



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:33 PM UTC

PDB ID : 5UW4 / pdb_00005uw4
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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

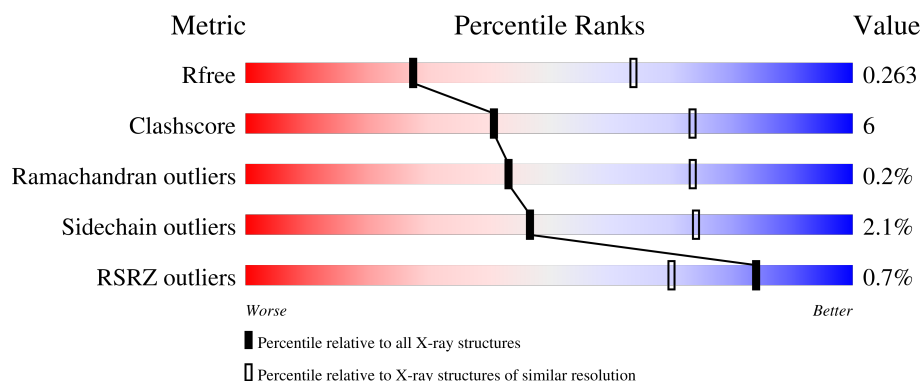
1 Overall quality at a glance

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}	180053	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (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	% 76% 11% • 11%
1	B	720	75% 12% • 12%
1	C	720	75% 12% • 11%
1	D	720	% 73% 14% • 12%

2 Entry composition

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

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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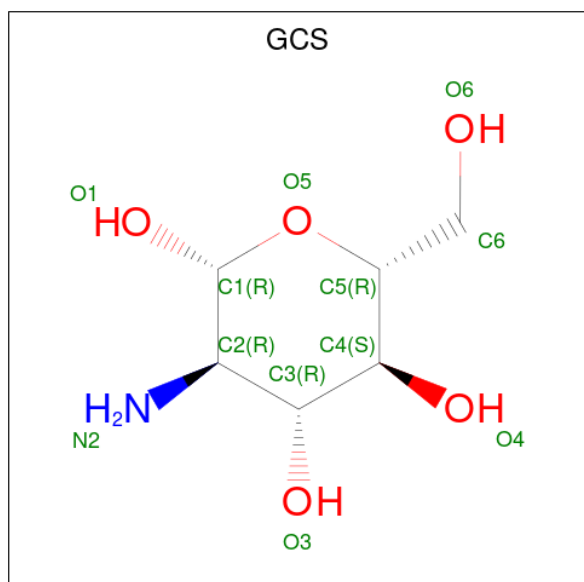
Chain	Residue	Modelled	Actual	Comment	Reference
C	535	SER	ALA	conflict	UNP P27472
B	-19	MET	-	initiating methionine	UNP P27472
B	-18	GLY	-	expression tag	UNP P27472
B	-17	SER	-	expression tag	UNP P27472
B	-16	SER	-	expression tag	UNP P27472
B	-15	HIS	-	expression tag	UNP P27472
B	-14	HIS	-	expression 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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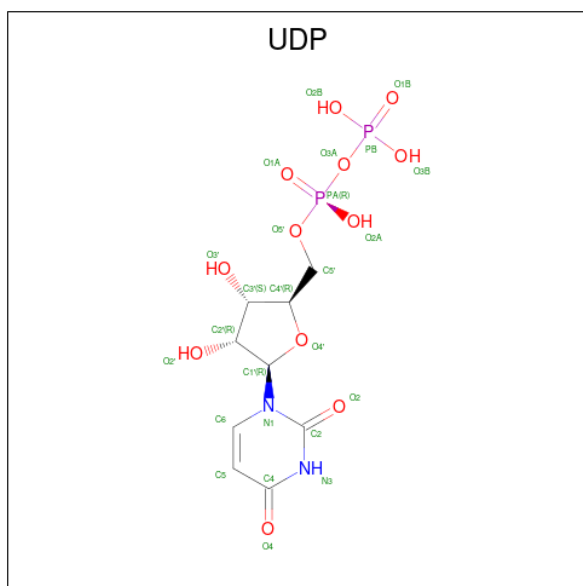
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	expression tag	UNP P27472
A	169	GLN	GLU	engineered mutation	UNP P27472
A	535	SER	ALA	conflict	UNP P27472
D	-19	MET	-	initiating methionine	UNP P27472
D	-18	GLY	-	expression tag	UNP P27472
D	-17	SER	-	expression tag	UNP P27472
D	-16	SER	-	expression 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

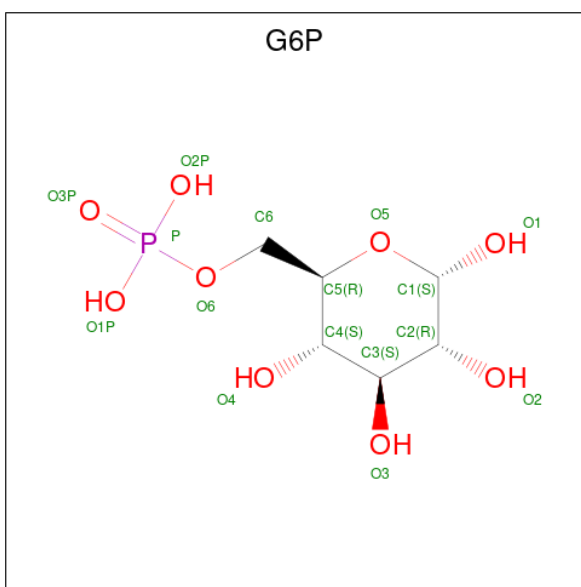
- Molecule 2 is 2-amino-2-deoxy-beta-D-glucopyranose (CCD ID: GCS) (formula: $C_6H_{13}NO_5$).



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

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





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

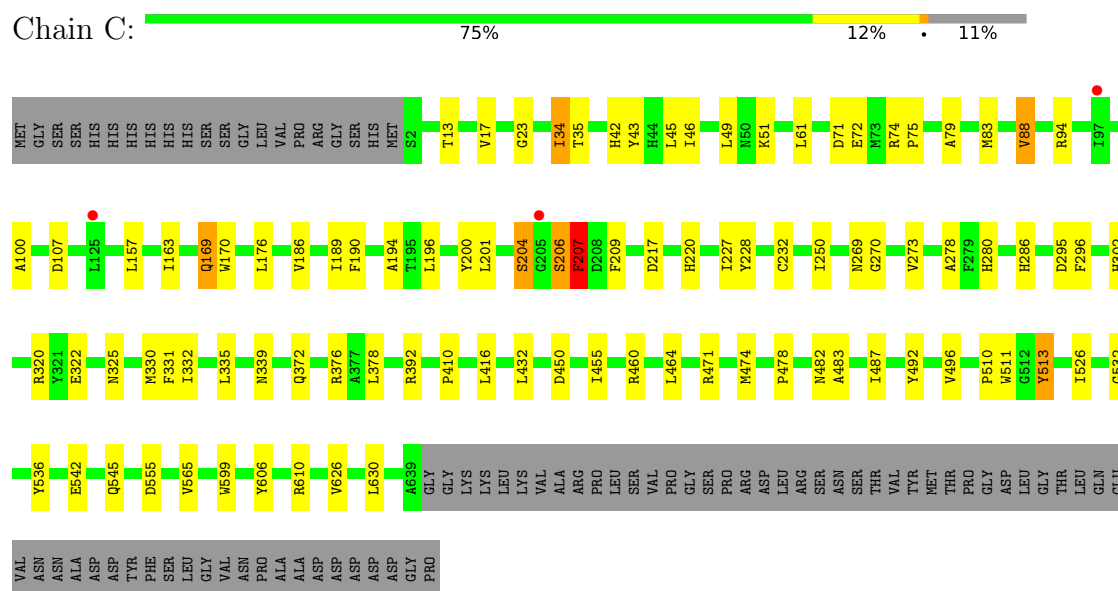
- Molecule 5 is water.

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

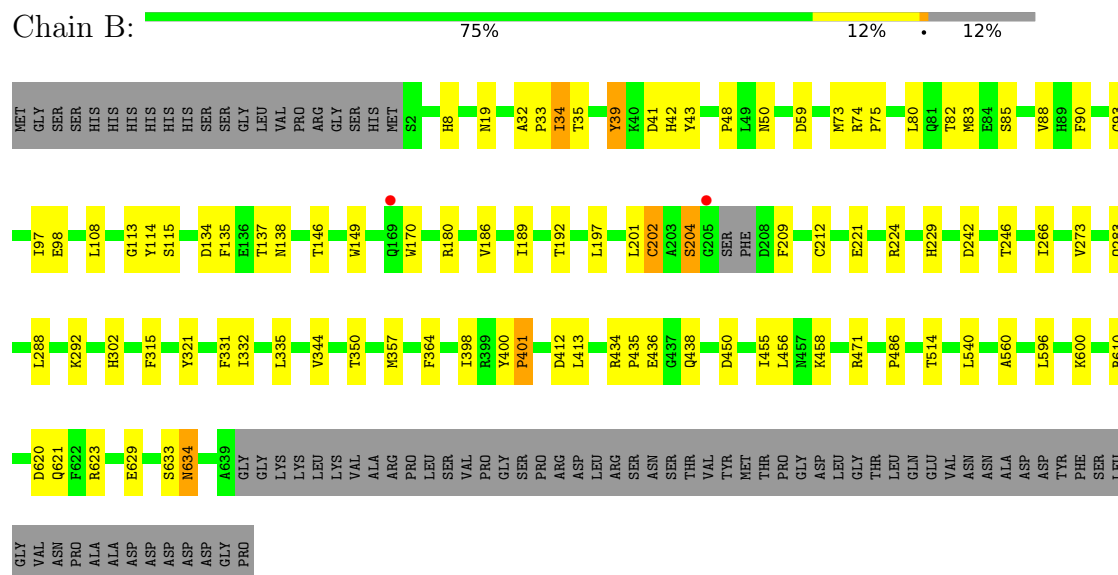
3 Residue-property plots [i](#)

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.

