



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:36 PM UTC

PDB ID : 9LQL / pdb_00009lql
Title : Structure of STG-hydrolyzing beta-glucosidase 1 (PSTG1) complexed with heptyl 1-thio-beta-D-glucopyranoside
Authors : Yanai, T.; Arai, A.; Takahashi, Y.; Imaizumi, R.; Takeshita, K.; Matsuura, H.; Sakai, N.; Takahashi, S.; Yamamoto, M.; Kataoka, K.; Nakayama, T.; Yamashita, S.
Deposited on : 2025-01-28
Resolution : 3.35 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](#) ①) 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

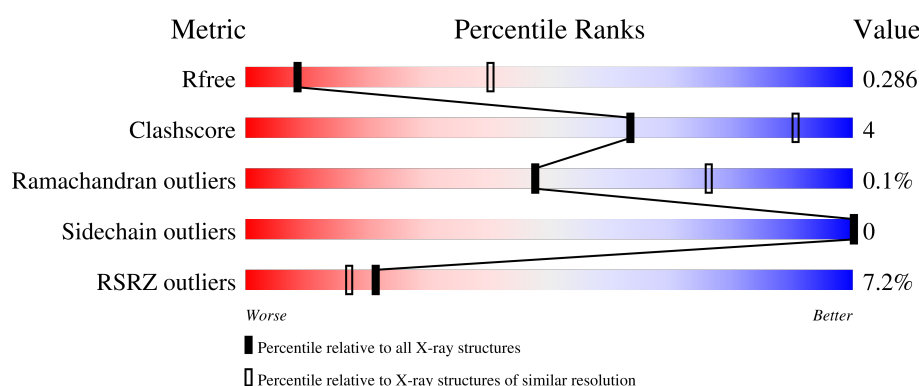
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.35 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	1099 (3.40-3.32)
Clashscore	190562	1116 (3.40-3.32)
Ramachandran outliers	187476	1101 (3.40-3.32)
Sidechain outliers	187428	1101 (3.40-3.32)
RSRZ outliers	180081	1099 (3.40-3.32)

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	780	 6% 83% 11% 6%
1	B	780	 7% 84% 11% 5%
1	C	780	 6% 87% 8% 5%
1	D	780	 7% 84% 11% 5%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 23162 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 Beta-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	732	Total	C	N	O	S	0	0	0
			5698	3603	983	1087	25			
1	B	740	Total	C	N	O	S	0	0	0
			5759	3643	991	1100	25			
1	C	738	Total	C	N	O	S	0	0	0
			5745	3633	992	1095	25			
1	D	741	Total	C	N	O	S	0	0	0
			5765	3647	994	1099	25			

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-26	MET	-	initiating methionine	UNP D6RVX0
A	-25	ASN	-	expression tag	UNP D6RVX0
A	-24	HIS	-	expression tag	UNP D6RVX0
A	-23	LYS	-	expression tag	UNP D6RVX0
A	-22	VAL	-	expression tag	UNP D6RVX0
A	-21	HIS	-	expression tag	UNP D6RVX0
A	-20	HIS	-	expression tag	UNP D6RVX0
A	-19	HIS	-	expression tag	UNP D6RVX0
A	-18	HIS	-	expression tag	UNP D6RVX0
A	-17	HIS	-	expression tag	UNP D6RVX0
A	-16	HIS	-	expression tag	UNP D6RVX0
A	-15	ILE	-	expression tag	UNP D6RVX0
A	-14	GLU	-	expression tag	UNP D6RVX0
A	-13	GLY	-	expression tag	UNP D6RVX0
A	-12	ARG	-	expression tag	UNP D6RVX0
A	-11	HIS	-	expression tag	UNP D6RVX0
A	-10	MET	-	expression tag	UNP D6RVX0
A	-9	GLU	-	expression tag	UNP D6RVX0
A	-8	LEU	-	expression tag	UNP D6RVX0
A	-7	GLY	-	expression tag	UNP D6RVX0
A	-6	THR	-	expression tag	UNP D6RVX0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	LEU	-	expression tag	UNP D6RVX0
A	-4	GLU	-	expression tag	UNP D6RVX0
A	-3	GLY	-	expression tag	UNP D6RVX0
A	-2	SER	-	expression tag	UNP D6RVX0
A	-1	GLU	-	expression tag	UNP D6RVX0
A	0	PHE	-	expression tag	UNP D6RVX0
B	-26	MET	-	initiating methionine	UNP D6RVX0
B	-25	ASN	-	expression tag	UNP D6RVX0
B	-24	HIS	-	expression tag	UNP D6RVX0
B	-23	LYS	-	expression tag	UNP D6RVX0
B	-22	VAL	-	expression tag	UNP D6RVX0
B	-21	HIS	-	expression tag	UNP D6RVX0
B	-20	HIS	-	expression tag	UNP D6RVX0
B	-19	HIS	-	expression tag	UNP D6RVX0
B	-18	HIS	-	expression tag	UNP D6RVX0
B	-17	HIS	-	expression tag	UNP D6RVX0
B	-16	HIS	-	expression tag	UNP D6RVX0
B	-15	ILE	-	expression tag	UNP D6RVX0
B	-14	GLU	-	expression tag	UNP D6RVX0
B	-13	GLY	-	expression tag	UNP D6RVX0
B	-12	ARG	-	expression tag	UNP D6RVX0
B	-11	HIS	-	expression tag	UNP D6RVX0
B	-10	MET	-	expression tag	UNP D6RVX0
B	-9	GLU	-	expression tag	UNP D6RVX0
B	-8	LEU	-	expression tag	UNP D6RVX0
B	-7	GLY	-	expression tag	UNP D6RVX0
B	-6	THR	-	expression tag	UNP D6RVX0
B	-5	LEU	-	expression tag	UNP D6RVX0
B	-4	GLU	-	expression tag	UNP D6RVX0
B	-3	GLY	-	expression tag	UNP D6RVX0
B	-2	SER	-	expression tag	UNP D6RVX0
B	-1	GLU	-	expression tag	UNP D6RVX0
B	0	PHE	-	expression tag	UNP D6RVX0
C	-10	MET	-	initiating methionine	UNP D6RVX0
C	-9	ASN	-	expression tag	UNP D6RVX0
C	-8	HIS	-	expression tag	UNP D6RVX0
C	-7	LYS	-	expression tag	UNP D6RVX0
C	-6	VAL	-	expression tag	UNP D6RVX0
C	-5	HIS	-	expression tag	UNP D6RVX0
C	-4	HIS	-	expression tag	UNP D6RVX0
C	-3	HIS	-	expression tag	UNP D6RVX0
C	-2	HIS	-	expression tag	UNP D6RVX0

