



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2025 – 01:15 pm BST

PDB ID : 9G4N / pdb_00009g4n
Title : Glycoside Hydrolase Family 157 from Labilicabulum antarcticum (LaGH157)
E224A mutant in complex with Laminaritriose and Glucose
Authors : Caseiro, C.; Alves, V.D.; Carvalho, A.L.; Bule, P.
Deposited on : 2024-07-15
Resolution : 2.32 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 Phenix2.0rc1
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

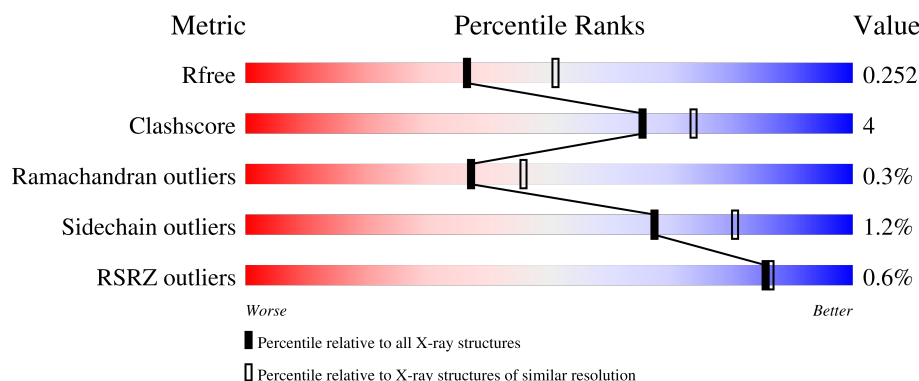
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.32 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.




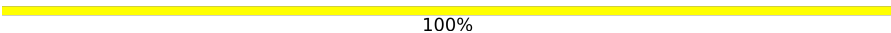

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	7250 (2.34-2.30)
Clashscore	180529	8063 (2.34-2.30)
Ramachandran outliers	177936	7993 (2.34-2.30)
Sidechain outliers	177891	7993 (2.34-2.30)
RSRZ outliers	164620	7250 (2.34-2.30)

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	537	 88% 9% ..
1	B	537	 89% 8% ..
1	C	537	 87% 11% ..
1	D	537	 89% 9% .
2	E	3	 67% 33%

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Mol	Chain	Length	Quality of chain
2	F	3	 33% 67%
2	G	3	 100%
2	H	3	 67% 33%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	A	601	-	-	X	-
3	EDO	B	604	-	-	X	-
3	EDO	C	601	-	-	X	-
3	EDO	D	603	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 18186 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 Glycoside hydrolase family 2 catalytic domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	528	Total	C	N	O	S	0	0	0
			4245	2717	685	828	15			
1	B	526	Total	C	N	O	S	0	0	0
			4231	2708	683	825	15			
1	C	527	Total	C	N	O	S	0	0	0
			4237	2711	684	827	15			
1	D	527	Total	C	N	O	S	0	1	0
			4242	2714	685	828	15			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A1Y1CQ89
A	224	ALA	GLU	engineered mutation	UNP A0A1Y1CQ89
A	530	LEU	-	expression tag	UNP A0A1Y1CQ89
A	531	GLU	-	expression tag	UNP A0A1Y1CQ89
A	532	HIS	-	expression tag	UNP A0A1Y1CQ89
A	533	HIS	-	expression tag	UNP A0A1Y1CQ89
A	534	HIS	-	expression tag	UNP A0A1Y1CQ89
A	535	HIS	-	expression tag	UNP A0A1Y1CQ89
A	536	HIS	-	expression tag	UNP A0A1Y1CQ89
A	537	HIS	-	expression tag	UNP A0A1Y1CQ89
B	1	MET	-	initiating methionine	UNP A0A1Y1CQ89
B	224	ALA	GLU	engineered mutation	UNP A0A1Y1CQ89
B	530	LEU	-	expression tag	UNP A0A1Y1CQ89
B	531	GLU	-	expression tag	UNP A0A1Y1CQ89
B	532	HIS	-	expression tag	UNP A0A1Y1CQ89
B	533	HIS	-	expression tag	UNP A0A1Y1CQ89
B	534	HIS	-	expression tag	UNP A0A1Y1CQ89
B	535	HIS	-	expression tag	UNP A0A1Y1CQ89
B	536	HIS	-	expression tag	UNP A0A1Y1CQ89
B	537	HIS	-	expression tag	UNP A0A1Y1CQ89