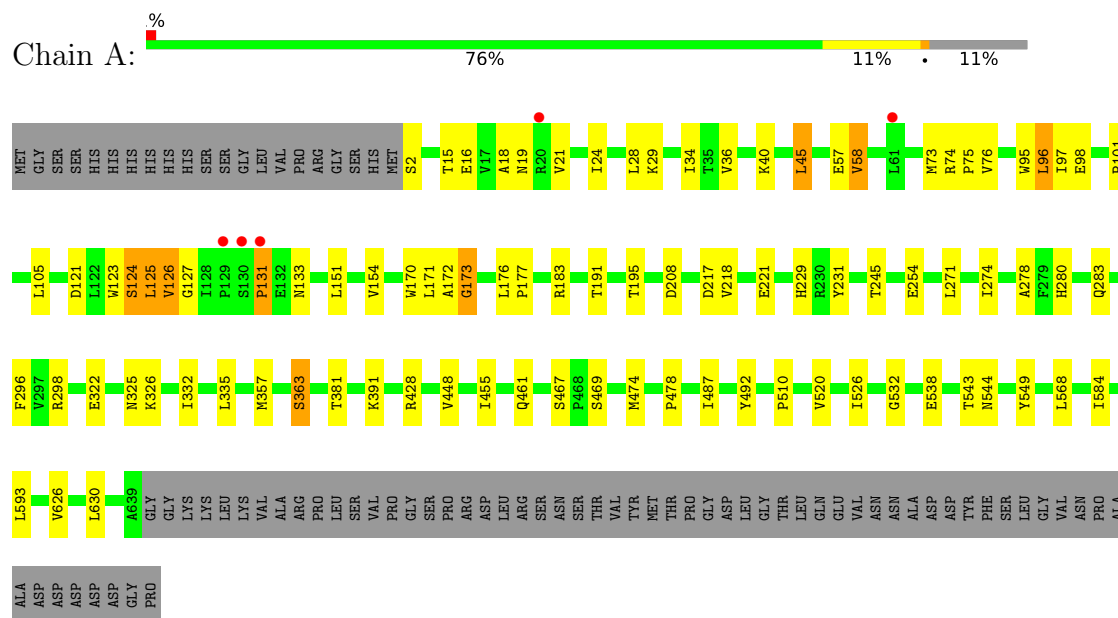
• Molecule 1: Glycogen [starch] synthase isoform 2



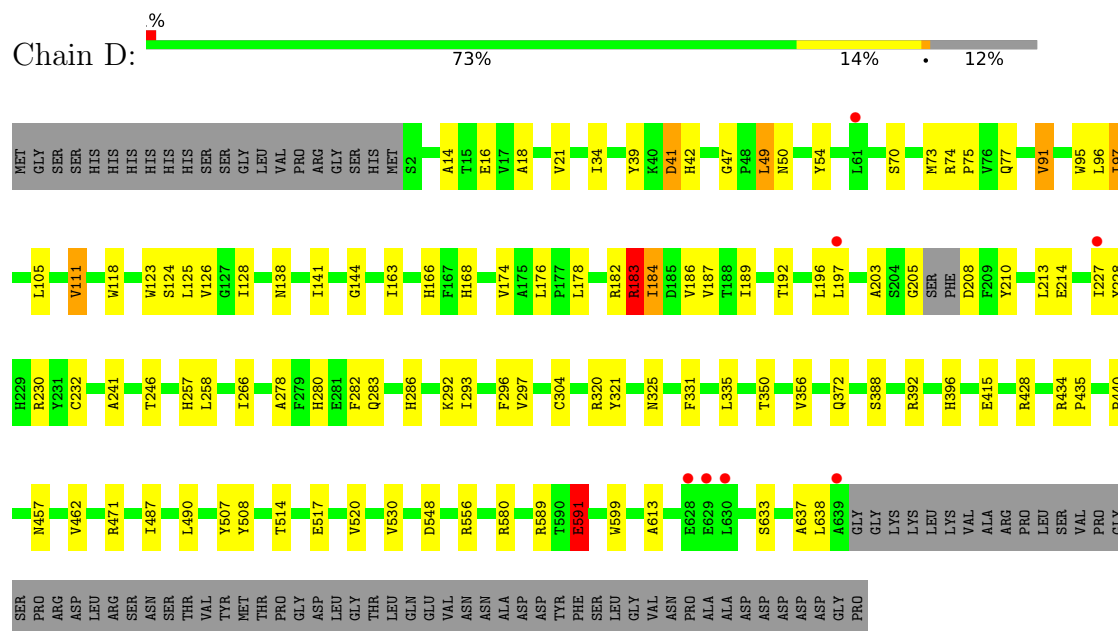
• Molecule 1: Glycogen [starch] synthase isoform 2



- Molecule 1: Glycogen [starch] synthase isoform 2



- Molecule 1: Glycogen [starch] synthase isoform 2



4 Data and refinement statistics

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 – 2.98 145.16 – 2.98	Depositor EDS
% Data completeness (in resolution range)	99.6 (145.16-2.98) 99.6 (145.16-2.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 3.00Å)	Xtriage
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 (5.40%)	wwPDB-VP
Wilson B-factor (Å ²)	74.6	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.003 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20374	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage'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	0.79	0/5101	1.02	8/6943 (0.1%)
1	B	0.81	0/5206	1.07	15/7061 (0.2%)
1	C	0.79	0/5251	1.04	11/7119 (0.2%)
1	D	0.79	1/5123 (0.0%)	1.04	7/6962 (0.1%)
All	All	0.80	1/20681 (0.0%)	1.04	41/28085 (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	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	124	SER	CA-C	6.48	1.61	1.52

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	125	LEU	N-CA-C	13.19	127.28	111.33
1	C	206	SER	N-CA-C	12.34	124.47	111.14
1	B	202	CYS	N-CA-C	-8.55	101.77	108.78
1	C	207	PHE	N-CA-C	8.40	120.86	108.60
1	A	131	PRO	N-CA-C	-8.12	95.73	112.47
1	B	170	TRP	N-CA-C	7.82	120.79	111.33
1	B	32	ALA	N-CA-C	7.49	123.79	113.16
1	C	17	VAL	N-CA-C	7.43	117.56	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	34	ILE	N-CA-CB	7.13	125.78	111.24
1	D	183	ARG	N-CA-C	6.91	120.93	111.39
1	A	124	SER	CA-C-N	6.82	129.91	120.63
1	A	124	SER	C-N-CA	6.82	129.91	120.63
1	D	124	SER	N-CA-C	6.79	125.26	110.80
1	A	124	SER	O-C-N	6.70	132.20	122.43
1	C	478	PRO	N-CA-C	6.57	121.34	112.48
1	C	372	GLN	N-CA-C	-6.46	103.86	111.03
1	B	33	PRO	CB-CA-C	-6.43	102.30	111.68
1	B	212	CYS	N-CA-C	6.33	120.46	112.87
1	C	88	VAL	N-CA-C	5.87	117.03	108.58
1	D	203	ALA	N-CA-C	-5.86	100.28	109.25
1	B	344	VAL	N-CA-C	5.78	116.43	110.36
1	A	96	LEU	N-CA-C	5.75	119.61	110.70
1	A	58	VAL	N-CA-C	5.68	116.30	108.12
1	C	478	PRO	CB-CA-C	-5.55	105.62	112.89
1	B	560	ALA	CA-C-N	-5.54	113.19	119.28
1	B	560	ALA	C-N-CA	-5.54	113.19	119.28
1	C	204	SER	O-C-N	-5.49	116.18	122.93
1	D	633	SER	N-CA-C	5.43	119.53	111.87
1	B	634	ASN	CB-CA-C	-5.43	100.53	110.62
1	A	126	VAL	N-CA-C	5.43	120.62	109.34
1	B	39	TYR	N-CA-C	5.38	120.71	114.04
1	B	93	GLY	N-CA-C	5.31	118.46	110.18
1	C	471	ARG	CG-CD-NE	5.25	123.55	112.00
1	A	173	GLY	N-CA-C	5.24	125.60	113.18
1	B	315	PHE	N-CA-C	-5.24	101.74	110.17
1	D	184	ILE	N-CA-C	5.23	116.76	109.80
1	B	209	PHE	N-CA-C	5.18	117.33	111.11
1	D	591	GLU	N-CA-C	5.14	116.57	111.07
1	C	555	ASP	N-CA-C	5.13	116.66	108.41
1	C	209	PHE	N-CA-C	5.09	117.49	111.33
1	B	59	ASP	N-CA-C	-5.07	102.55	110.10

