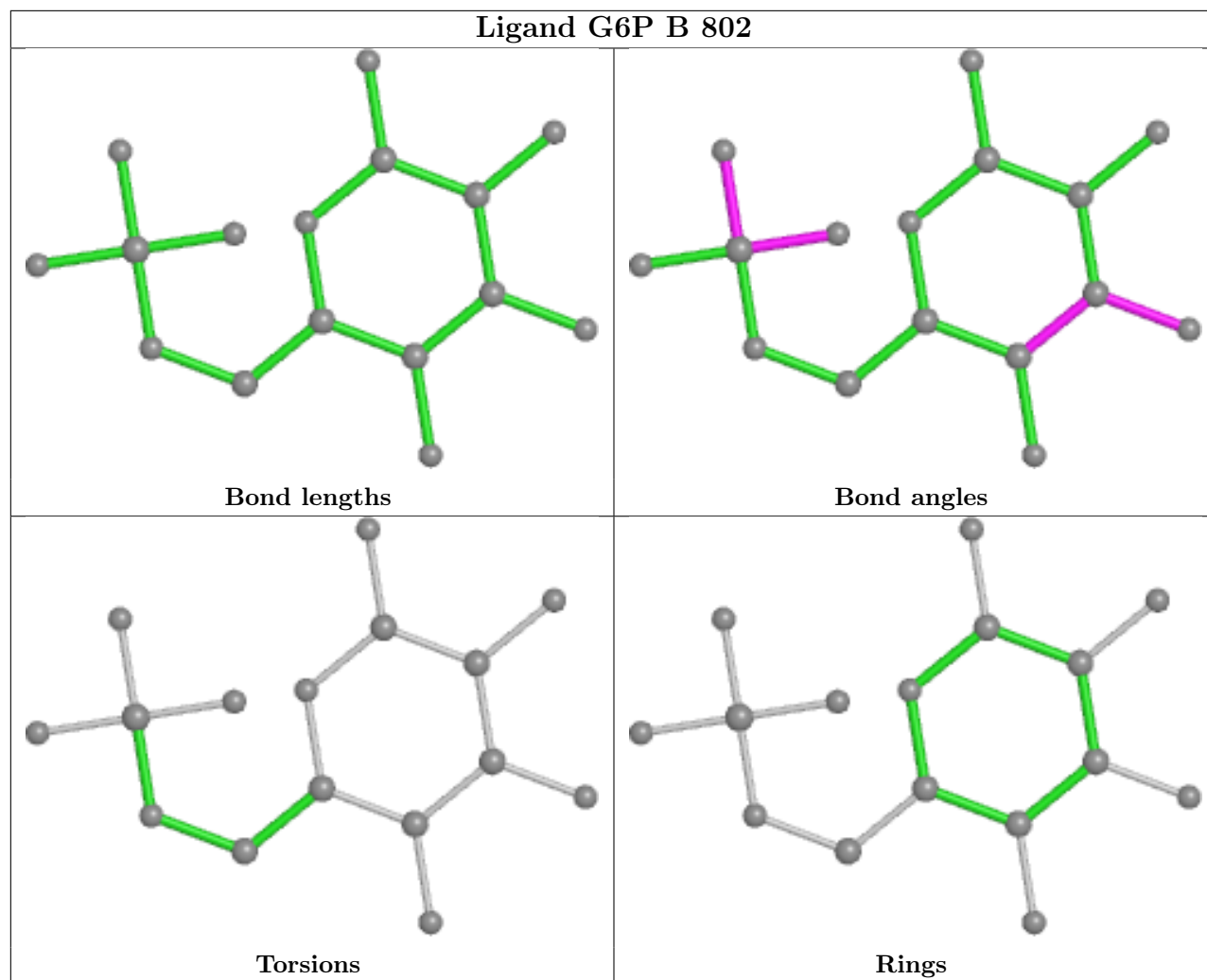
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