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1	MET	-	initiating methionine	UNP A0A1Y1CQ89
C	224	ALA	GLU	engineered mutation	UNP A0A1Y1CQ89
C	530	LEU	-	expression tag	UNP A0A1Y1CQ89
C	531	GLU	-	expression tag	UNP A0A1Y1CQ89
C	532	HIS	-	expression tag	UNP A0A1Y1CQ89
C	533	HIS	-	expression tag	UNP A0A1Y1CQ89
C	534	HIS	-	expression tag	UNP A0A1Y1CQ89
C	535	HIS	-	expression tag	UNP A0A1Y1CQ89
C	536	HIS	-	expression tag	UNP A0A1Y1CQ89
C	537	HIS	-	expression tag	UNP A0A1Y1CQ89
D	1	MET	-	initiating methionine	UNP A0A1Y1CQ89
D	224	ALA	GLU	engineered mutation	UNP A0A1Y1CQ89
D	530	LEU	-	expression tag	UNP A0A1Y1CQ89
D	531	GLU	-	expression tag	UNP A0A1Y1CQ89
D	532	HIS	-	expression tag	UNP A0A1Y1CQ89
D	533	HIS	-	expression tag	UNP A0A1Y1CQ89
D	534	HIS	-	expression tag	UNP A0A1Y1CQ89
D	535	HIS	-	expression tag	UNP A0A1Y1CQ89
D	536	HIS	-	expression tag	UNP A0A1Y1CQ89
D	537	HIS	-	expression tag	UNP A0A1Y1CQ89

- Molecule 2 is an oligosaccharide called beta-D-glucopyranose-(1-3)-beta-D-glucopyranose-(1-3)-beta-D-glucopyranose.



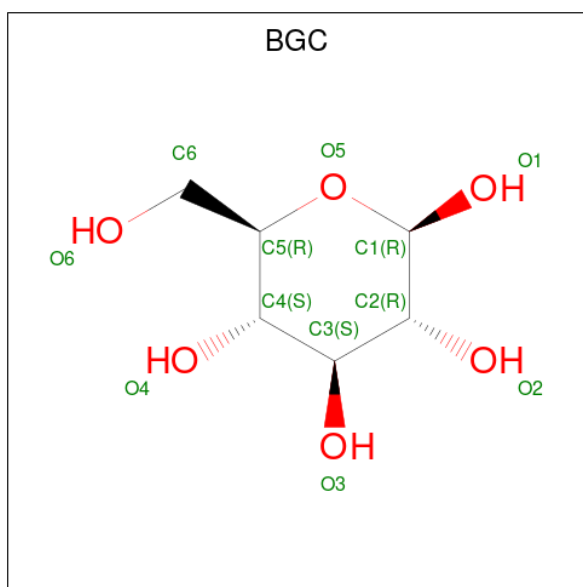
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	3	Total	C	O	0	0	0
			34	18	16			
2	F	3	Total	C	O	0	0	0
			34	18	16			
2	G	3	Total	C	O	0	0	0
			34	18	16			
2	H	3	Total	C	O	0	0	0
			34	18	16			