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 ⓘ

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	65	0
1	B	5084	0	4955	57	0
1	C	5127	0	5015	69	0
1	D	5005	0	4827	72	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	258	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 (258) 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
1:C:217:ASP:CG	1:C:220:HIS:CD2	2.20	1.19
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:LEU:HD12	1:D:50:ASN:N	1.88	0.88
1:C:217:ASP:CG	1:C:220:HIS:HD2	1.66	0.87
1:D:41:ASP:OD1	1:D:73:MET:HG3	1.75	0.85
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:D:49:LEU:HD12	1:D:49:LEU:C	2.19	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:B:114:TYR:O	1:B:115:SER:C	2.42	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:202:CYS:C	1:B:204:SER:H	2.09	0.60
1:B:80:LEU:HD22	1:B:90:PHE:CE2	2.37	0.60
1:B:398:ILE:CD1	1:A:298:ARG:HD3	2.31	0.60
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:A:170:TRP:O	1:A:171:LEU:C	2.46	0.58
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:270:GLY:O	1:C:513:TYR:OH	2.20	0.57
1:D:41:ASP:CG	1:D:73:MET:HG3	2.29	0.56
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: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:B:435:PRO:O	1:B:436:GLU:C	2.47	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: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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:74:ARG:O	1:A:75:PRO:C	2.53	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: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:B:202:CYS:C	1:B:204:SER:N	2.65	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:A:231:TYR:CD1	1:A:231:TYR:C	2.89	0.50
1:C:335:LEU:HD13	1:C:474:MET:HE1	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:B:620:ASP:CG	1:B:621:GLN:N	2.68	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:335:LEU:CD1	1:A:474:MET:HE1	2.42	0.48
1:A:170:TRP:CH2	1:A:171:LEU:HD23	2.48	0.48
1:A:335:LEU:HD13	1:A:474:MET:HE1	1.94	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:B:42:HIS:CD2	1:B:73:MET:HE1	2.49	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:B:412:ASP:C	1:B:413:LEU:O	2.54	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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:D:49:LEU:C	1:D:49:LEU:CD1	2.85	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:C:227:ILE:O	1:C:228:TYR:C	2.59	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:C:217:ASP:CG	1:C:220:HIS:NE2	2.71	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: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:83:MET:HE2	1:C:88:VAL:HG21	1.97	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:113:GLY:O	1:B:114:TYR:C	2.60	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:195:THR:OG1	1:A:254:GLU:OE1	2.29	0.45
1:B:620:ASP:CG	1:B:621:GLN:HG3	2.40	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:D:321:TYR:HD2	1:D:321:TYR:C	2.25	0.44
1:C:536:TYR:CD1	1:C:536:TYR:C	2.95	0.44
1:B:192:THR:CG2	1:B:246:THR:HG22	2.47	0.44
1:B:357:MET:HE3	1:B:455:ILE:HD11	1.98	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:D:637:ALA:C	1:D:638:LEU:HD12	2.43	0.43
1:B:8:HIS:NE2	1:B:39:TYR:OH	2.42	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: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:D:321:TYR:C	1:D:321:TYR:CD2	2.94	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: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:97:ILE:O	1:B:98:GLU:C	2.61	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:41:ASP:OD1	1:B:73:MET:HE3	2.20	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:83:MET:HE2	1:B:88:VAL:HG11	2.02	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
1:C:410:PRO:HG2	1:C:416:LEU:HD21	2.02	0.41
1:C:335:LEU:CD1	1:C:474:MET:HE1	2.50	0.41
1:A:208:ASP:OD1	1:A:208:ASP:C	2.64	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:B:357:MET:HE2	1:B:455:ILE:HD11	2.00	0.41
1:A:19:ASN:OD1	1:A:19:ASN:N	2.54	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:B:288:LEU:C	1:B:288:LEU:HD23	2.46	0.41
1:D:178:LEU:O	1:D:182:ARG:CB	2.69	0.41
1:A:626:VAL:O	1:A:626:VAL:HG12	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: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%)	24	58
1	B	632/720 (88%)	585 (93%)	46 (7%)	1 (0%)	43	73
1	C	636/720 (88%)	586 (92%)	50 (8%)	0	100	100
1	D	632/720 (88%)	579 (92%)	52 (8%)	1 (0%)	43	73
All	All	2536/2880 (88%)	2325 (92%)	206 (8%)	5 (0%)	43	73

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%)	502 (97%)	14 (3%)	39	70
1	B	539/621 (87%)	533 (99%)	6 (1%)	65	82
1	C	547/621 (88%)	540 (99%)	7 (1%)	61	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	D	523/621 (84%)	505 (97%)	18 (3%)	32 64
All	All	2125/2484 (86%)	2080 (98%)	45 (2%)	47 74

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	34	ILE
1	C	49	LEU
1	C	169	GLN
1	C	207	PHE
1	C	250	ILE
1	C	513	TYR
1	C	526	ILE
1	B	186	VAL
1	B	204	SER
1	B	266	ILE
1	B	273	VAL
1	B	292	LYS
1	B	514	THR
1	A	2	SER
1	A	34	ILE
1	A	45	LEU
1	A	58	VAL
1	A	124	SER
1	A	125	LEU
1	A	274	ILE
1	A	326	LYS
1	A	363	SER
1	A	391	LYS
1	A	448	VAL
1	A	461	GLN
1	A	467	SER
1	A	469	SER
1	D	16	GLU
1	D	21	VAL
1	D	41	ASP
1	D	49	LEU
1	D	91	VAL
1	D	97	ILE
1	D	111	VAL
1	D	123	TRP
1	D	183	ARG

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Mol	Chain	Res	Type
1	D	184	ILE
1	D	196	LEU
1	D	266	ILE
1	D	440	PRO
1	D	457	ASN
1	D	462	VAL
1	D	517	GLU
1	D	520	VAL
1	D	591	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	38	GLN
1	C	44	HIS
1	C	160	GLN
1	C	220	HIS
1	C	362	ASN
1	C	372	GLN
1	B	19	ASN
1	B	42	HIS
1	B	160	GLN
1	B	211	ASN
1	B	402	HIS
1	B	403	ASN
1	B	452	ASN
1	B	621	GLN
1	B	634	ASN
1	A	257	HIS
1	A	325	ASN
1	A	461	GLN
1	D	283	GLN
1	D	325	ASN
1	D	372	GLN
1	D	457	ASN
1	D	582	GLN

5.3.3 RNA ⓘ

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 oligosaccharides 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	$\# Z > 2$
4	G6P	C	803	-	16,16,16	0.78	0	23,24,24	1.33	4 (17%)
4	G6P	B	802	-	16,16,16	0.60	0	23,24,24	1.29	2 (8%)
2	GCS	C	801	-	12,12,12	0.73	0	16,17,17	1.62	4 (25%)
3	UDP	C	802	-	25,26,26	1.19	4 (16%)	38,40,40	1.79	9 (23%)
3	UDP	B	801	-	25,26,26	1.55	4 (16%)	38,40,40	1.67	5 (13%)
3	UDP	D	801	-	25,26,26	1.25	4 (16%)	38,40,40	1.69	7 (18%)
4	G6P	A	802	-	16,16,16	0.71	0	23,24,24	1.21	1 (4%)
4	G6P	D	802	-	16,16,16	0.64	0	23,24,24	1.40	4 (17%)
3	UDP	A	801	-	25,26,26	1.21	5 (20%)	38,40,40	1.78	6 (15%)

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
4	G6P	C	803	-	-	0/6/26/26	0/1/1/1
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	C	802	-	-	4/16/32/32	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	UDP	B	801	-	-	5/16/32/32	0/2/2/2
3	UDP	D	801	-	-	1/16/32/32	0/2/2/2
4	G6P	A	802	-	-	2/6/26/26	0/1/1/1
4	G6P	D	802	-	-	1/6/26/26	0/1/1/1
3	UDP	A	801	-	-	6/16/32/32	0/2/2/2

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	801	UDP	PA-O3A	5.60	1.65	1.59
3	D	801	UDP	PA-O3A	3.33	1.63	1.59
3	B	801	UDP	C2-N1	2.74	1.42	1.38
3	A	801	UDP	C4-N3	-2.62	1.34	1.38
3	A	801	UDP	C2-N1	2.59	1.42	1.38
3	B	801	UDP	C6-C5	2.55	1.41	1.35
3	D	801	UDP	C6-C5	2.39	1.40	1.35
3	C	802	UDP	C2-N1	2.31	1.42	1.38
3	D	801	UDP	C2-N1	2.25	1.42	1.38
3	D	801	UDP	C4-N3	-2.25	1.34	1.38
3	C	802	UDP	C6-C5	2.20	1.40	1.35
3	B	801	UDP	C4-N3	-2.16	1.34	1.38
3	A	801	UDP	C2-N3	-2.15	1.34	1.38
3	C	802	UDP	C4-N3	-2.15	1.34	1.38
3	A	801	UDP	C6-C5	2.11	1.40	1.35
3	A	801	UDP	PA-O3A	2.08	1.61	1.59
3	C	802	UDP	PA-O3A	2.03	1.61	1.59