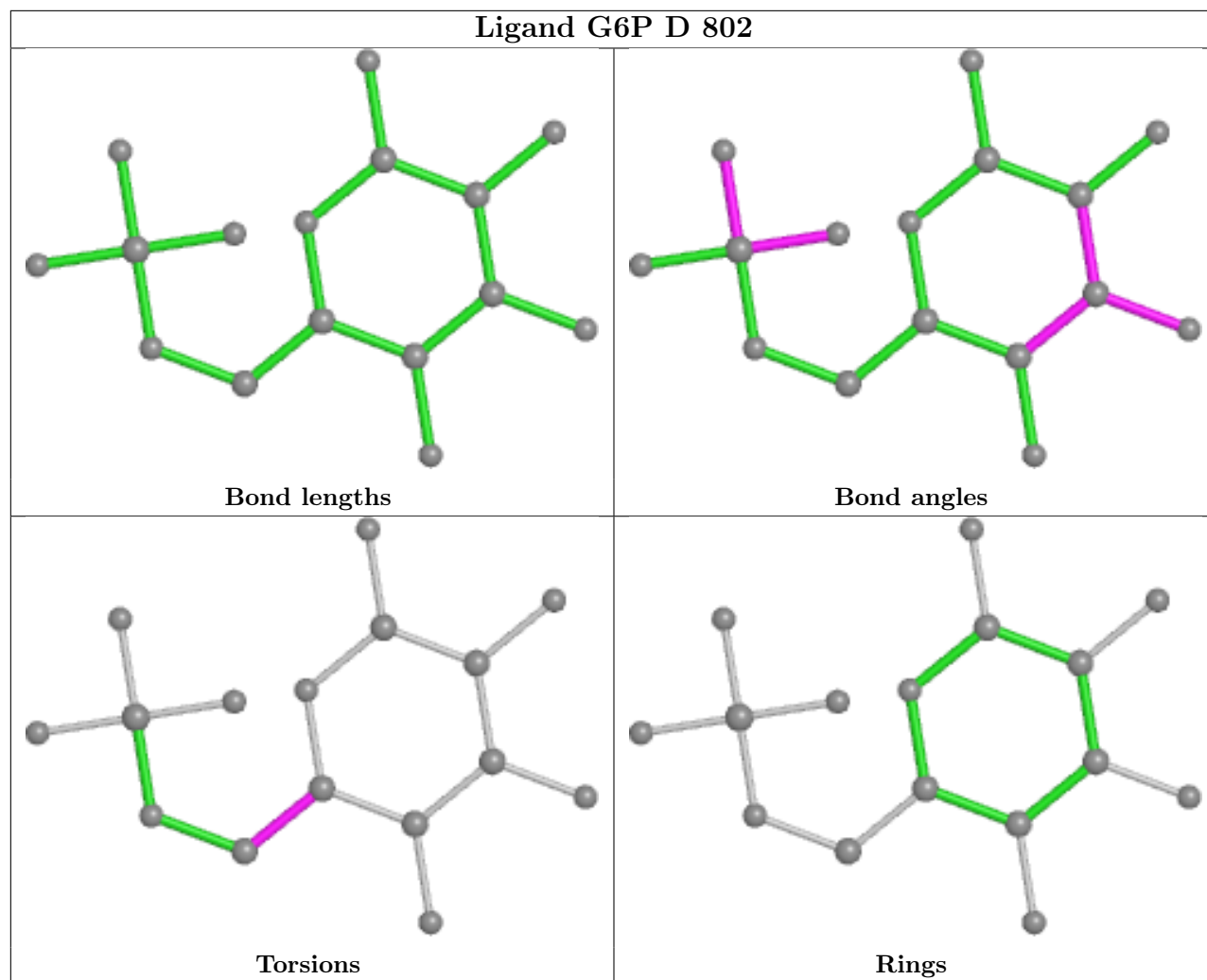
in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	638/720 (88%)	0.20	28 (4%) 34 20	50, 86, 141, 178	0
1	B	636/720 (88%)	-0.02	20 (3%) 49 30	39, 71, 115, 166	0
1	C	638/720 (88%)	0.18	41 (6%) 19 10	45, 80, 129, 169	0
1	D	636/720 (88%)	0.25	44 (6%) 16 9	42, 87, 146, 164	0
All	All	2548/2880 (88%)	0.15	133 (5%) 27 16	39, 80, 138, 178	0