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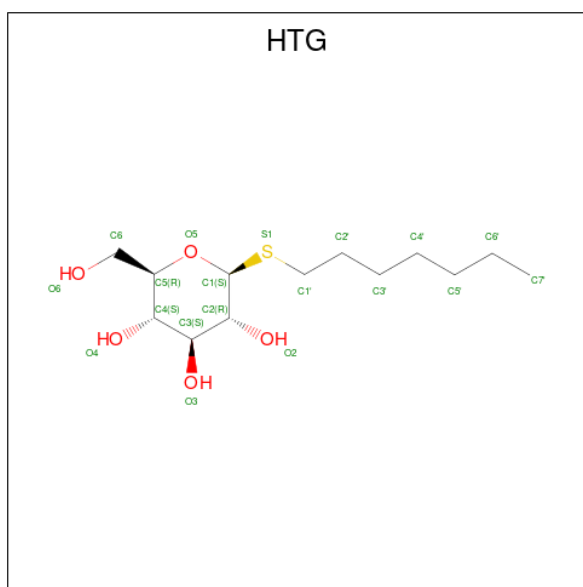
Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	HIS	-	expression tag	UNP D6RVX0
C	0	HIS	-	expression tag	UNP D6RVX0
C	1	ILE	-	expression tag	UNP D6RVX0
C	2	GLU	-	expression tag	UNP D6RVX0
C	3	GLY	-	expression tag	UNP D6RVX0
C	4	ARG	-	expression tag	UNP D6RVX0
C	5	HIS	-	expression tag	UNP D6RVX0
C	6	MET	-	expression tag	UNP D6RVX0
C	7	GLU	-	expression tag	UNP D6RVX0
C	8	LEU	-	expression tag	UNP D6RVX0
C	9	GLY	-	expression tag	UNP D6RVX0
C	10	THR	-	expression tag	UNP D6RVX0
C	11	LEU	-	expression tag	UNP D6RVX0
C	12	GLU	-	expression tag	UNP D6RVX0
C	13	GLY	-	expression tag	UNP D6RVX0
C	14	SER	-	expression tag	UNP D6RVX0
C	15	GLU	-	expression tag	UNP D6RVX0
C	16	PHE	-	expression tag	UNP D6RVX0
D	-26	MET	-	initiating methionine	UNP D6RVX0
D	-25	ASN	-	expression tag	UNP D6RVX0
D	-24	HIS	-	expression tag	UNP D6RVX0
D	-23	LYS	-	expression tag	UNP D6RVX0
D	-22	VAL	-	expression tag	UNP D6RVX0
D	-21	HIS	-	expression tag	UNP D6RVX0
D	-20	HIS	-	expression tag	UNP D6RVX0
D	-19	HIS	-	expression tag	UNP D6RVX0
D	-18	HIS	-	expression tag	UNP D6RVX0
D	-17	HIS	-	expression tag	UNP D6RVX0
D	-16	HIS	-	expression tag	UNP D6RVX0
D	-15	ILE	-	expression tag	UNP D6RVX0
D	-14	GLU	-	expression tag	UNP D6RVX0
D	-13	GLY	-	expression tag	UNP D6RVX0
D	-12	ARG	-	expression tag	UNP D6RVX0
D	-11	HIS	-	expression tag	UNP D6RVX0
D	-10	MET	-	expression tag	UNP D6RVX0
D	-9	GLU	-	expression tag	UNP D6RVX0
D	-8	LEU	-	expression tag	UNP D6RVX0
D	-7	GLY	-	expression tag	UNP D6RVX0
D	-6	THR	-	expression tag	UNP D6RVX0
D	-5	LEU	-	expression tag	UNP D6RVX0
D	-4	GLU	-	expression tag	UNP D6RVX0
D	-3	GLY	-	expression tag	UNP D6RVX0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	SER	-	expression tag	UNP D6RVX0
D	-1	GLU	-	expression tag	UNP D6RVX0
D	0	PHE	-	expression tag	UNP D6RVX0

- Molecule 2 is heptyl 1-thio-beta-D-glucopyranoside (CCD ID: HTG) (formula: $C_{13}H_{26}O_5S$) (labeled as "Ligand of Interest" by depositor).



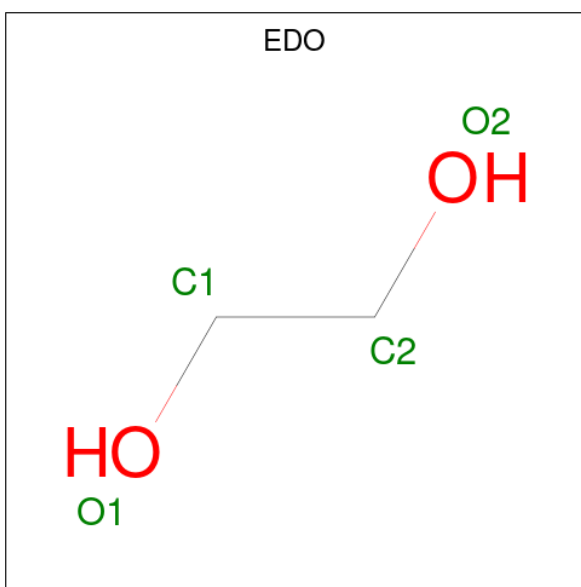
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			19	13	5	1		
2	B	1	Total	C	O	S	0	0
			19	13	5	1		
2	C	1	Total	C	O	S	0	0
			19	13	5	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



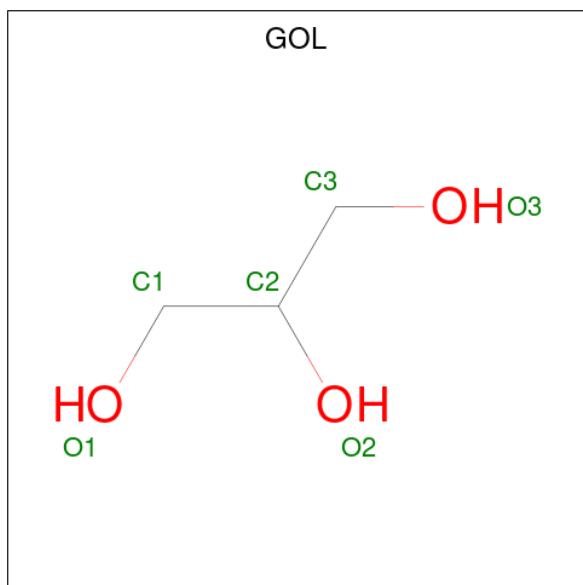
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			4	2	2		

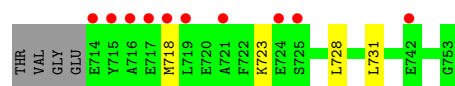
- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



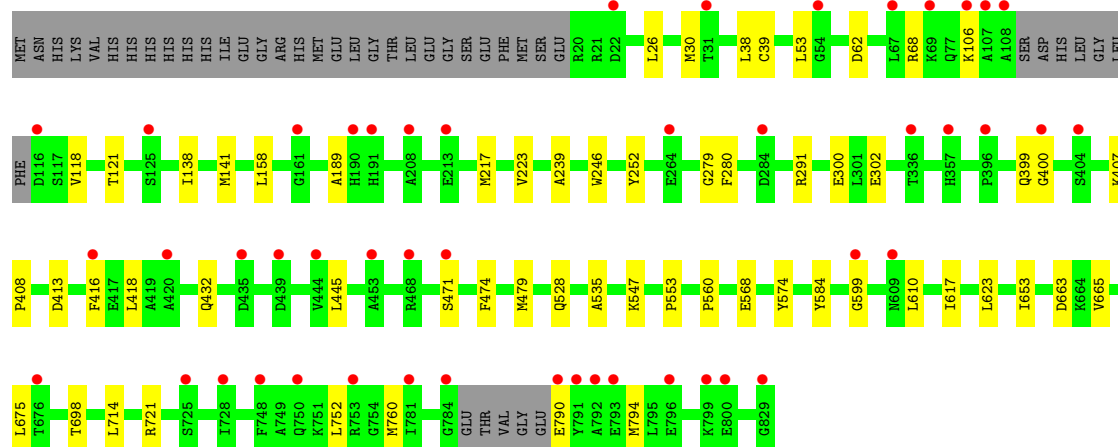
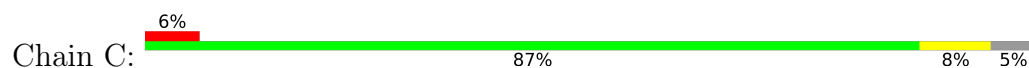
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

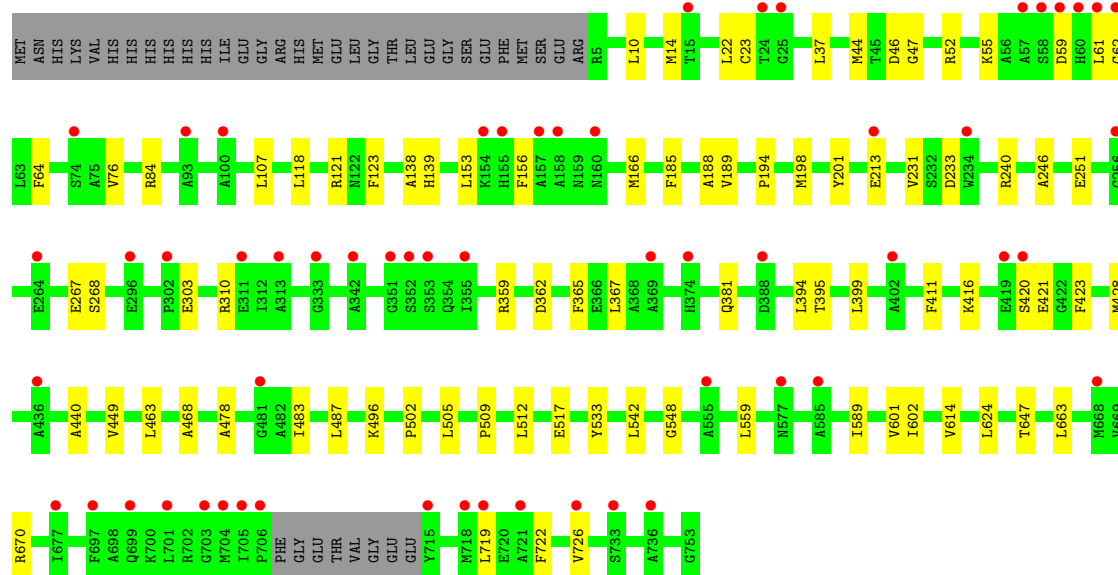
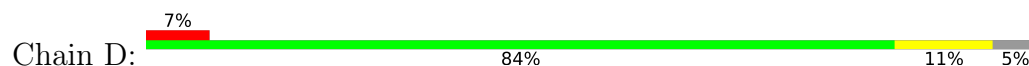
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	22	Total	O	0	0
			22	22		
6	B	34	Total	O	0	0
			34	34		
6	C	23	Total	O	0	0
			23	23		
6	D	30	Total	O	0	0
			30	30		