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	UDP	N3-C2-N1	5.61	122.19	114.89
3	C	802	UDP	N3-C2-N1	5.28	121.77	114.89
3	B	801	UDP	N3-C2-N1	5.28	121.77	114.89
3	A	801	UDP	C4-N3-C2	-5.22	120.13	126.61
3	B	801	UDP	C4-N3-C2	-5.04	120.35	126.61
3	D	801	UDP	N3-C2-N1	5.04	121.45	114.89
3	D	801	UDP	C4-N3-C2	-4.74	120.72	126.61
3	C	802	UDP	C4-N3-C2	-4.37	121.18	126.61
2	C	801	GCS	O5-C1-C2	4.03	114.21	109.51
3	D	801	UDP	C5-C4-N3	3.62	119.86	114.80
4	D	802	G6P	O3-C3-C2	-3.49	102.15	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801	UDP	C5-C4-N3	3.40	119.56	114.80
3	A	801	UDP	C5-C4-N3	3.26	119.36	114.80
3	C	802	UDP	C5-C4-N3	3.02	119.03	114.80
3	C	802	UDP	C6-N1-C2	-2.96	117.40	121.00
3	C	802	UDP	O4'-C1'-N1	2.87	114.87	108.36
4	B	802	G6P	O2P-P-O1P	2.86	118.51	107.80
3	D	801	UDP	O4-C4-C5	-2.81	120.31	125.16
3	A	801	UDP	O3B-PB-O2B	2.78	118.24	107.80
4	C	803	G6P	O2P-P-O1P	2.61	117.57	107.80
3	C	802	UDP	O4-C4-C5	-2.57	120.73	125.16
4	D	802	G6P	O3-C3-C4	-2.53	104.41	110.38
3	A	801	UDP	O2-C2-N3	-2.46	116.96	121.49
4	C	803	G6P	C4-C3-C2	2.44	115.11	110.83
3	D	801	UDP	O2-C2-N1	-2.43	119.63	122.80
2	C	801	GCS	C3-C2-N2	-2.42	106.10	111.05
4	D	802	G6P	C4-C3-C2	2.39	115.03	110.83
3	C	802	UDP	O3A-PB-O1B	-2.32	98.81	111.04
3	B	801	UDP	O2-C2-N1	-2.31	119.79	122.80
4	B	802	G6P	O3-C3-C4	-2.29	104.98	110.38
3	D	801	UDP	C3'-C2'-C1'	2.25	105.72	101.46
4	C	803	G6P	O3-C3-C2	-2.23	105.12	110.38
3	B	801	UDP	O4-C4-C5	-2.19	121.39	125.16
3	D	801	UDP	C6-N1-C2	-2.16	118.37	121.00
3	A	801	UDP	O2A-PA-O1A	2.14	122.42	112.44
4	A	802	G6P	O6-C6-C5	2.05	115.96	108.99
3	C	802	UDP	O3B-PB-O2B	2.03	115.43	107.80
2	C	801	GCS	C1-C2-C3	-2.03	107.93	110.60
4	D	802	G6P	O2P-P-O1P	2.02	115.39	107.80
3	C	802	UDP	O2-C2-N1	-2.02	120.16	122.80
4	C	803	G6P	O5-C1-C2	-2.02	106.75	110.30
2	C	801	GCS	O1-C1-O5	-2.00	104.47	110.41

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	801	UDP	PA-O3A-PB-O3B
3	A	801	UDP	C5'-O5'-PA-O1A
4	A	802	G6P	O5-C5-C6-O6
3	B	801	UDP	C3'-C4'-C5'-O5'
3	B	801	UDP	O4'-C4'-C5'-O5'
3	C	802	UDP	O4'-C4'-C5'-O5'

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

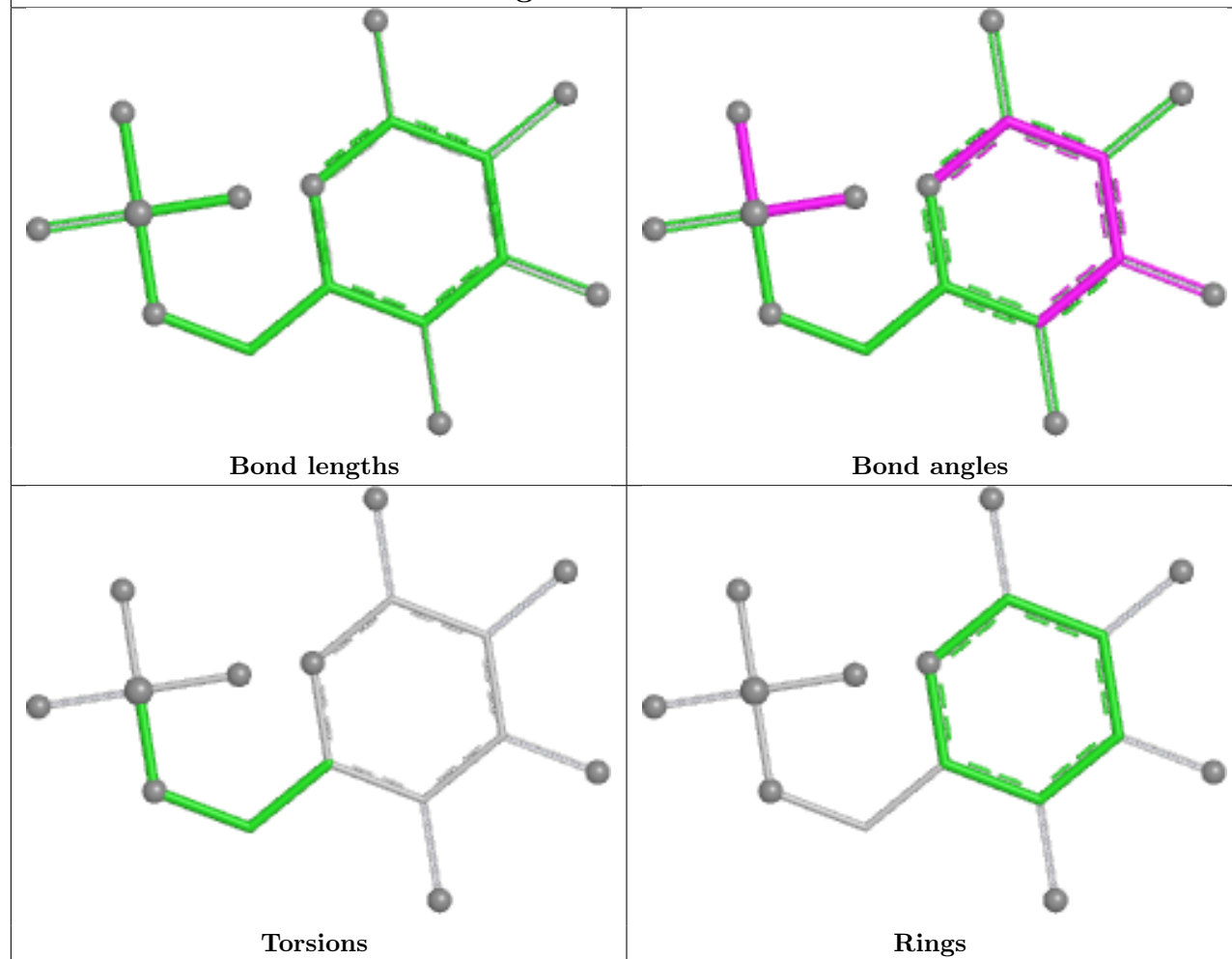
There are no ring outliers.

8 monomers are involved in 10 short contacts:

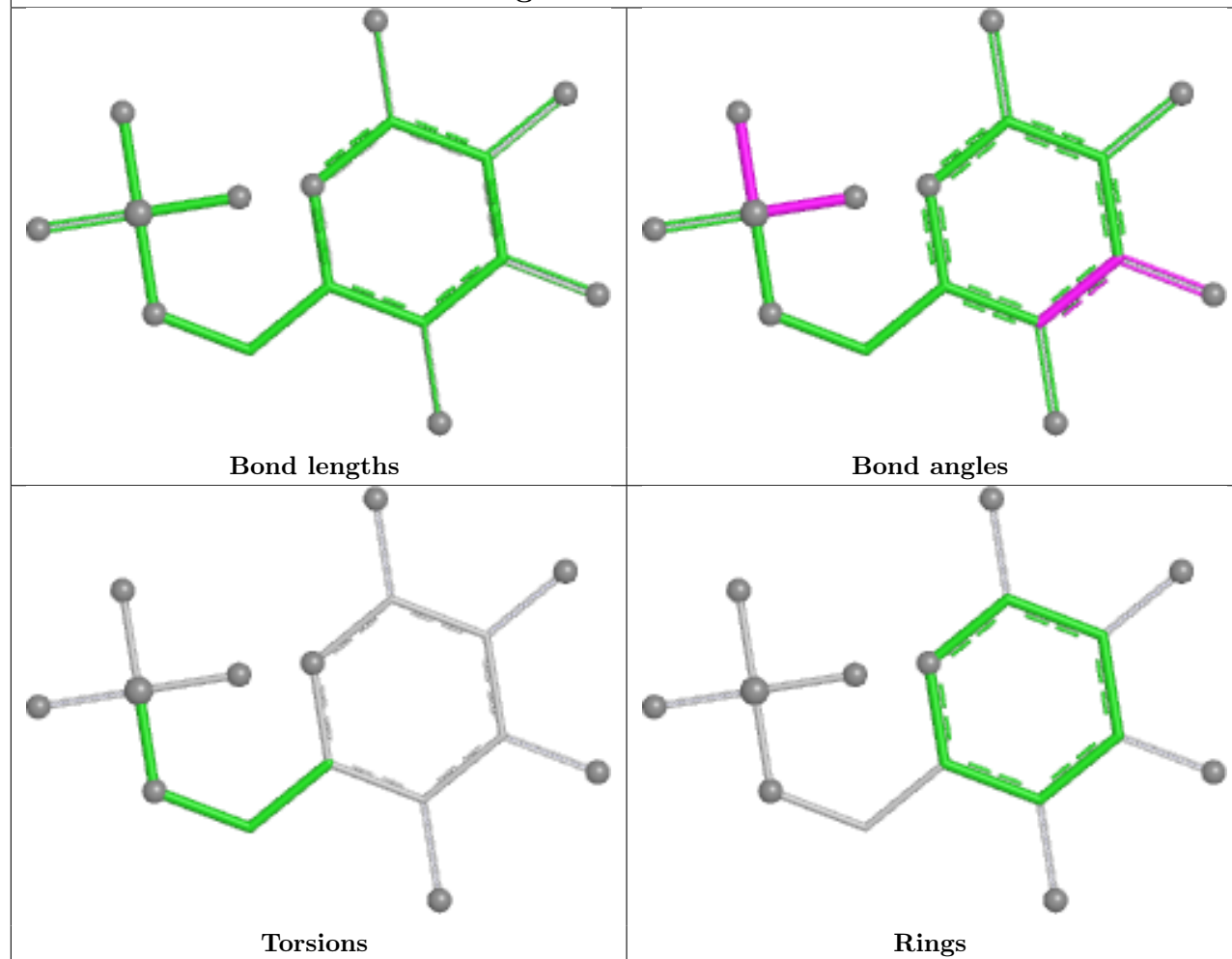
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	803	G6P	1	0
4	B	802	G6P	1	0
2	C	801	GCS	1	0
3	C	802	UDP	1	0
3	D	801	UDP	2	0
4	A	802	G6P	1	0
4	D	802	G6P	2	0
3	A	801	UDP	1	0