All (133) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	124	SER	5.7
1	D	630	LEU	5.4
1	C	125	LEU	5.1
1	D	2	SER	5.0
1	A	52	ALA	4.9
1	A	96	LEU	4.9
1	D	13	THR	4.9
1	C	52	ALA	4.8
1	D	167	PHE	4.8
1	B	630	LEU	4.6
1	D	620	ASP	4.4
1	D	639	ALA	4.3
1	D	124	SER	4.1
1	D	68	ALA	4.1
1	A	133	ASN	4.0
1	D	628	GLU	4.0
1	D	61	LEU	3.9
1	D	125	LEU	3.9
1	C	525	SER	3.8
1	C	624	GLU	3.7
1	D	619	PRO	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	61	LEU	3.7
1	C	626	VAL	3.7
1	B	166	HIS	3.7
1	B	631	ASN	3.7
1	C	66	PRO	3.6
1	A	53	THR	3.5
1	D	166	HIS	3.5
1	A	135	PHE	3.5
1	A	525	SER	3.4
1	C	13	THR	3.4
1	D	11	PHE	3.4
1	D	629	GLU	3.4
1	A	502	GLY	3.4
1	C	123	TRP	3.3
1	A	56	ASN	3.3
1	B	628	GLU	3.3
1	A	21	VAL	3.3
1	A	627	GLY	3.3
1	D	165	ALA	3.2
1	A	274	ILE	3.2
1	D	622	PHE	3.1
1	C	630	LEU	3.1
1	D	64	LYS	3.1
1	D	6	GLN	3.1
1	A	156	HIS	3.1
1	C	183	ARG	3.1
1	C	544	ASN	3.1
1	C	68	ALA	3.0
1	C	128	ILE	3.0
1	C	526	ILE	3.0
1	D	52	ALA	3.0
1	C	524	PRO	3.0
1	D	133	ASN	3.0
1	A	504	PHE	3.0
1	D	618	TYR	3.0
1	A	13	THR	3.0
1	B	623	ARG	3.0
1	C	165	ALA	3.0
1	C	518	CYS	3.0
1	D	525	SER	3.0
1	C	220	HIS	2.9
1	B	11	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	98	GLU	2.8
1	C	11	PHE	2.8
1	A	625	LEU	2.8
1	B	165	ALA	2.8
1	D	96	LEU	2.8
1	D	4	ASP	2.8
1	A	85	SER	2.8
1	D	85	SER	2.8
1	A	626	VAL	2.8
1	C	129	PRO	2.7
1	A	558	PHE	2.7
1	B	624	GLU	2.7
1	B	10	LEU	2.7
1	C	126	VAL	2.7
1	D	632	ASP	2.7
1	D	274	ILE	2.7
1	D	67	GLU	2.6
1	D	197	LEU	2.6
1	C	166	HIS	2.6
1	D	172	ALA	2.6
1	A	503	VAL	2.6
1	B	13	THR	2.6
1	D	173	GLY	2.6
1	C	623	ARG	2.6
1	C	53	THR	2.6
1	A	78	HIS	2.6
1	B	633	SER	2.5
1	D	159	SER	2.5
1	A	214	GLU	2.5
1	B	499	CYS	2.5
1	D	501	LEU	2.5
1	D	631	ASN	2.5
1	D	502	GLY	2.5
1	D	225	PHE	2.4
1	A	452	ASN	2.4
1	C	167	PHE	2.4
1	C	625	LEU	2.4
1	B	44	HIS	2.4
1	B	46	ILE	2.4
1	C	558	PHE	2.4
1	C	502	GLY	2.4
1	A	524	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	503	VAL	2.3
1	D	117	GLU	2.3
1	A	12	GLU	2.3
1	C	211	ASN	2.3
1	C	627	GLY	2.3
1	B	183	ARG	2.3
1	A	557	ARG	2.3
1	A	514	THR	2.3
1	C	218	VAL	2.3
1	D	120	GLY	2.3
1	D	615	ARG	2.2
1	C	77	GLN	2.2
1	A	317	ILE	2.2
1	D	276	PHE	2.2
1	B	626	VAL	2.2
1	D	209	PHE	2.1
1	B	159	SER	2.1
1	D	638	LEU	2.1
1	C	2	SER	2.1
1	C	181	LYS	2.1
1	D	21	VAL	2.1
1	C	62	ASP	2.1
1	C	116	ASN	2.1
1	C	452	ASN	2.1
1	B	61	LEU	2.1
1	B	156	HIS	2.0
1	B	172	ALA	2.0
1	C	172	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