● Molecule 1: Beta-glucosidase



● Molecule 1: Beta-glucosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	64.60Å 112.04Å 156.59Å 71.50° 82.83° 79.85°	Depositor
Resolution (Å)	45.85 – 3.35 45.85 – 3.35	Depositor EDS
% Data completeness (in resolution range)	98.2 (45.85-3.35) 98.2 (45.85-3.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 3.32Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.245 , 0.286 0.245 , 0.286	Depositor DCC
R_{free} test set	2872 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	1.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	23162	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.34% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, HTG, SO4, EDO

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	0.08	0/5804	0.25	0/7841
1	B	0.08	0/5868	0.25	0/7928
1	C	0.07	0/5853	0.24	0/7907
1	D	0.09	0/5875	0.25	0/7940
All	All	0.08	0/23400	0.25	0/31616

There are no bond length outliers.

There are no bond angle outliers.

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	5698	0	5691	52	0
1	B	5759	0	5739	47	0
1	C	5745	0	5734	36	0
1	D	5765	0	5754	54	0
2	A	19	0	26	2	0
2	B	19	0	26	1	0
2	C	19	0	26	3	0
3	A	5	0	0	0	0
3	C	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	5	0	0	0	0
4	B	4	0	6	0	0
4	C	4	0	6	1	0
5	D	6	0	8	0	0
6	A	22	0	0	1	0
6	B	34	0	0	0	0
6	C	23	0	0	0	0
6	D	30	0	0	0	0
All	All	23162	0	23016	182	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 4.