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.

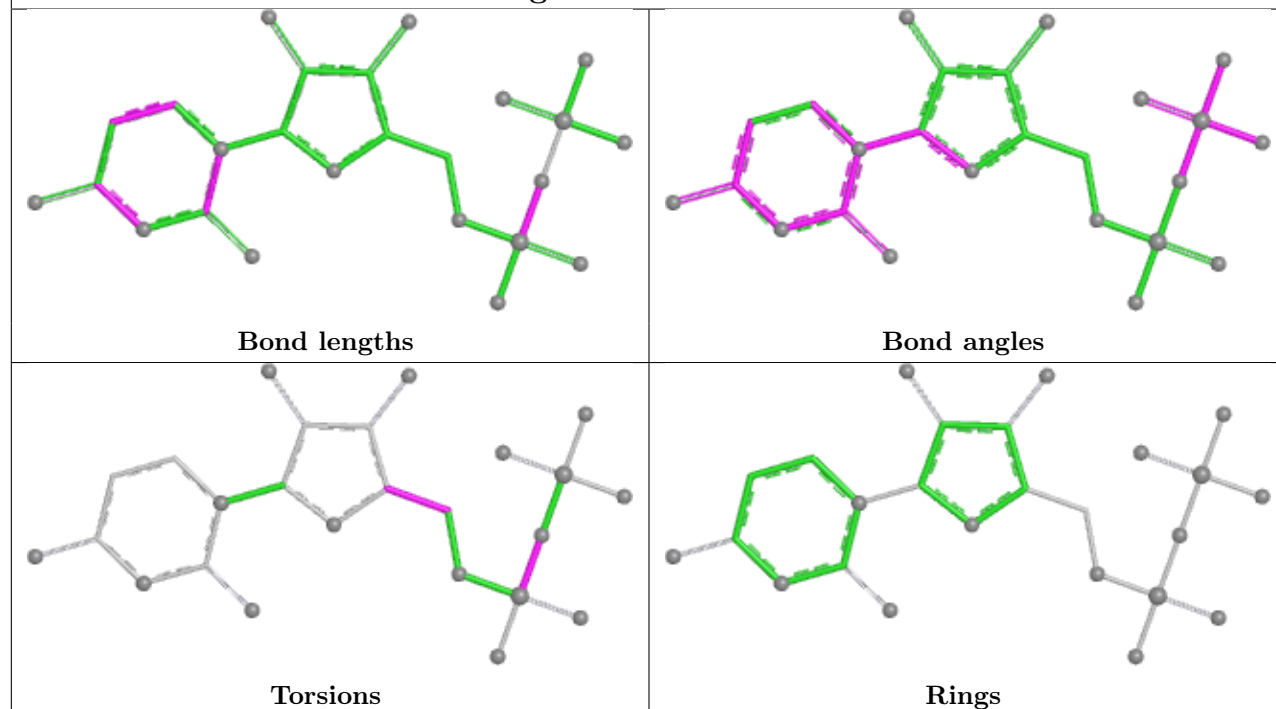
Ligand G6P C 803

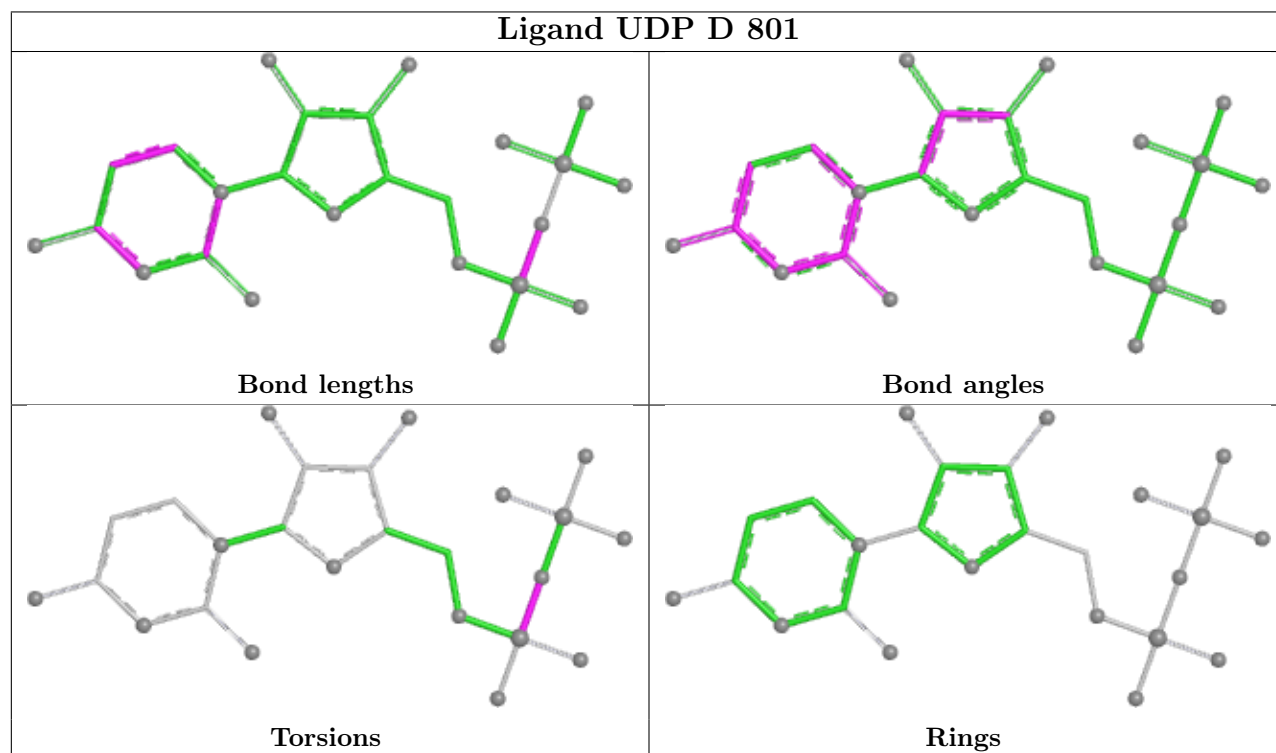
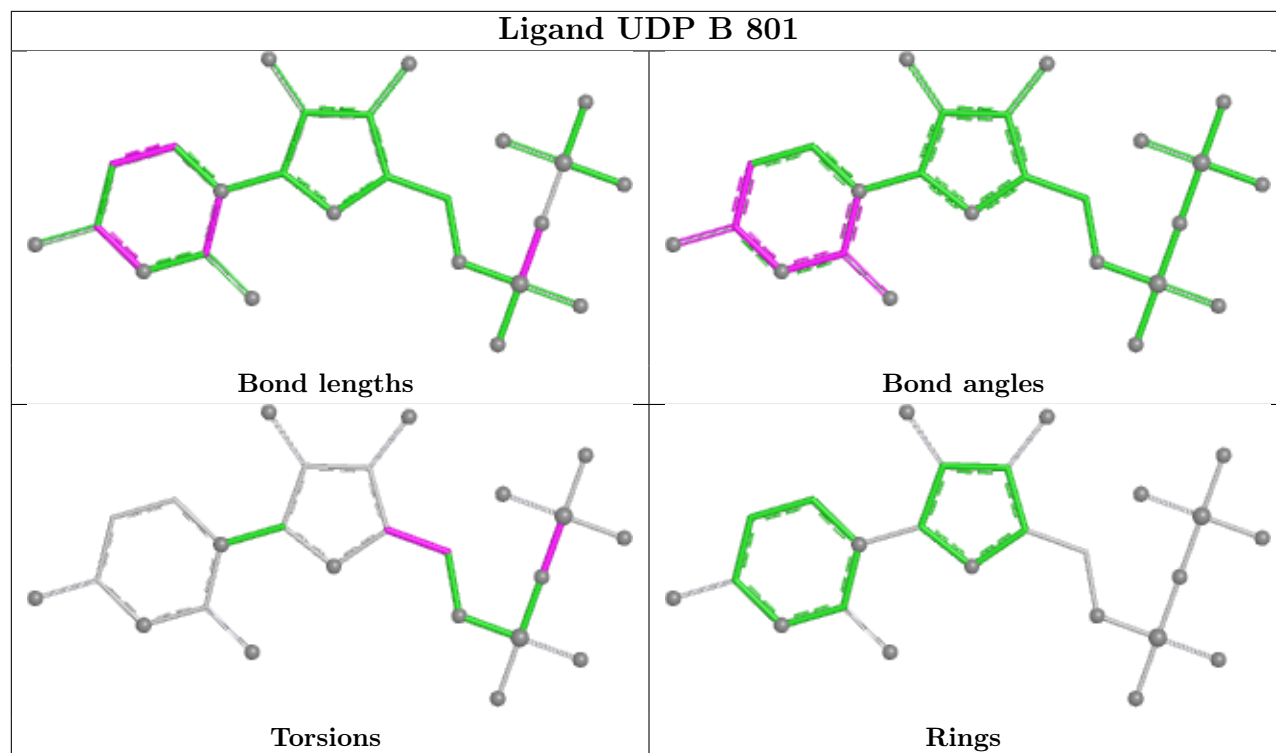