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



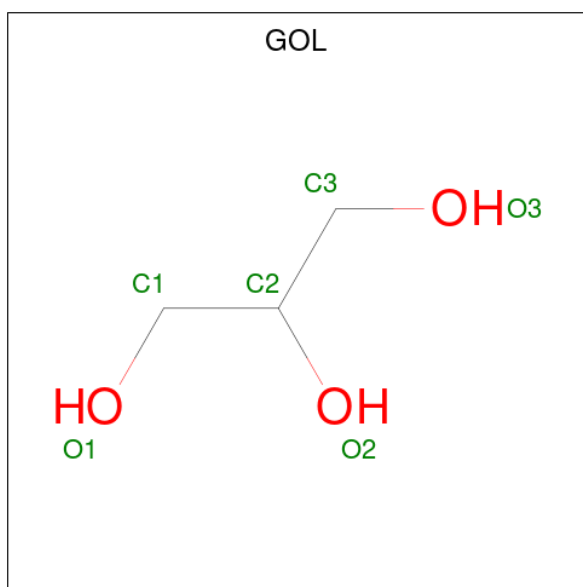
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			12	6	6		
4	B	1	Total	C	O	0	0
			12	6	6		
4	D	1	Total	C	O	0	0
			12	6	6		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



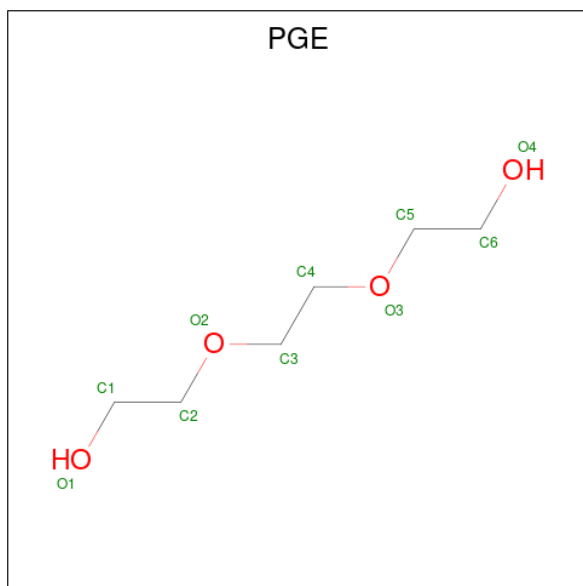
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			10	6	4		

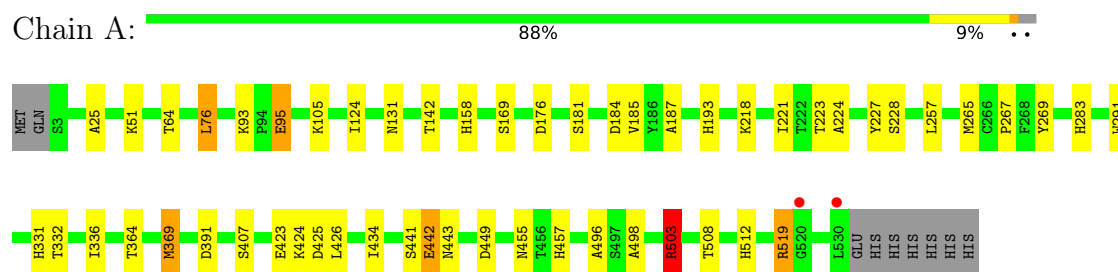
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	275	Total 275	O 275	0	0
7	B	242	Total 242	O 242	0	0
7	C	188	Total 188	O 188	0	0
7	D	264	Total 264	O 264	0	0

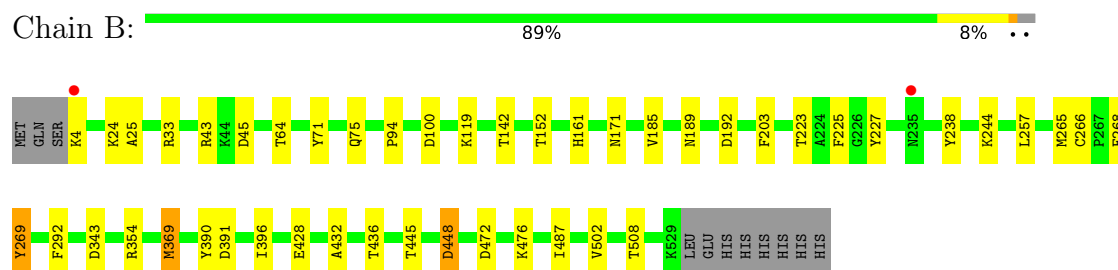
3 Residue-property plots

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.