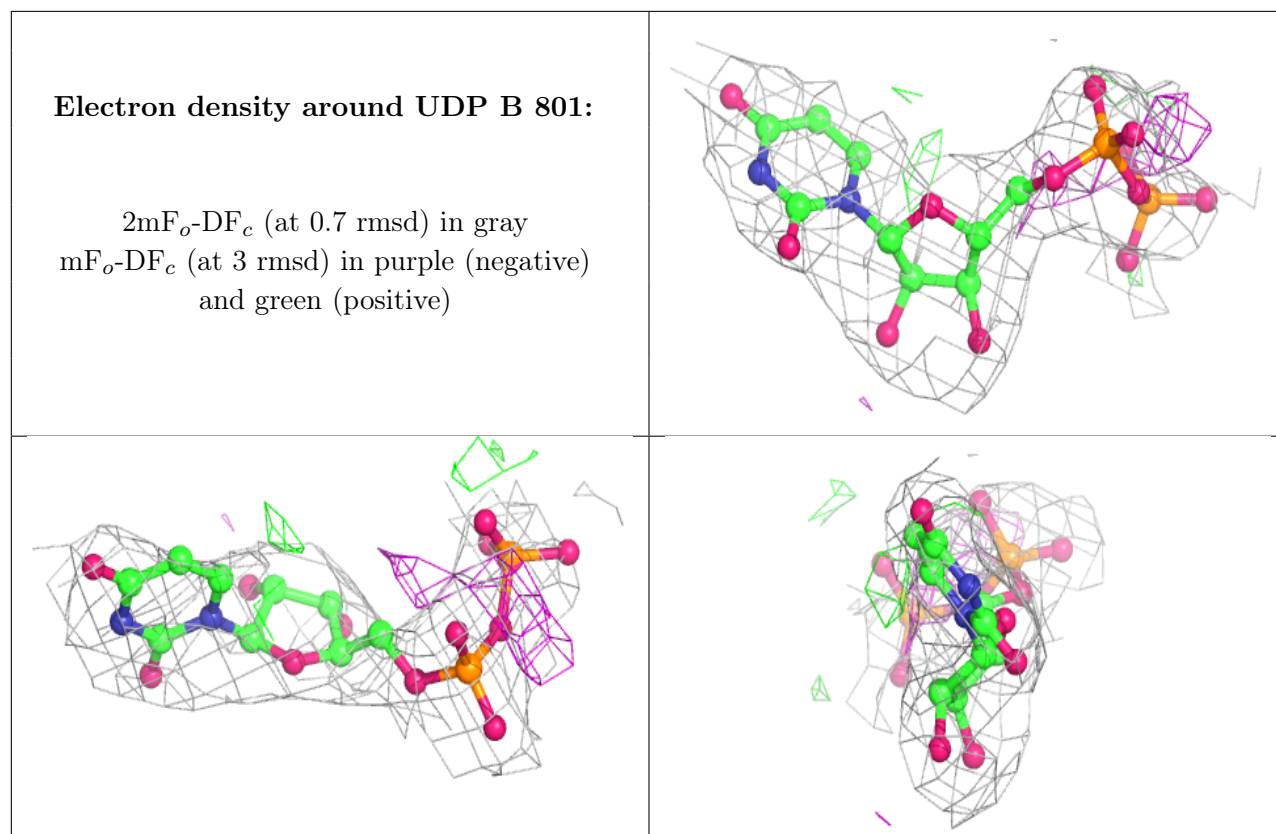
6.4 Ligands [i](#)