Ligand G6P B 802

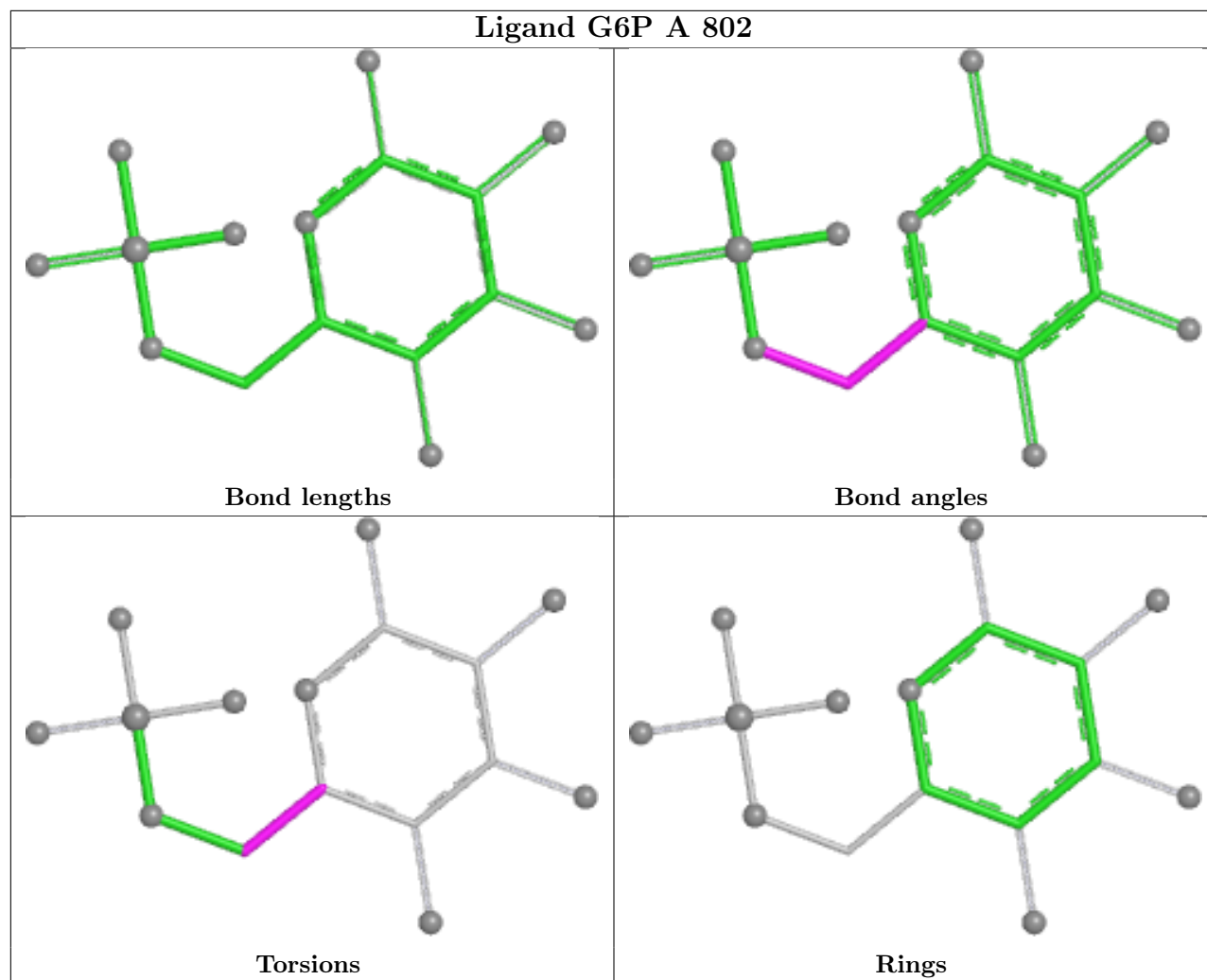


Ligand UDP C 802

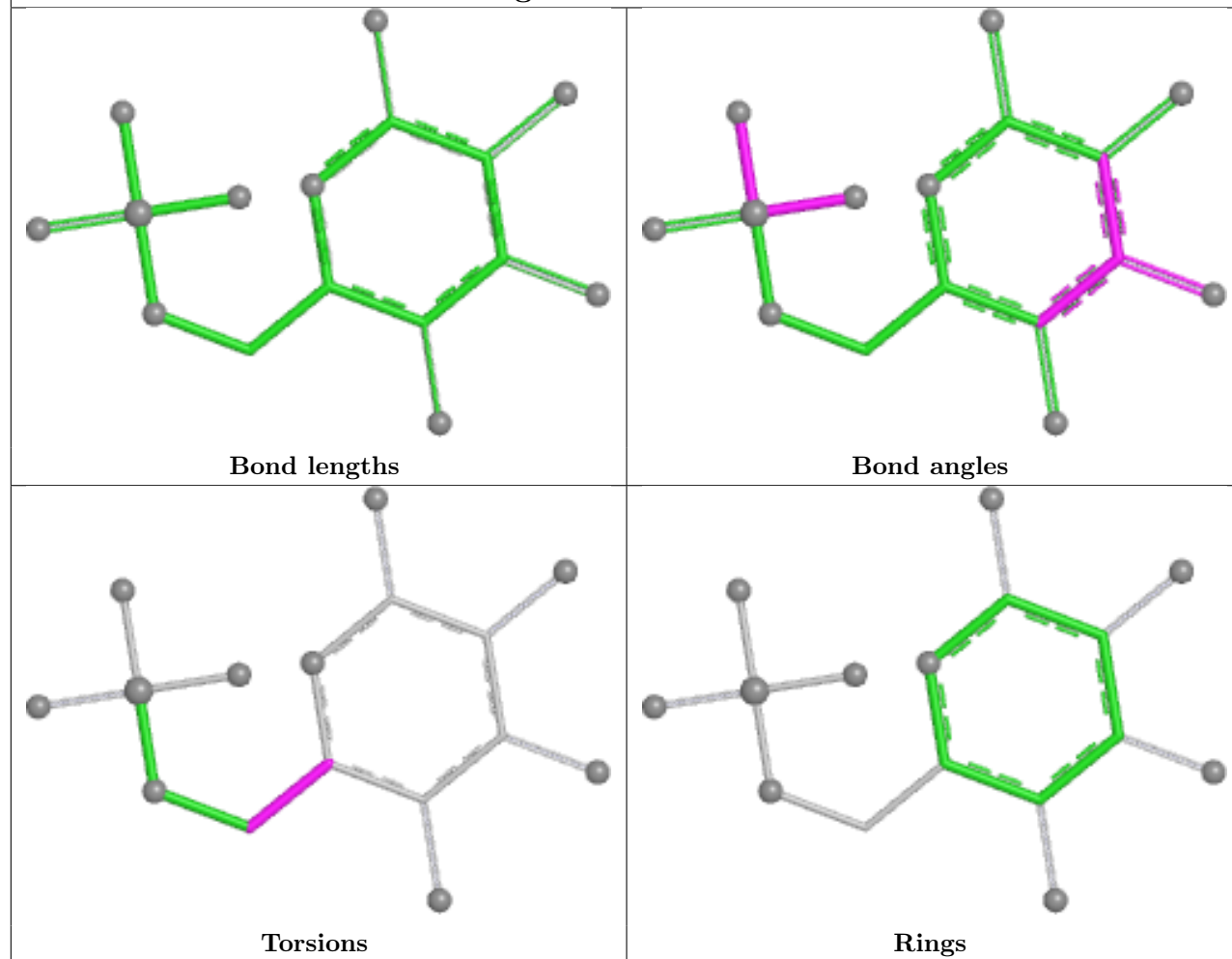




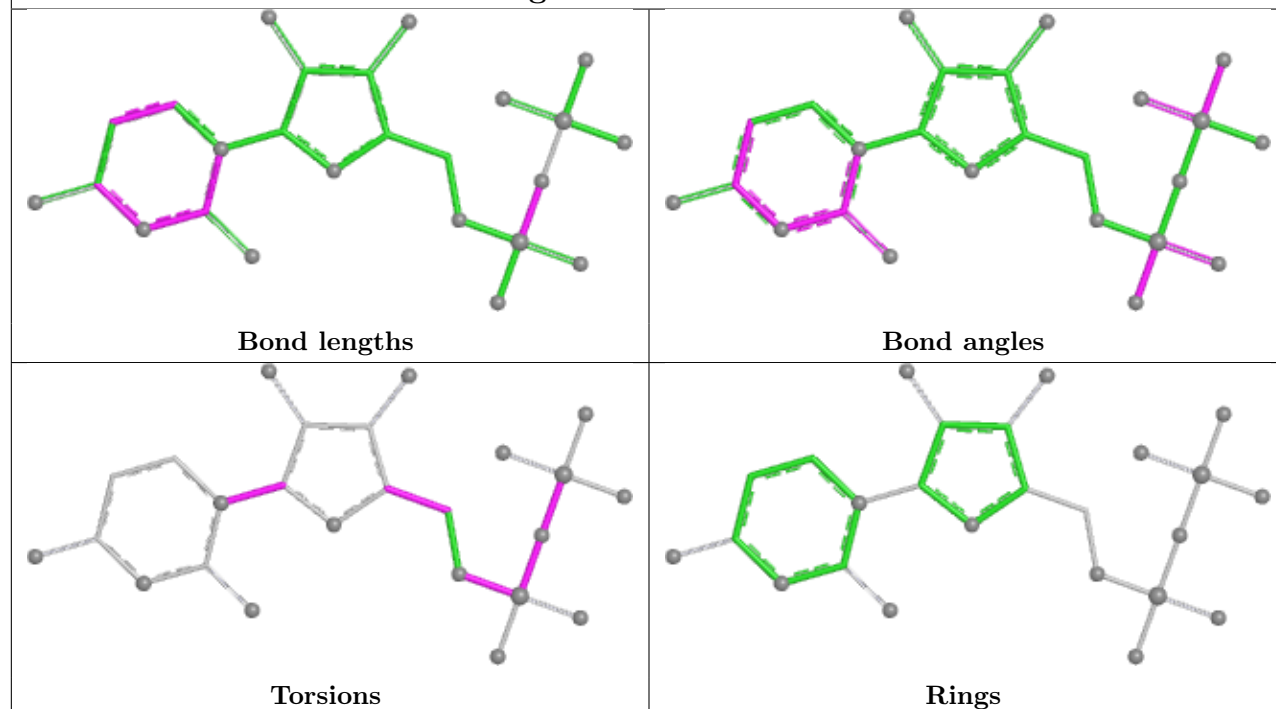
Ligand G6P A 802



Ligand G6P D 802



Ligand UDP A 801



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 [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, 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%)	-0.17	5 (0%) 82 67	50, 86, 141, 178	0
1	B	636/720 (88%)	-0.41	2 (0%) 90 82	39, 71, 115, 166	0
1	C	638/720 (88%)	-0.25	3 (0%) 87 75	45, 80, 129, 169	0
1	D	636/720 (88%)	-0.20	7 (1%) 78 61	42, 87, 146, 164	0
All	All	2548/2880 (88%)	-0.26	17 (0%) 84 70	39, 80, 138, 178	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	61	LEU	3.8
1	C	125	LEU	3.2
1	A	131	PRO	2.9
1	D	628	GLU	2.7
1	A	130	SER	2.6
1	D	61	LEU	2.6
1	A	129	PRO	2.6
1	D	639	ALA	2.5
1	C	97	ILE	2.4
1	C	205	GLY	2.4
1	D	630	LEU	2.4