All (182) 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:D:722:PHE:O	1:D:726:VAL:HG23	1.83	0.77
1:A:46:ASP:HA	1:A:107:LEU:HB2	1.75	0.68
1:B:14:MET:HE1	1:B:284:LEU:HD11	1.76	0.66
1:B:701:LEU:HD12	1:B:731:LEU:HD13	1.77	0.66
1:B:141:ILE:HD11	1:B:153:LEU:HD13	1.78	0.66
1:A:141:ILE:HD11	1:A:153:LEU:HD13	1.78	0.65
1:D:46:ASP:HA	1:D:107:LEU:HB2	1.79	0.65
1:C:752:LEU:HA	1:C:760:MET:HE2	1.80	0.63
1:D:10:LEU:HD11	1:D:37:LEU:HB3	1.80	0.63
1:C:217:MET:HA	1:C:252:TYR:HB3	1.81	0.63
1:D:23:CYS:HB3	1:D:240:ARG:HH21	1.64	0.62
1:C:617:ILE:HD13	1:C:623:LEU:HB2	1.82	0.62
1:A:166:MET:HA	1:A:201:TYR:HB3	1.83	0.61
1:C:106:LYS:HE3	1:C:118:VAL:HG22	1.81	0.61
1:B:166:MET:HA	1:B:201:TYR:HB3	1.84	0.60
1:A:14:MET:HE1	1:A:22:LEU:HD12	1.84	0.60
1:B:46:ASP:HA	1:B:107:LEU:HB2	1.84	0.59
1:B:612:ASP:HB3	1:B:624:LEU:HD23	1.83	0.59
1:B:55:LYS:HE3	1:B:67:VAL:HG22	1.87	0.57
1:C:62:ASP:HA	1:C:158:LEU:HB2	1.87	0.56
1:D:647:THR:HG23	1:D:670:ARG:HG2	1.88	0.56
1:B:647:THR:HG23	1:B:670:ARG:HG2	1.88	0.55
1:D:496:LYS:HG2	1:D:548:GLY:HA3	1.87	0.55
1:A:362:ASP:HB3	1:A:365:PHE:HB3	1.88	0.55
1:A:352:SER:HB3	2:A:801:HTG:H1	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:610:LEU:HB2	1:C:714:LEU:HD13	1.89	0.54
1:C:26:LEU:HD11	1:C:53:LEU:HB3	1.89	0.54
1:A:23:CYS:HB3	1:A:240:ARG:HH21	1.72	0.54
1:D:153:LEU:HD22	1:D:185:PHE:HD2	1.73	0.54
1:A:559:LEU:HB2	1:A:663:LEU:HD13	1.91	0.53
1:A:18:GLU:HB3	1:A:34:ILE:HD13	1.90	0.53
1:D:166:MET:HA	1:D:201:TYR:HB3	1.91	0.53
1:B:10:LEU:HD11	1:B:37:LEU:HB3	1.91	0.53
1:B:559:LEU:HB2	1:B:663:LEU:HD13	1.91	0.53
1:D:14:MET:HE1	1:D:22:LEU:HD12	1.90	0.53
1:D:399:LEU:HD21	1:D:440:ALA:HB1	1.90	0.52
1:C:413:ASP:HB3	1:C:416:PHE:HB3	1.90	0.52
1:A:24:THR:HG21	1:A:255:GLU:H	1.74	0.51
1:D:449:VAL:HG22	1:D:468:ALA:HB3	1.92	0.51
1:A:44:MET:HE2	1:A:234:TRP:HH2	1.74	0.51
1:A:496:LYS:HG2	1:A:548:GLY:HA3	1.92	0.51
1:A:647:THR:HG23	1:A:670:ARG:HG2	1.92	0.51
1:B:449:VAL:HG22	1:B:468:ALA:HB3	1.93	0.51
1:D:559:LEU:HB2	1:D:663:LEU:HD13	1.92	0.51
1:B:24:THR:HG21	1:B:255:GLU:H	1.75	0.50
1:D:44:MET:HE3	1:D:107:LEU:HD21	1.92	0.50
4:C:903:EDO:H12	1:D:512:LEU:HD12	1.93	0.50
1:C:68:ARG:HH21	2:C:901:HTG:H5'2	1.77	0.50
1:D:428:MET:HE1	1:D:502:PRO:HD2	1.94	0.49
1:D:362:ASP:HB3	1:D:365:PHE:HB3	1.94	0.49
1:B:362:ASP:HB3	1:B:365:PHE:HB3	1.94	0.49
1:D:153:LEU:CD2	1:D:185:PHE:HD2	2.25	0.49
1:C:432:GLN:HB3	1:C:445:LEU:HD22	1.94	0.49
1:A:399:LEU:HD21	1:A:440:ALA:HB1	1.94	0.49
1:B:237:VAL:HG11	1:B:251:GLU:HG3	1.94	0.49
2:A:801:HTG:H6'2	1:B:718:MET:HE2	1.93	0.49
1:C:30:MET:HE1	1:C:38:LEU:HD12	1.94	0.49
1:B:702:ARG:HH21	1:B:723:LYS:HB3	1.78	0.49
1:C:479:MET:HE1	1:C:553:PRO:HD2	1.95	0.49
1:D:153:LEU:CD2	1:D:185:PHE:CD2	2.96	0.48
1:B:57:ALA:HB1	1:B:64:PHE:HD2	1.79	0.48
1:C:698:THR:HG23	1:C:721:ARG:HG2	1.96	0.48
1:D:411:PHE:HE2	1:D:483:ILE:HD13	1.79	0.48
1:A:428:MET:HE1	1:A:502:PRO:HD2	1.95	0.48
1:B:728:LEU:HA	1:B:731:LEU:HD12	1.96	0.48
1:D:84:ARG:HD3	1:D:139:HIS:NE2	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:ARG:HD3	1:B:139:HIS:NE2	2.29	0.47
1:C:547:LYS:HG2	1:C:599:GLY:HA3	1.95	0.47
1:D:198:MET:HA	1:D:231:VAL:O	2.15	0.47
1:B:23:CYS:HB3	1:B:240:ARG:HH21	1.79	0.47
1:C:653:ILE:HG13	1:D:517:GLU:HG3	1.97	0.47
1:D:153:LEU:HD22	1:D:185:PHE:CD2	2.50	0.47
1:D:213:GLU:HG2	1:D:246:ALA:HA	1.96	0.46
1:A:420:SER:HB3	1:A:423:PHE:CZ	2.50	0.46
1:A:322:ASN:HD21	1:A:327:LEU:HB2	1.81	0.46
1:A:719:LEU:HD22	1:A:723:LYS:NZ	2.30	0.46
1:B:399:LEU:HD21	1:B:440:ALA:HB1	1.97	0.46
1:C:399:GLN:HG2	1:C:400:GLY:H	1.81	0.46
1:A:355:ILE:HD11	6:A:905:HOH:O	2.15	0.46
1:B:702:ARG:NH2	1:B:723:LYS:HB3	2.31	0.46
1:B:76:VAL:HB	1:B:123:PHE:HA	1.99	0.45
1:D:381:GLN:HB3	1:D:394:LEU:HD22	1.98	0.45
1:A:138:ALA:HB2	1:A:188:ALA:HA	1.97	0.45
1:B:155:HIS:HA	1:B:200:SER:HB3	1.99	0.45
1:D:118:LEU:HD11	1:D:505:LEU:HD21	1.98	0.45
1:A:84:ARG:HB3	1:A:139:HIS:CG	2.52	0.45
1:B:197:VAL:O	1:B:230:VAL:HA	2.17	0.45
1:A:322:ASN:ND2	1:A:327:LEU:HB2	2.31	0.45
1:A:587:LYS:HG2	1:A:615:PHE:HD1	1.82	0.45
1:D:614:VAL:HB	1:D:624:LEU:HD21	1.98	0.45
1:C:39:CYS:HB3	1:C:291:ARG:HH21	1.82	0.45
1:A:449:VAL:HG22	1:A:468:ALA:HB3	1.99	0.44
1:B:677:ILE:HG13	1:B:690:LEU:HD13	1.99	0.44
1:C:68:ARG:NH2	2:C:901:HTG:H5'2	2.33	0.44
1:B:26:ARG:HD3	1:B:33:PRO:HD3	2.00	0.44
1:A:692:GLU:H	1:A:692:GLU:HG2	1.63	0.44
1:B:107:LEU:HA	1:B:152:SER:HB2	2.00	0.44
1:D:240:ARG:HD3	1:D:251:GLU:OE1	2.17	0.44
1:B:496:LYS:HG2	1:B:548:GLY:HA3	1.99	0.44
1:A:240:ARG:HD3	1:A:251:GLU:OE1	2.18	0.44
1:C:794:MET:SD	1:D:61:LEU:HB3	2.58	0.44
1:D:310:ARG:HG3	1:D:478:ALA:HB1	1.99	0.44
1:A:556:TYR:HE2	1:A:590:VAL:HG13	1.83	0.44
1:C:121:THR:HG23	1:C:528:GLN:NE2	2.33	0.44
1:C:418:LEU:HD12	1:C:535:ALA:HB2	1.98	0.44
1:A:84:ARG:HB3	1:A:139:HIS:CD2	2.53	0.43
1:C:223:VAL:HG22	1:C:574:TYR:CG	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:280:PHE:HB2	1:C:300:GLU:HB2	2.00	0.43
1:A:748:LEU:HD13	1:D:268:SER:HB2	1.99	0.43
1:A:509:PRO:HB3	1:A:533:TYR:CG	2.53	0.43
1:B:156:PHE:O	1:B:156:PHE:CG	2.71	0.43
1:C:407:LYS:HD2	1:C:408:PRO:HD2	2.00	0.43
1:A:602:ILE:HG13	1:B:517:GLU:HG3	2.00	0.43
1:D:55:LYS:H	1:D:55:LYS:HG2	1.57	0.43
1:C:138:ILE:HG13	1:C:141:MET:HE2	2.00	0.43
1:D:121:ARG:HD2	1:D:421:GLU:HG3	2.01	0.43
1:A:68:PRO:HA	1:A:356:LYS:HB3	2.01	0.43
1:A:198:MET:HA	1:A:231:VAL:O	2.19	0.43
1:B:420:SER:HB3	1:B:423:PHE:CZ	2.54	0.43
1:A:310:ARG:HG3	1:A:478:ALA:HB1	2.01	0.43
1:B:198:MET:HA	1:B:231:VAL:O	2.19	0.43
1:D:59:ASP:HB2	1:D:64:PHE:CD2	2.53	0.43
1:B:14:MET:HE2	1:B:39:ILE:HD11	2.01	0.43
1:B:322:ASN:ND2	1:B:327:LEU:HB2	2.33	0.43
1:C:568:GLU:HG2	1:D:601:VAL:HA	2.01	0.43
1:D:463:LEU:HD22	1:D:542:LEU:HD11	2.01	0.43
1:A:87:ILE:HG13	1:A:90:MET:HE2	2.00	0.42
1:C:560:PRO:HB3	1:C:584:TYR:CD1	2.54	0.42
1:A:10:LEU:O	1:A:14:MET:HG3	2.20	0.42
1:B:29:TRP:HZ2	1:B:52:ARG:HH22	1.66	0.42
1:D:420:SER:HB3	1:D:423:PHE:CZ	2.55	0.42
1:A:107:LEU:HA	1:A:152:SER:HB3	2.01	0.42
1:D:138:ALA:HB2	1:D:188:ALA:HA	2.01	0.42
1:A:614:VAL:HB	1:A:624:LEU:HD21	2.02	0.42
1:D:76:VAL:HB	1:D:123:PHE:HA	2.02	0.42
1:B:381:GLN:HB3	1:B:394:LEU:HD22	2.01	0.42
1:D:156:PHE:CG	1:D:156:PHE:O	2.73	0.42
1:B:5:ARG:HE	1:B:5:ARG:HB2	1.56	0.42
1:B:352:SER:HB3	2:B:802:HTG:H1	2.01	0.42
1:D:509:PRO:HB3	1:D:533:TYR:CG	2.55	0.42
1:A:396:GLU:O	1:A:400:GLN:HG2	2.19	0.42
1:A:678:VAL:HG11	1:D:267:GLU:HB3	2.02	0.42
1:B:234:TRP:HA	1:B:252:MET:O	2.20	0.42
1:C:568:GLU:HG3	1:D:602:ILE:HG13	2.01	0.42
1:A:172:VAL:HG22	1:A:523:TYR:CG	2.55	0.41
1:C:790:GLU:CD	1:D:416:LYS:HB2	2.45	0.41
1:A:189:VAL:HA	1:A:194:PRO:HD2	2.01	0.41
1:D:395:THR:O	1:D:399:LEU:HG	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:LEU:HD12	1:A:153:LEU:HA	1.88	0.41
1:A:556:TYR:CG	1:A:662:VAL:HG21	2.55	0.41
1:B:310:ARG:HG3	1:B:478:ALA:HB1	2.01	0.41
1:B:556:TYR:HE2	1:B:590:VAL:HG13	1.85	0.41
1:C:246:TRP:O	1:C:279:GLY:HA3	2.20	0.41
1:C:189:ALA:HB2	1:C:239:ALA:HA	2.03	0.41
1:C:471:SER:HB3	1:C:474:PHE:CZ	2.55	0.41
1:D:367:LEU:HD21	1:D:487:LEU:HD12	2.02	0.41
1:A:19:LYS:HG2	1:A:280:VAL:HG21	2.02	0.41
1:A:719:LEU:HD22	1:A:723:LYS:HZ2	1.85	0.41
1:B:346:ARG:HH22	1:B:419:GLU:HB2	1.86	0.41
1:C:399:GLN:HG2	1:C:400:GLY:N	2.35	0.41
1:A:157:ALA:HB3	1:A:200:SER:OG	2.20	0.41
1:A:697:PHE:CE2	1:A:747:LEU:HD22	2.56	0.41
1:B:633:PHE:HB3	1:B:650:PHE:CE2	2.56	0.41
1:D:153:LEU:HD21	1:D:185:PHE:CD2	2.56	0.41
1:D:496:LYS:HG3	1:D:589:ILE:HD12	2.03	0.41
1:B:384:ASP:HB2	1:B:387:SER:HB3	2.03	0.41
1:D:52:ARG:HE	1:D:62:GLY:HA2	1.85	0.41
1:D:509:PRO:HB3	1:D:533:TYR:CD1	2.56	0.41
1:B:138:ALA:HB2	1:B:188:ALA:HA	2.02	0.41
1:A:33:PRO:HB3	1:A:41:SER:HB2	2.04	0.40
1:A:234:TRP:HA	1:A:252:MET:O	2.20	0.40
1:B:454:ASN:O	1:B:473:TYR:HA	2.21	0.40
1:C:663:ASP:HB3	1:C:675:LEU:HD23	2.03	0.40
1:D:303:GLU:HG2	1:D:359:ARG:HH12	1.87	0.40
1:C:665:VAL:HG11	1:C:675:LEU:HD11	2.03	0.40
2:C:901:HTG:H2'1	2:C:901:HTG:H1	1.88	0.40
1:C:291:ARG:HD3	1:C:302:GLU:OE1	2.21	0.40
1:D:719:LEU:HD23	1:D:719:LEU:HA	1.93	0.40
1:A:44:MET:HE2	1:A:234:TRP:CH2	2.55	0.40
1:A:381:GLN:HB3	1:A:394:LEU:HD22	2.04	0.40
1:D:10:LEU:O	1:D:14:MET:HG3	2.21	0.40
1:D:189:VAL:HA	1:D:194:PRO:HD2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	726/780 (93%)	701 (97%)	24 (3%)	1 (0%)	48	76
1	B	734/780 (94%)	706 (96%)	28 (4%)	0	100	100
1	C	732/780 (94%)	701 (96%)	31 (4%)	0	100	100
1	D	737/780 (94%)	710 (96%)	25 (3%)	2 (0%)	36	63
All	All	2929/3120 (94%)	2818 (96%)	108 (4%)	3 (0%)	48	76