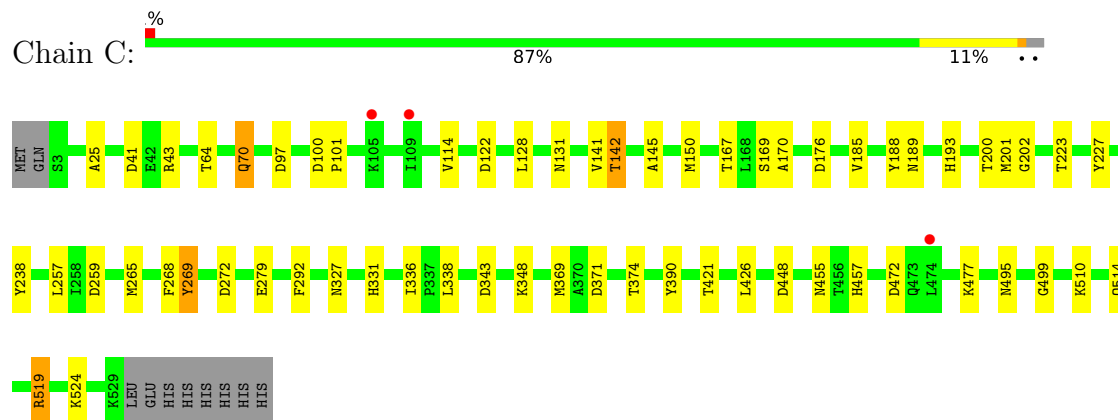
- Molecule 1: Glycoside hydrolase family 2 catalytic domain-containing protein



- Molecule 1: Glycoside hydrolase family 2 catalytic domain-containing protein

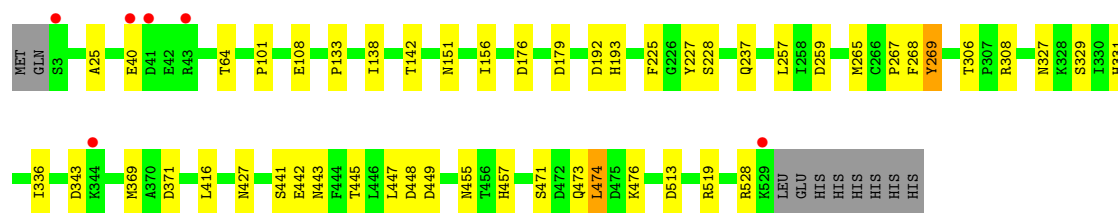


- Molecule 1: Glycoside hydrolase family 2 catalytic domain-containing protein



- Molecule 1: Glycoside hydrolase family 2 catalytic domain-containing protein





- Molecule 2: beta-D-glucopyranose-(1-3)-beta-D-glucopyranose-(1-3)-beta-D-glucopyranose

Chain E: 67% 33%



- Molecule 2: beta-D-glucopyranose-(1-3)-beta-D-glucopyranose-(1-3)-beta-D-glucopyranose

Chain F: 33% 67%



- Molecule 2: beta-D-glucopyranose-(1-3)-beta-D-glucopyranose-(1-3)-beta-D-glucopyranose

Chain G: 100%



- Molecule 2: beta-D-glucopyranose-(1-3)-beta-D-glucopyranose-(1-3)-beta-D-glucopyranose

Chain H: 67% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	114.85Å 127.55Å 171.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.71 – 2.32 47.71 – 2.32	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.71-2.32) 99.9 (47.71-2.32)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.32Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.205 , 0.247 0.212 , 0.252	Depositor DCC
R_{free} test set	5276 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.546	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 31.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	18186	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.50% 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: PGE, BGC, GOL, 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	0.66	0/4344	1.13	4/5874 (0.1%)
1	B	0.67	0/4330	1.12	12/5855 (0.2%)
1	C	0.65	0/4336	1.14	11/5863 (0.2%)
1	D	0.67	0/4344	1.14	6/5874 (0.1%)
All	All	0.66	0/17354	1.13	33/23466 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	4