1	D	197	LEU	2.3
1	B	169	GLN	2.3
1	D	227	ILE	2.2
1	B	205	GLY	2.2
1	A	20	ARG	2.1
1	D	629	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

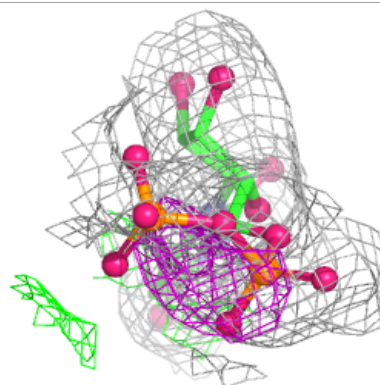
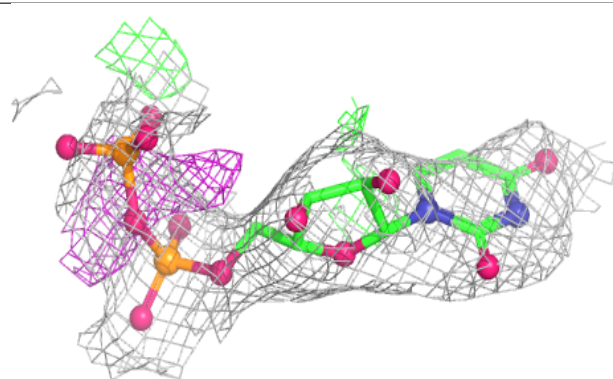
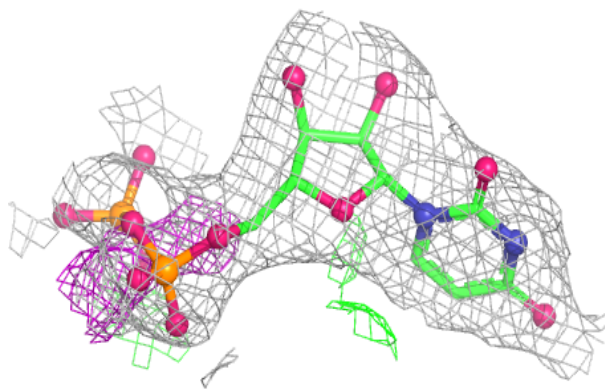
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	0.83	0.12	81,116,172,177	0
3	UDP	D	801	25/25	0.90	0.08	80,99,127,130	0
3	UDP	A	801	25/25	0.92	0.07	79,94,109,114	0
2	GCS	C	801	12/12	0.96	0.07	60,67,78,78	0
3	UDP	C	802	25/25	0.97	0.05	61,77,83,86	0
4	G6P	C	803	16/16	0.97	0.06	56,66,75,76	0
4	G6P	D	802	16/16	0.97	0.06	43,51,58,63	0
4	G6P	A	802	16/16	0.98	0.05	50,58,65,66	0
4	G6P	B	802	16/16	0.98	0.05	39,49,55,58	0

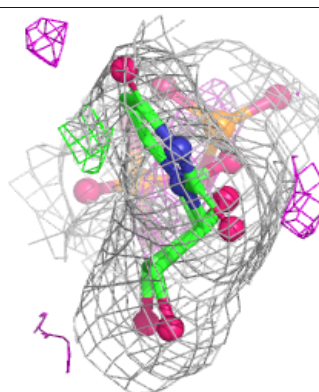
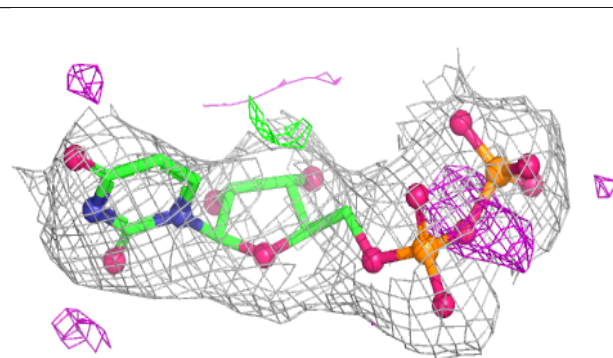
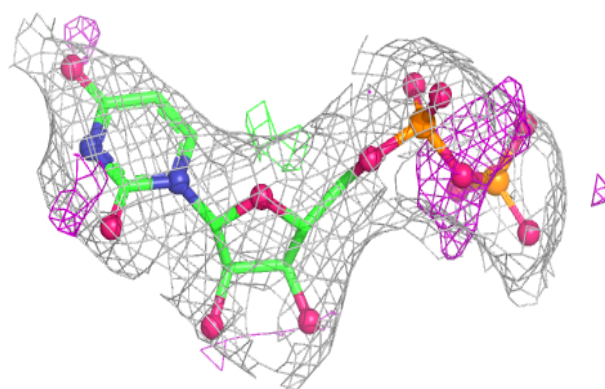
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around UDP B 801:

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

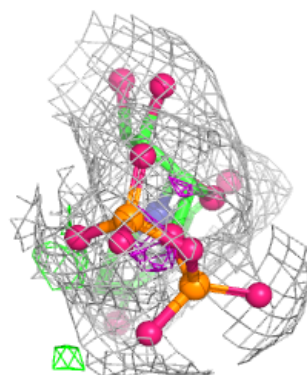
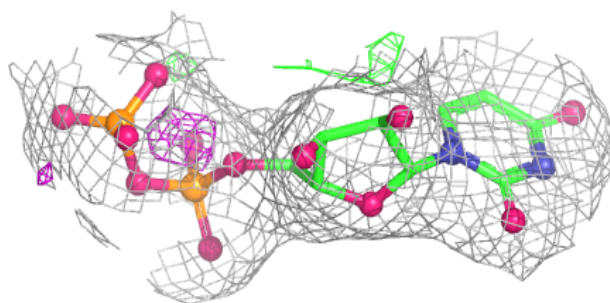
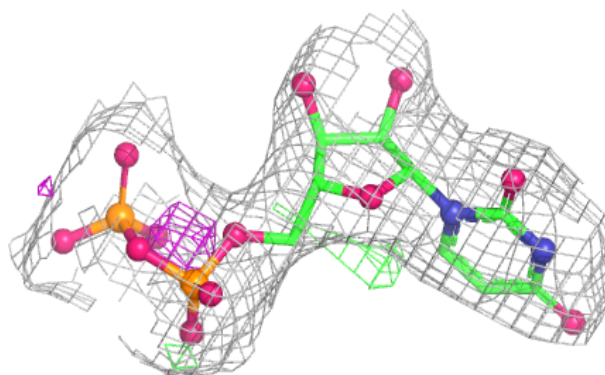
**Electron density around UDP D 801:**

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

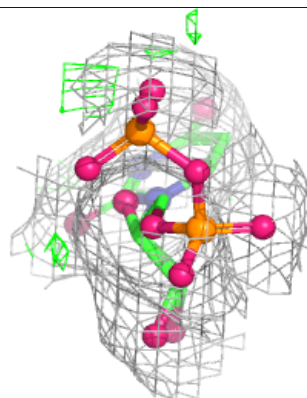
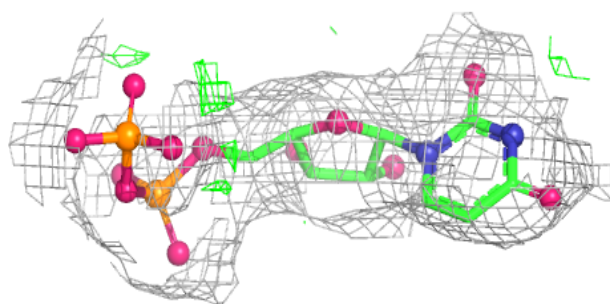
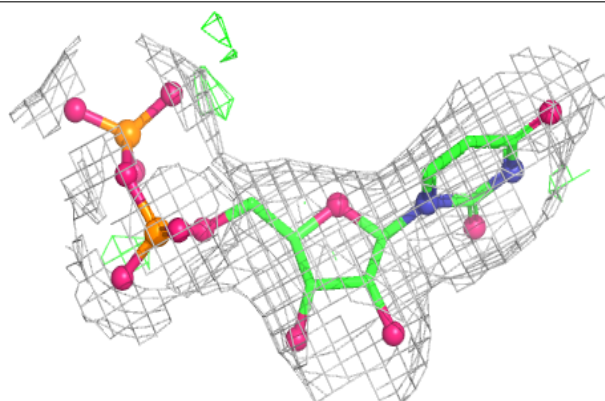


Electron density around UDP A 801:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

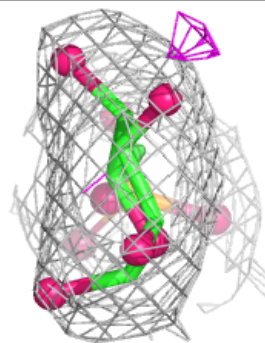
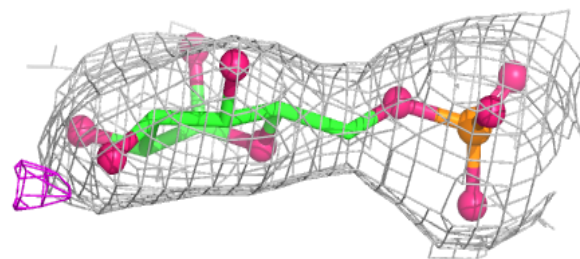
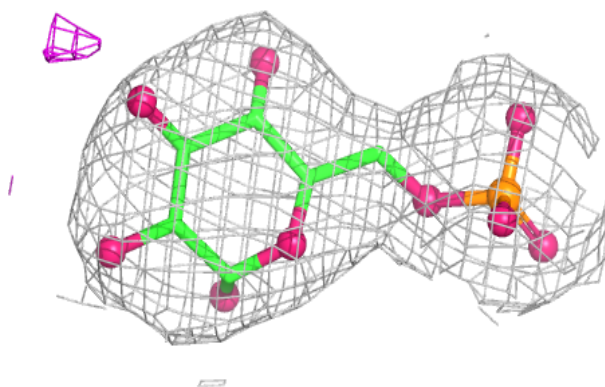
**Electron density around UDP C 802:**

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

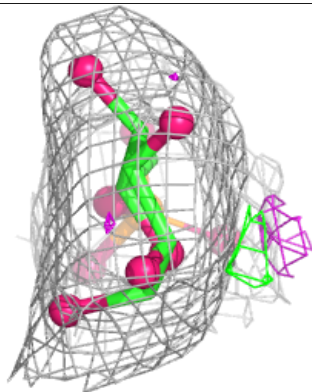
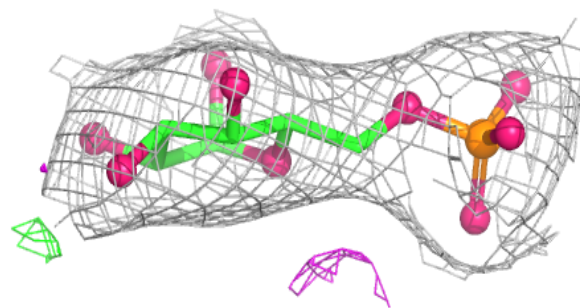
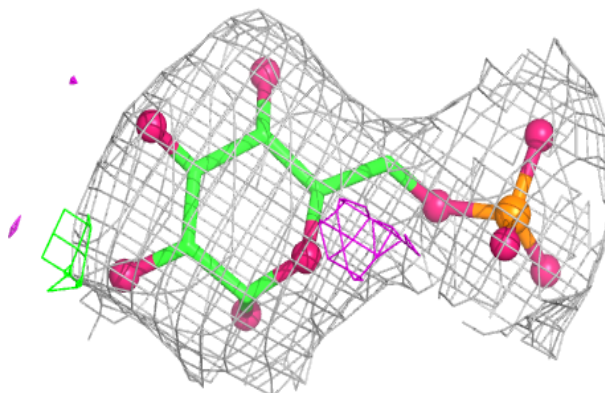


Electron density around G6P C 803:

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

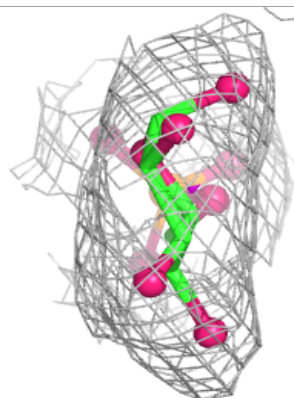
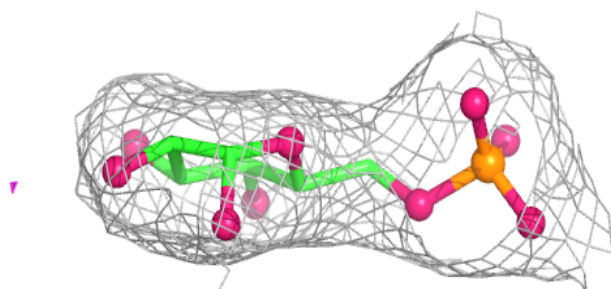
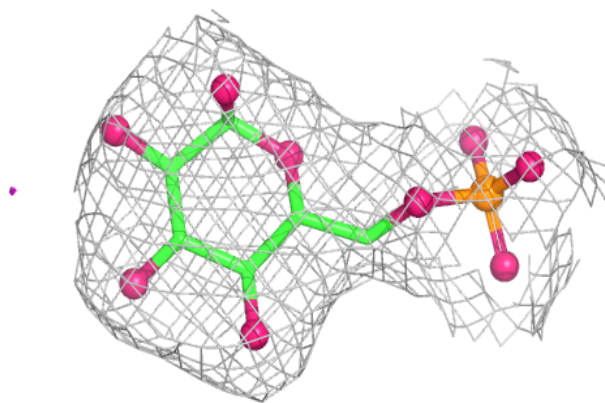
**Electron density around G6P D 802:**

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

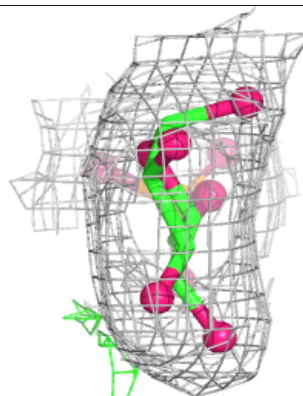
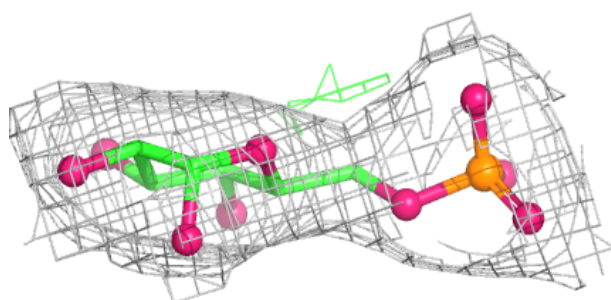


Electron density around G6P A 802:

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

**Electron density around G6P B 802:**

$2mF_o-DF_c$ (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.