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

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

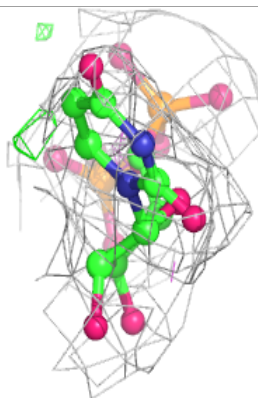
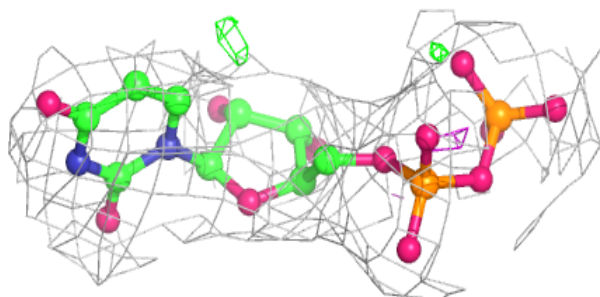
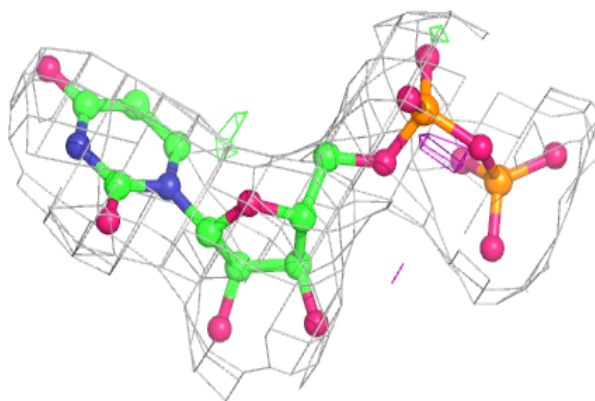
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	UDP	B	801	25/25	0.88	0.20	81,116,172,177	0
3	UDP	A	801	25/25	0.91	0.15	79,94,109,114	0
3	UDP	D	801	25/25	0.92	0.14	80,99,127,130	0
3	UDP	C	802	25/25	0.96	0.16	61,77,83,86	0
4	G6P	B	802	16/16	0.96	0.14	39,49,55,58	0
2	GCS	C	801	12/12	0.97	0.18	60,67,78,78	0
4	G6P	A	802	16/16	0.97	0.12	50,58,65,66	0
4	G6P	C	803	16/16	0.98	0.16	56,66,75,76	0
4	G6P	D	802	16/16	0.98	0.13	43,51,58,63	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.