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	233	ASP
1	A	47	GLY
1	D	47	GLY

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	607/648 (94%)	607 (100%)	0	100	100
1	B	613/648 (95%)	613 (100%)	0	100	100
1	C	611/648 (94%)	611 (100%)	0	100	100
1	D	614/648 (95%)	614 (100%)	0	100	100
All	All	2445/2592 (94%)	2445 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	49	HIS
1	A	147	GLN
1	A	173	ASN
1	B	147	GLN
1	B	160	ASN
1	B	354	GLN
1	C	345	HIS
1	C	405	GLN
1	C	451	GLN
1	C	528	GLN
1	C	564	ASN
1	C	598	HIS
1	D	155	HIS
1	D	160	ASN
1	D	354	GLN
1	D	400	GLN
1	D	547	HIS

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 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	EDO	C	903	-	3,3,3	0.25	0	2,2,2	0.34	0
3	SO4	A	802	-	4,4,4	0.67	0	6,6,6	0.08	0
3	SO4	C	902	-	4,4,4	0.67	0	6,6,6	0.08	0
3	SO4	D	801	-	4,4,4	0.67	0	6,6,6	0.07	0
2	HTG	B	802	-	19,19,19	0.55	0	23,24,24	0.95	1 (4%)
5	GOL	D	802	-	5,5,5	0.34	0	5,5,5	0.41	0
4	EDO	B	801	-	3,3,3	0.25	0	2,2,2	0.33	0
2	HTG	C	901	-	19,19,19	0.54	0	23,24,24	0.83	0
2	HTG	A	801	-	19,19,19	0.55	0	23,24,24	1.08	1 (4%)

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	EDO	C	903	-	-	0/1/1/1	-
5	GOL	D	802	-	-	1/4/4/4	-
2	HTG	B	802	-	-	3/10/30/30	0/1/1/1
4	EDO	B	801	-	-	0/1/1/1	-
2	HTG	C	901	-	-	7/10/30/30	0/1/1/1
2	HTG	A	801	-	-	2/10/30/30	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HTG	C1-C2-C3	-3.19	104.31	110.55
2	B	802	HTG	C1-C2-C3	-2.58	105.50	110.55

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	901	HTG	O5-C1-S1-C1'
2	C	901	HTG	O5-C5-C6-O6

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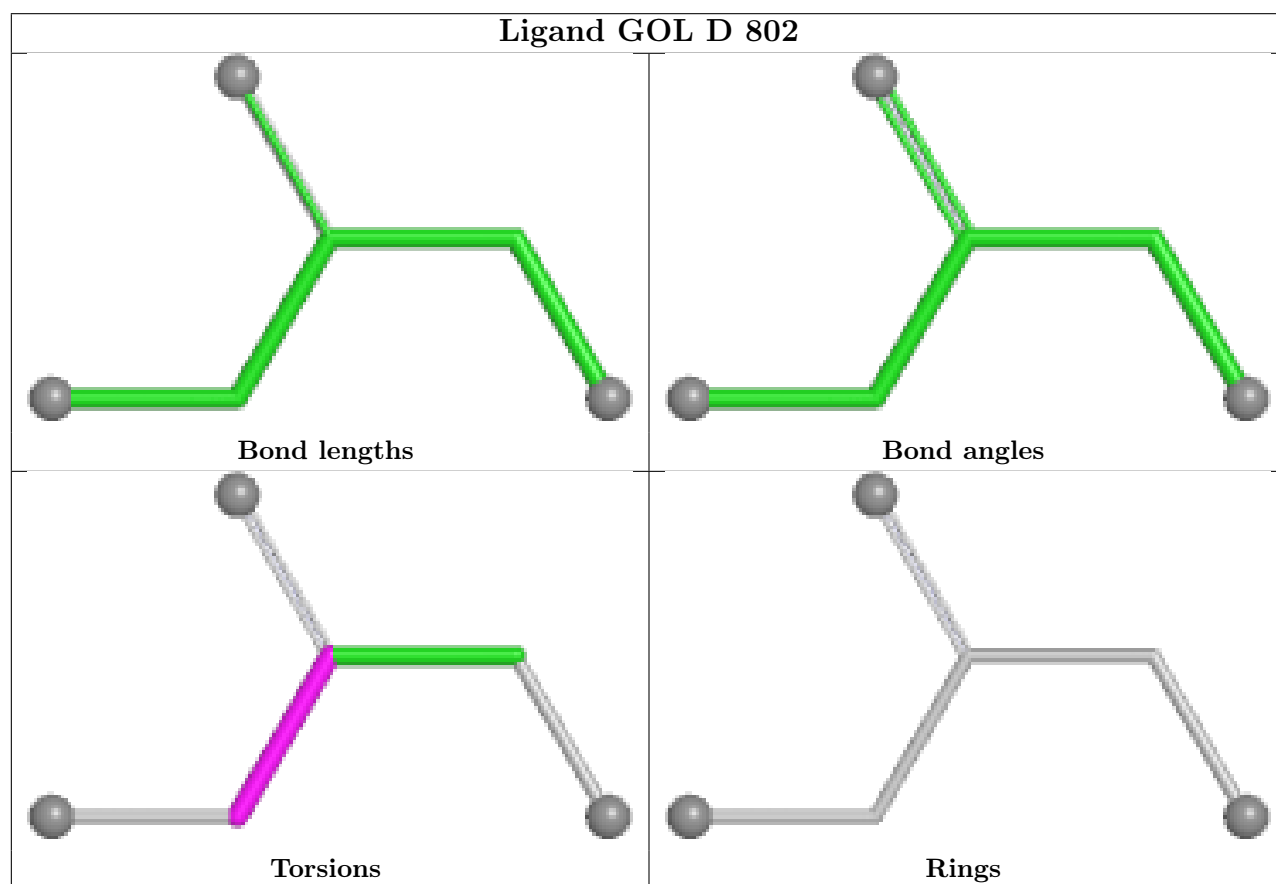
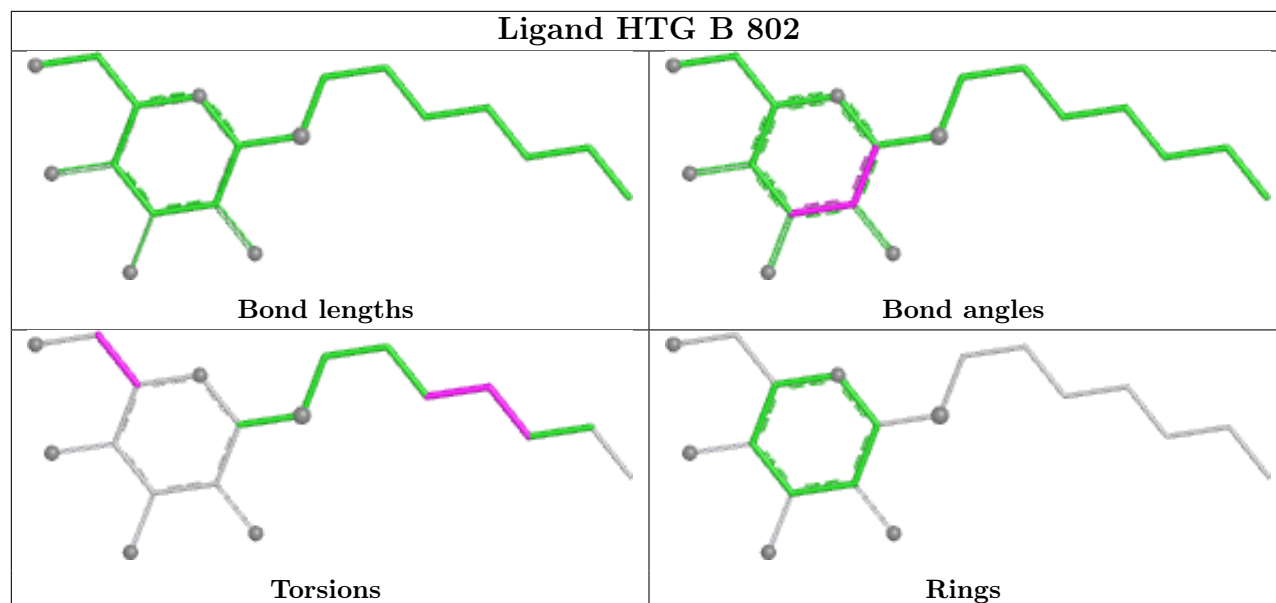
Mol	Chain	Res	Type	Atoms
2	C	901	HTG	C4-C5-C6-O6
2	A	801	HTG	O5-C5-C6-O6
2	C	901	HTG	C1'-C2'-C3'-C4'
2	B	802	HTG	C3'-C4'-C5'-C6'
2	C	901	HTG	S1-C1'-C2'-C3'
2	B	802	HTG	O5-C5-C6-O6
2	A	801	HTG	S1-C1'-C2'-C3'
2	B	802	HTG	C2'-C3'-C4'-C5'
2	C	901	HTG	C2-C1-S1-C1'
5	D	802	GOL	O2-C2-C3-O3
2	C	901	HTG	C4'-C5'-C6'-C7'