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	421	THR	CA-CB-OG1	-7.28	98.68	109.60
1	D	371	ASP	CA-CB-CG	6.93	119.53	112.60
1	B	343	ASP	CA-CB-CG	6.66	119.26	112.60
1	C	374	THR	CA-CB-OG1	-6.60	99.70	109.60
1	B	448	ASP	CA-CB-CG	6.50	119.10	112.60
1	A	425	ASP	CA-CB-CG	6.28	118.88	112.60
1	C	448	ASP	CA-CB-CG	6.20	118.80	112.60
1	D	343	ASP	CA-CB-CG	6.09	118.69	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	391	ASP	CA-CB-CG	6.07	118.67	112.60
1	C	343	ASP	CA-CB-CG	6.06	118.66	112.60
1	C	122	ASP	CA-CB-CG	5.98	118.58	112.60
1	B	445	THR	CA-CB-OG1	-5.96	100.66	109.60
1	D	306	THR	CA-CB-OG1	-5.79	100.92	109.60
1	C	272	ASP	CA-CB-CG	5.72	118.33	112.60
1	C	472	ASP	CA-CB-CG	5.72	118.32	112.60
1	D	449	ASP	CA-CB-CG	5.70	118.30	112.60
1	C	41	ASP	CA-CB-CG	5.68	118.28	112.60
1	B	472	ASP	CA-CB-CG	5.65	118.25	112.60
1	D	445	THR	CA-CB-OG1	-5.62	101.17	109.60
1	B	152	THR	CA-CB-OG1	-5.59	101.21	109.60
1	C	524	LYS	N-CA-CB	5.52	118.72	110.22
1	B	94	PRO	N-CA-C	5.36	120.81	113.84
1	A	442	GLU	CB-CA-C	5.35	119.38	109.54
1	B	100	ASP	CA-CB-CG	5.33	117.93	112.60
1	A	95	GLU	CB-CA-C	5.31	118.89	109.65
1	C	97	ASP	CA-CB-CG	5.31	117.91	112.60
1	B	436	THR	CA-CB-OG1	-5.30	101.66	109.60
1	B	45	ASP	CA-CB-CG	5.25	117.86	112.60
1	D	513	ASP	CA-CB-CG	5.23	117.83	112.60
1	C	202	GLY	CA-C-O	-5.12	118.14	122.29
1	B	225	PHE	CA-CB-CG	5.06	118.86	113.80
1	A	51	LYS	N-CA-CB	5.03	117.31	110.01
1	B	428	GLU	CB-CG-CD	5.03	121.15	112.60

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	503	ARG	Sidechain
1	B	33	ARG	Sidechain
1	C	43	ARG	Sidechain
1	D	308	ARG	Sidechain