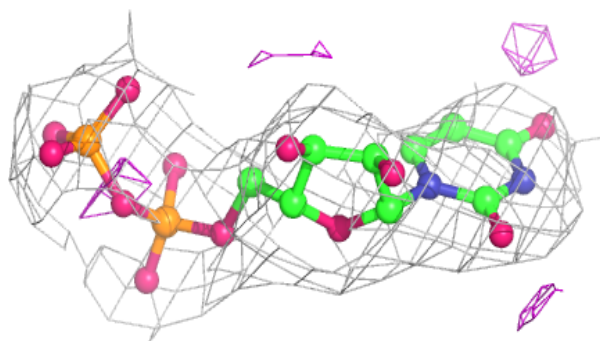
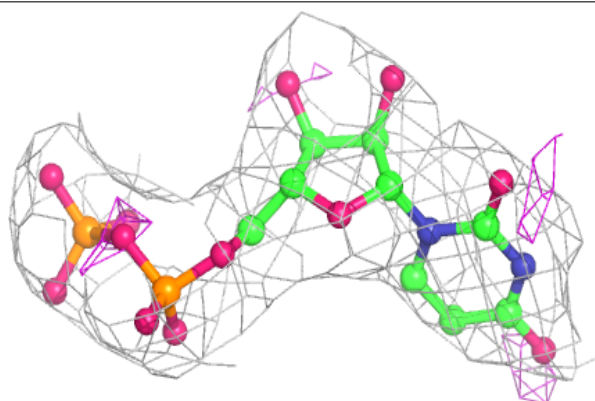


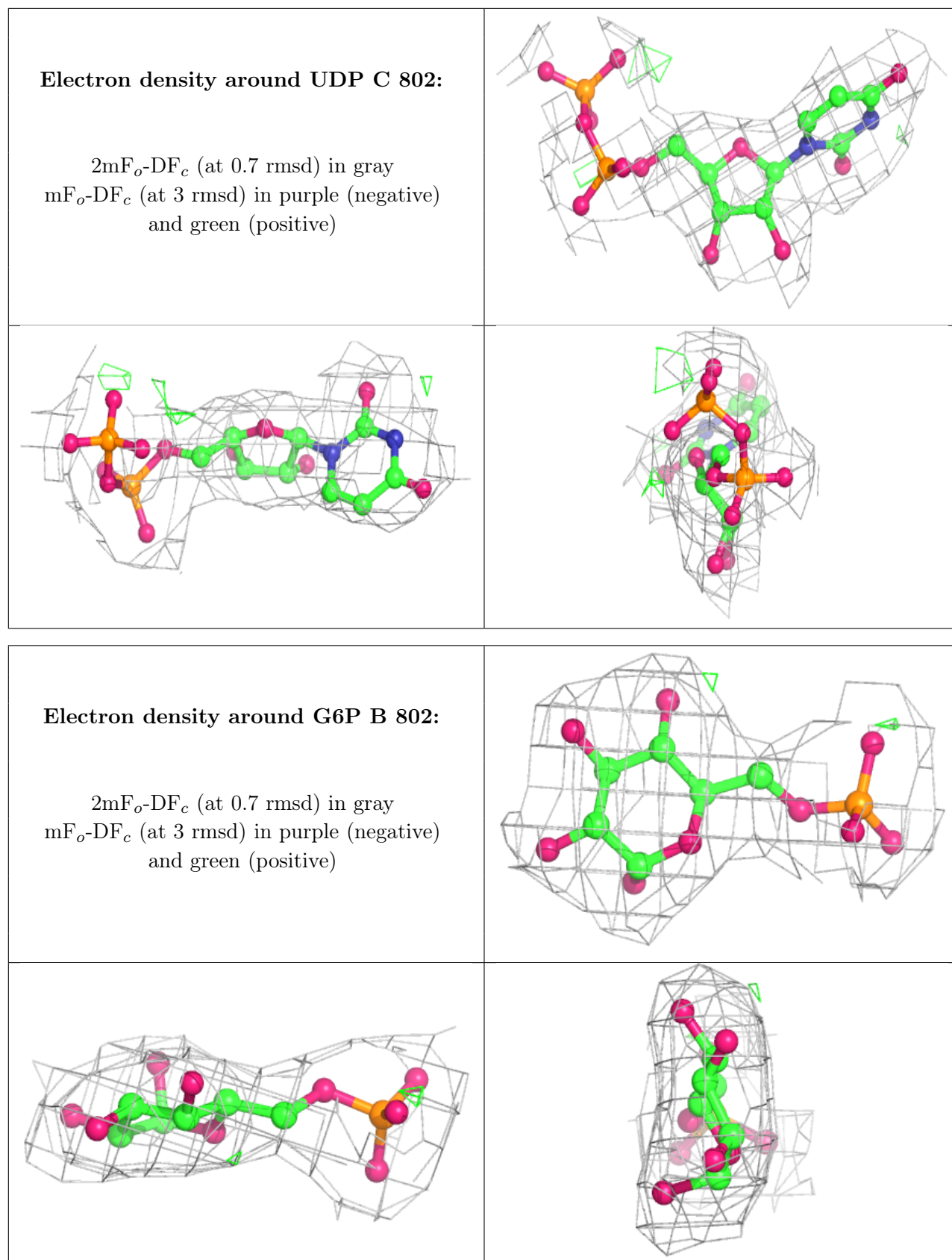
Electron density around UDP A 801:

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

**Electron density around UDP D 801:**

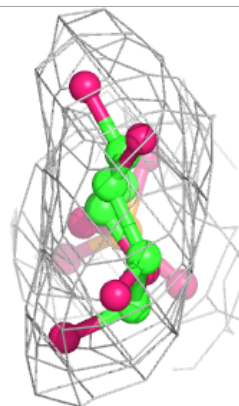
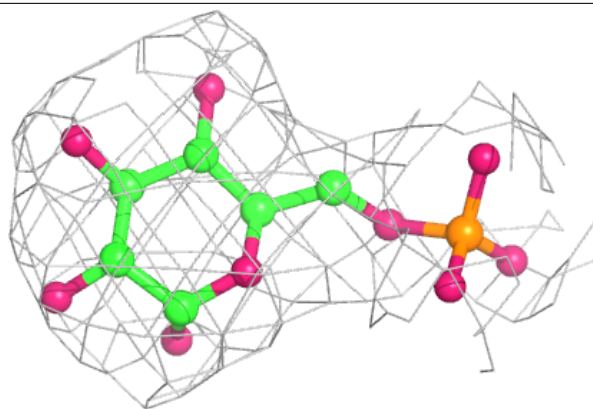
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



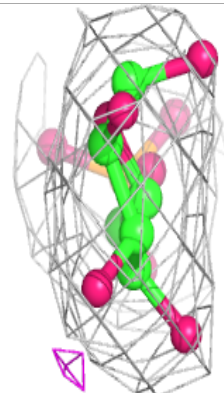
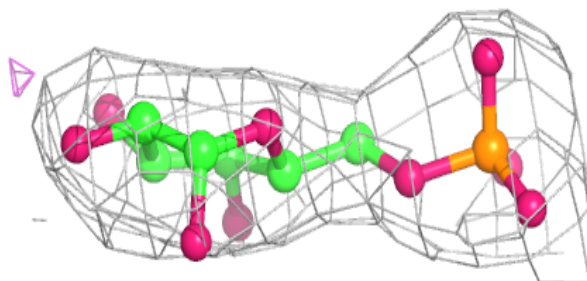
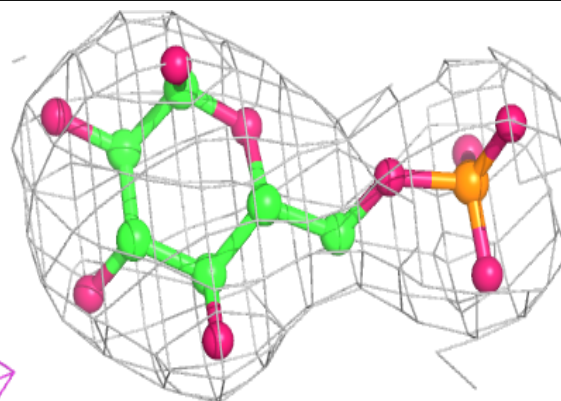