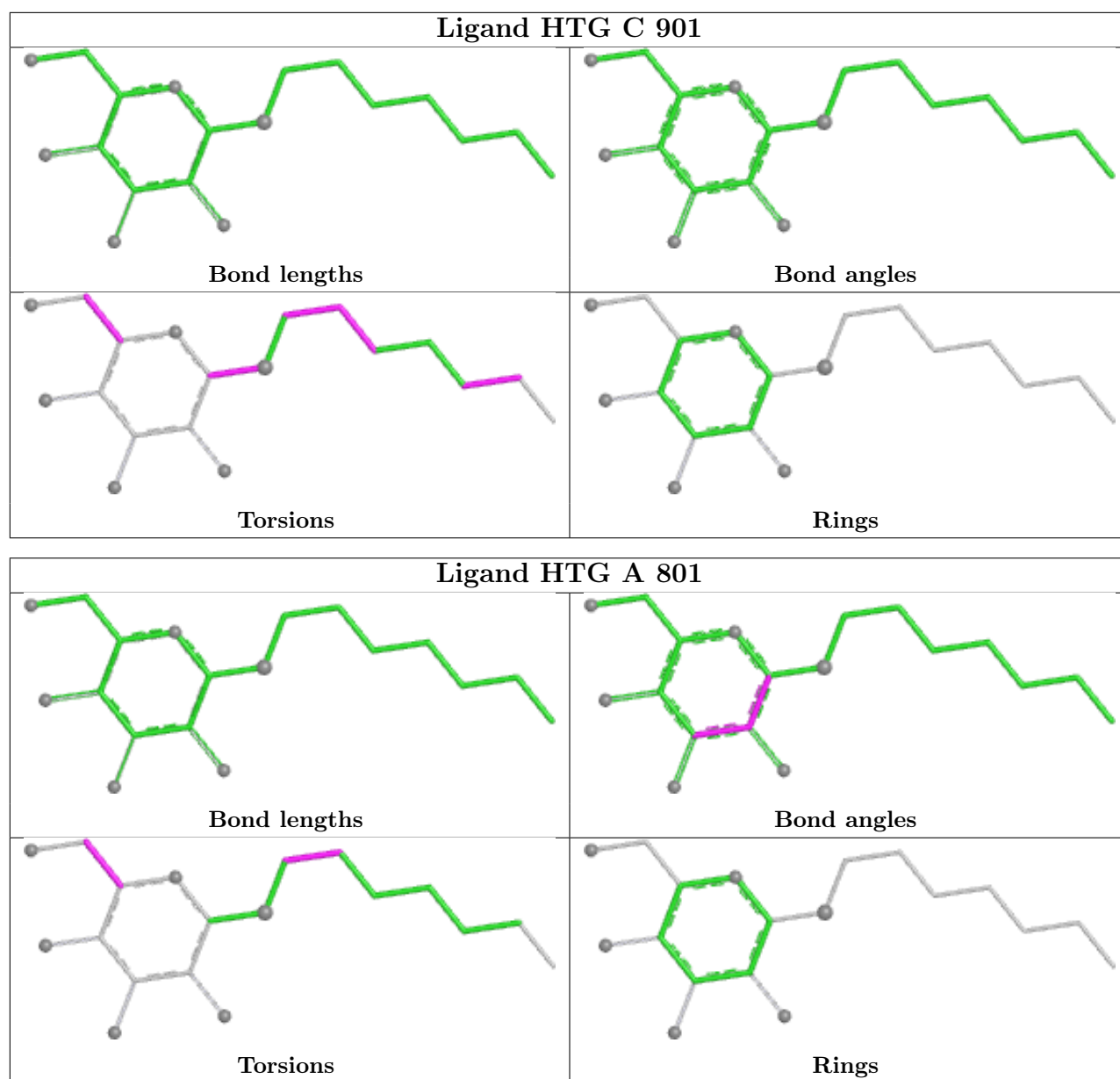
There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	903	EDO	1	0
2	B	802	HTG	1	0
2	C	901	HTG	3	0
2	A	801	HTG	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	732/780 (93%)	0.83	49 (6%) 24 18	12, 26, 50, 84	0
1	B	740/780 (94%)	0.91	58 (7%) 19 15	11, 26, 55, 95	0
1	C	738/780 (94%)	0.85	48 (6%) 25 18	12, 27, 54, 78	0
1	D	741/780 (95%)	0.82	58 (7%) 19 15	12, 24, 55, 80	0
All	All	2951/3120 (94%)	0.85	213 (7%) 21 17	11, 26, 54, 95	0

All (213) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	703	GLY	5.1
1	B	716	ALA	4.6
1	C	22	ASP	4.5
1	B	708	GLY	4.4
1	D	677	ILE	4.3
1	A	352	SER	4.2
1	A	373	GLU	4.2
1	D	311	GLU	4.2
1	B	715	TYR	4.1
1	C	125	SER	4.0
1	A	702	ARG	4.0
1	B	699	GLN	4.0
1	A	369	ALA	4.0
1	C	791	TYR	3.9
1	A	715	TYR	3.9
1	B	353	SER	3.7
1	A	721	ALA	3.7
1	C	404	SER	3.7
1	B	152	SER	3.7
1	C	792	ALA	3.6
1	A	704	MET	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	742	GLU	3.5
1	B	59	ASP	3.5
1	B	392	ALA	3.5
1	C	796	GLU	3.5
1	D	352	SER	3.4
1	D	60	HIS	3.4
1	C	439	ASP	3.3
1	D	59	ASP	3.3
1	A	716	ALA	3.3
1	B	707	PHE	3.3
1	B	697	PHE	3.3
1	B	701	LEU	3.3
1	D	61	LEU	3.3
1	D	296	GLU	3.3
1	C	107	ALA	3.2
1	D	62	GLY	3.2
1	B	57	ALA	3.2
1	B	686	ALA	3.2
1	B	274	GLU	3.2
1	D	388	ASP	3.2
1	B	419	GLU	3.1
1	B	388	ASP	3.1
1	D	313	ALA	3.1
1	A	718	MET	3.1
1	C	750	GLN	3.0
1	A	705	ILE	3.0
1	D	704	MET	3.0
1	C	784	GLY	3.0
1	B	56	ALA	3.0
1	B	64	PHE	3.0
1	B	58	SER	3.0
1	A	384	ASP	3.0
1	D	697	PHE	2.9
1	A	419	GLU	2.9
1	B	352	SER	2.9
1	A	518	GLY	2.9
1	D	420	SER	2.9
1	B	65	ASP	2.9
1	B	233	ASP	2.9
1	D	234	TRP	2.9
1	D	715	TYR	2.9
1	B	157	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	296	GLU	2.8
1	A	699	GLN	2.8
1	B	725	SER	2.8
1	C	725	SER	2.8
1	D	419	GLU	2.8
1	C	609	ASN	2.8
1	B	51	LEU	2.8
1	C	161	GLY	2.8
1	D	736	ALA	2.8
1	A	199	CYS	2.8
1	A	701	LEU	2.8
1	B	475	GLY	2.7
1	A	719	LEU	2.7
1	D	705	ILE	2.7
1	B	15	THR	2.7
1	D	668	MET	2.7
1	D	58	SER	2.7
1	A	155	HIS	2.7
1	B	705	ILE	2.7
1	D	703	GLY	2.7
1	D	719	LEU	2.6
1	B	718	MET	2.6
1	D	93	ALA	2.6
1	C	471	SER	2.6
1	B	704	MET	2.6
1	D	157	ALA	2.6
1	B	350	SER	2.6
1	C	599	GLY	2.6
1	D	24	THR	2.6
1	D	154	LYS	2.6
1	D	302	PRO	2.6
1	C	793	GLU	2.6
1	D	402	ALA	2.5
1	B	698	ALA	2.5
1	C	420	ALA	2.5
1	C	435	ASP	2.5
1	C	800	GLU	2.5
1	D	699	GLN	2.5
1	B	724	GLU	2.5
1	C	264	GLU	2.5
1	A	157	ALA	2.5
1	D	353	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	728	ILE	2.5
1	B	160	ASN	2.5
1	A	120	GLY	2.5
1	B	714	GLU	2.4
1	D	213	GLU	2.4
1	A	361	ASP	2.4
1	C	31	THR	2.4
1	B	393	VAL	2.4
1	A	56	ALA	2.4
1	A	392	ALA	2.4
1	C	213	GLU	2.4
1	C	748	PHE	2.4
1	B	18	GLU	2.4
1	C	468	ARG	2.4
1	B	81	SER	2.4
1	D	555	ALA	2.4
1	B	156	PHE	2.4
1	B	22	LEU	2.4
1	A	623	THR	2.4
1	B	623	THR	2.4
1	C	106	LYS	2.4
1	D	718	MET	2.4
1	A	503	VAL	2.4
1	B	279	ALA	2.3
1	B	721	ALA	2.3
1	C	284	ASP	2.3
1	A	724	GLU	2.3
1	D	374	HIS	2.3
1	C	781	ILE	2.3
1	B	20	ALA	2.3
1	B	369	ALA	2.3
1	A	697	PHE	2.3
1	C	116	ASP	2.3
1	C	829	GLY	2.3
1	D	436	ALA	2.3
1	B	719	LEU	2.3
1	D	155	HIS	2.3
1	A	400	GLN	2.3
1	B	702	ARG	2.3
1	A	76	VAL	2.3
1	C	400	GLY	2.2
1	D	701	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	15	THR	2.2
1	D	706	PRO	2.2
1	D	721	ALA	2.2
1	A	156	PHE	2.2
1	B	181	TYR	2.2
1	A	311	GLU	2.2
1	B	158	ALA	2.2
1	C	453	ALA	2.2
1	D	342	ALA	2.2
1	C	69	LYS	2.2
1	A	309	ALA	2.2
1	A	336	ALA	2.2
1	A	686	ALA	2.2
1	D	369	ALA	2.2
1	D	481	GLY	2.2
1	A	612	ASP	2.2
1	A	89	ARG	2.2
1	C	336	THR	2.2
1	C	676	THR	2.2
1	D	74	SER	2.2
1	C	67	LEU	2.2
1	B	717	GLU	2.2
1	D	158	ALA	2.2
1	A	548	GLY	2.2
1	C	416	PHE	2.2
1	C	190	HIS	2.2
1	A	191	LYS	2.1
1	A	342	ALA	2.1
1	A	256	GLY	2.1
1	B	258	GLY	2.1
1	D	256	GLY	2.1
1	D	57	ALA	2.1
1	D	585	ALA	2.1
1	B	580	ASN	2.1
1	D	160	ASN	2.1
1	C	108	ALA	2.1
1	A	351	GLY	2.1
1	A	598	GLU	2.1
1	D	25	GLY	2.1
1	D	733	SER	2.1
1	A	485	ASP	2.1
1	C	191	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	358	THR	2.1
1	C	208	ALA	2.1
1	D	333	GLY	2.1
1	B	35	GLU	2.1
1	A	80	SER	2.1
1	C	799	LYS	2.1
1	D	264	GLU	2.1
1	D	726	VAL	2.0
1	B	118	LEU	2.0
1	C	753	ARG	2.0
1	D	577	ASN	2.0
1	C	357	HIS	2.0
1	D	351	GLY	2.0
1	C	790	GLU	2.0
1	C	444	VAL	2.0
1	A	420	SER	2.0
1	D	355	ILE	2.0
1	D	100	ALA	2.0
1	C	54	GLY	2.0
1	B	53	LYS	2.0
1	C	396	PRO	2.0
1	A	444	VAL	2.0
1	A	499	GLU	2.0
1	B	124	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](#)

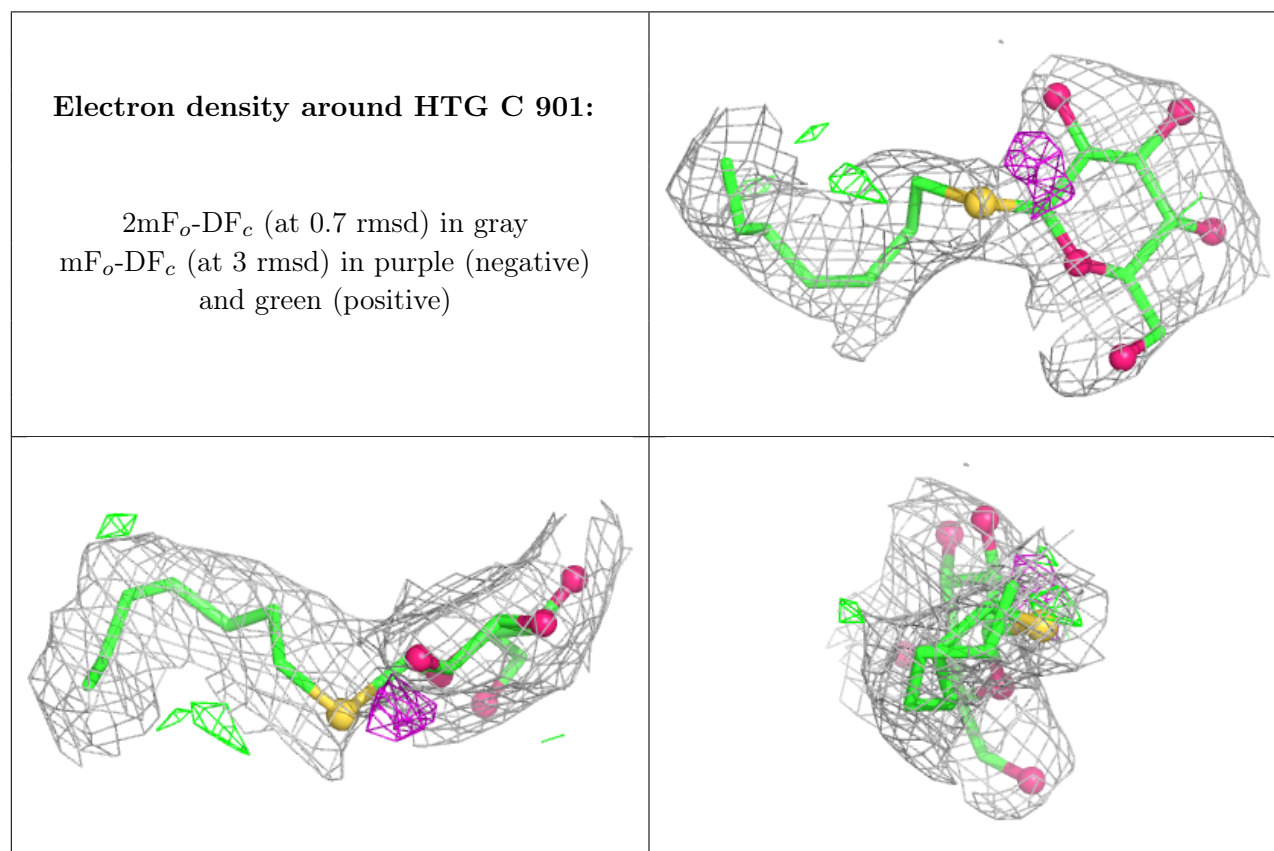
There are no oligosaccharides in this entry.