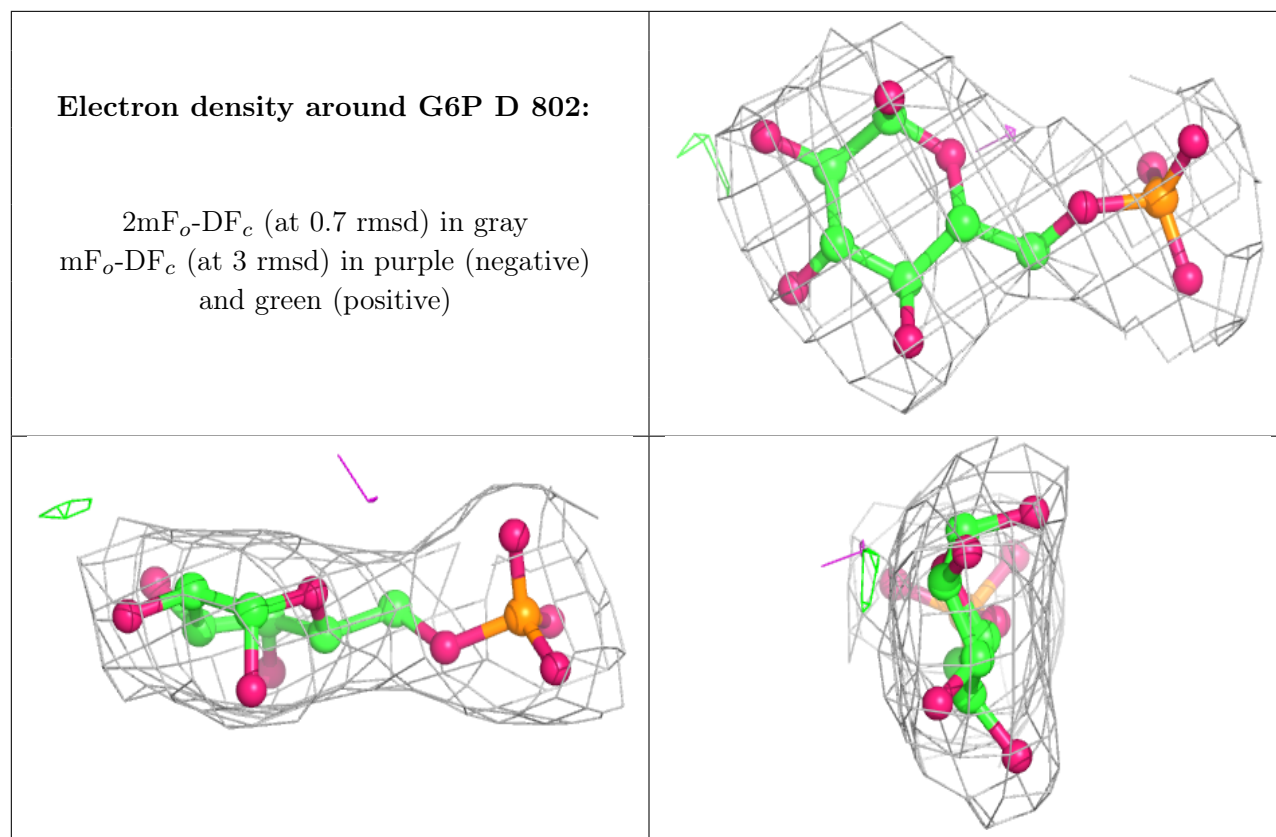
Electron density around G6P A 802:

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

**Electron density around G6P C 803:**

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





6.5 Other polymers [i](#)

There are no such residues in this entry.