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.

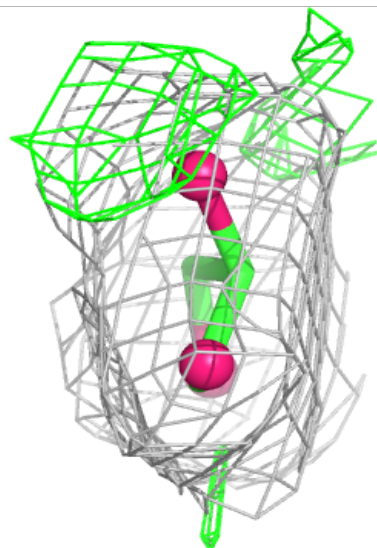
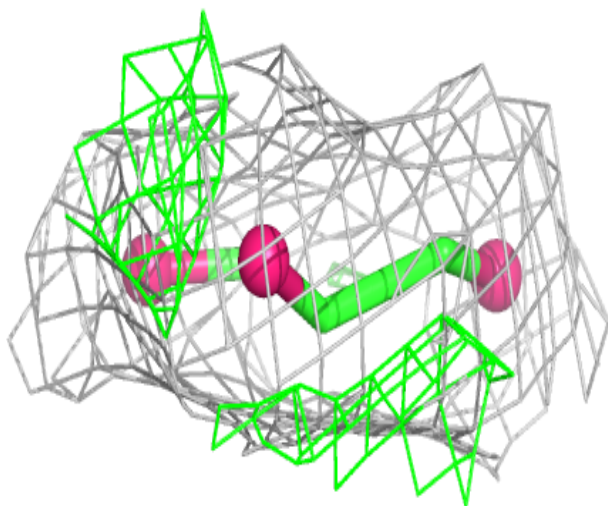
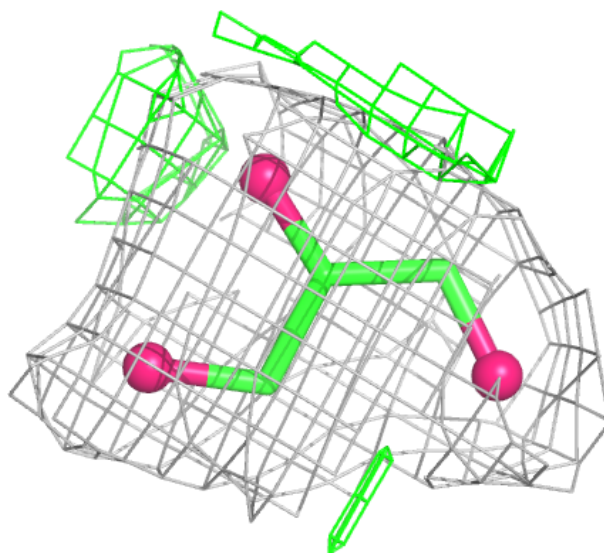
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	HTG	C	901	19/19	0.73	0.22	18,36,54,60	0
3	SO4	C	902	5/5	0.73	0.21	56,71,83,89	0
5	GOL	D	802	6/6	0.78	0.27	20,30,39,42	0
2	HTG	B	802	19/19	0.79	0.23	25,40,51,63	0
3	SO4	A	802	5/5	0.80	0.19	36,56,68,104	0
2	HTG	A	801	19/19	0.82	0.20	29,40,54,61	0
4	EDO	B	801	4/4	0.85	0.18	3,7,13,13	0
3	SO4	D	801	5/5	0.91	0.17	33,44,57,70	0
4	EDO	C	903	4/4	0.93	0.14	2,4,5,9	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.



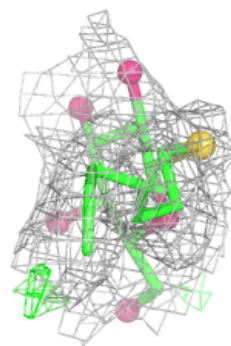
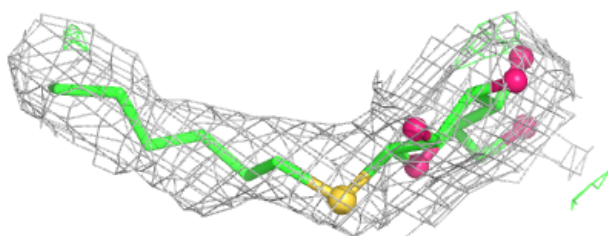
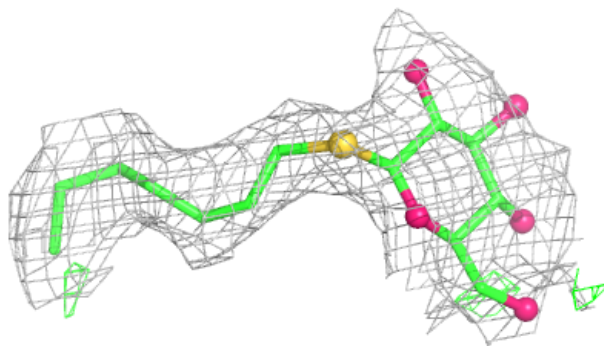
Electron density around GOL D 802:

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

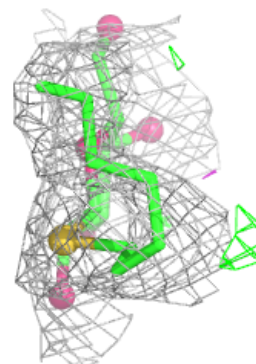
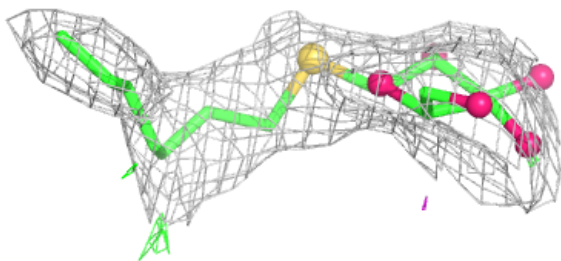
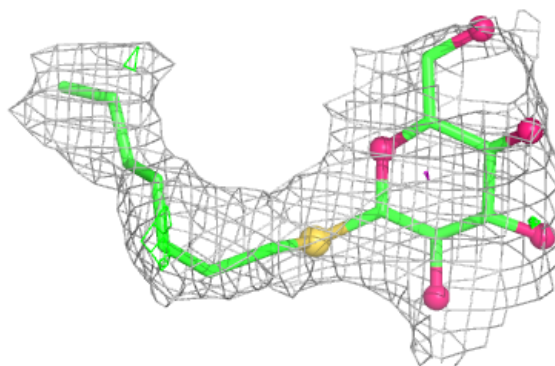


Electron density around HTG B 802:

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

**Electron density around HTG 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.