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	4245	0	4138	36	0
1	B	4231	0	4122	23	0
1	C	4237	0	4127	36	0
1	D	4242	0	4133	30	0
2	E	34	0	30	0	0
2	F	34	0	30	0	0
2	G	34	0	30	1	0
2	H	34	0	30	0	0
3	A	4	0	6	7	0
3	B	8	0	12	5	0
3	C	4	0	6	9	0
3	D	4	0	6	4	0
4	A	12	0	12	0	0
4	B	12	0	12	0	0
4	D	12	0	12	0	0
5	A	6	0	8	0	0
5	B	18	0	24	1	0
5	C	6	0	8	0	0
5	D	30	0	40	2	0
6	C	10	0	14	0	0
7	A	275	0	0	2	0
7	B	242	0	0	4	0
7	C	188	0	0	1	0
7	D	264	0	0	0	0
All	All	18186	0	16800	124	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 (124) 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:141:VAL:O	1:C:142:THR:HG22	1.69	0.92
1:A:227:TYR:HA	3:A:601:EDO:H11	1.56	0.88
1:B:238:TYR:HA	3:B:604:EDO:H11	1.68	0.76
1:B:43:ARG:HD3	7:B:807:HOH:O	1.90	0.70
1:C:142:THR:HG23	1:C:145:ALA:H	1.58	0.66
1:B:292:PHE:CE1	3:B:604:EDO:H21	2.31	0.64
1:A:176:ASP:OD1	1:A:457:HIS:HE1	1.81	0.64
1:D:227:TYR:HA	3:D:603:EDO:H22	1.81	0.63
1:C:176:ASP:OD1	1:C:457:HIS:HE1	1.82	0.62
1:A:193:HIS:HE1	3:A:601:EDO:H12	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:292:PHE:CE1	3:C:601:EDO:H12	2.36	0.60
1:D:176:ASP:OD1	1:D:457:HIS:HE1	1.86	0.58
1:C:227:TYR:CD1	3:C:601:EDO:H21	2.39	0.57
1:D:427:ASN:ND2	1:D:519:ARG:HH11	2.01	0.57
1:B:257:LEU:HB2	1:B:265:MET:HE1	1.88	0.56
1:A:193:HIS:HE1	3:A:601:EDO:C1	2.18	0.56
1:A:193:HIS:CE1	3:A:601:EDO:H12	2.42	0.55
1:D:331:HIS:CD2	1:D:336:ILE:HD12	2.42	0.55
1:A:158:HIS:HE1	1:A:184:ASP:OD2	1.89	0.54
1:C:331:HIS:CD2	1:C:336:ILE:HD12	2.43	0.54
1:A:105:LYS:HG2	1:D:101:PRO:HG3	1.88	0.54
1:B:390:TYR:CE2	1:B:396:ILE:HG13	2.42	0.53
1:A:227:TYR:HA	3:A:601:EDO:C1	2.35	0.53
1:C:141:VAL:O	1:C:142:THR:CG2	2.50	0.53
1:B:354:ARG:HA	5:B:605:GOL:H11	1.89	0.53
1:B:244:LYS:HD3	7:B:940:HOH:O	2.09	0.52
1:C:193:HIS:HE1	3:C:601:EDO:O1	1.93	0.52
1:B:25:ALA:HA	1:B:64:THR:O	2.10	0.51
1:D:193:HIS:NE2	3:D:603:EDO:H21	2.25	0.51
1:D:455:ASN:O	1:D:457:HIS:HD2	1.93	0.51
1:D:471:SER:HA	1:D:474:LEU:CD2	2.41	0.51
1:C:268:PHE:HA	1:C:269:TYR:HB3	1.93	0.51
1:D:257:LEU:HB2	1:D:265:MET:HE1	1.93	0.51
1:B:292:PHE:CZ	3:B:604:EDO:H21	2.46	0.51
1:D:179:ASP:HB2	1:D:528:ARG:HA	1.93	0.51
1:D:25:ALA:HA	1:D:64:THR:O	2.12	0.50
1:A:95:GLU:H	1:A:95:GLU:CD	2.20	0.50
1:A:93:LYS:HB3	1:A:95:GLU:OE2	2.11	0.50
1:C:70:GLN:HG2	1:C:114:VAL:HG21	1.94	0.50
1:A:503:ARG:HG3	1:A:508:THR:HG22	1.95	0.49
1:D:442:GLU:O	1:D:443:ASN:HB2	2.13	0.48
1:D:133:PRO:HB2	1:D:138:ILE:HD11	1.95	0.48
1:A:257:LEU:HB2	1:A:265:MET:HE1	1.95	0.48
1:C:257:LEU:HB2	1:C:265:MET:HE1	1.95	0.48
1:C:201:MET:HE1	1:C:514:GLN:HE21	1.77	0.47
1:B:268:PHE:HA	1:B:269:TYR:HB3	1.96	0.47
1:D:237:GLN:OE1	5:D:604:GOL:H12	2.13	0.47
1:D:447:LEU:O	1:D:448:ASP:C	2.57	0.47
1:C:193:HIS:CE1	3:C:601:EDO:O1	2.68	0.47
1:A:181:SER:HB3	7:A:912:HOH:O	2.15	0.47
1:C:227:TYR:HA	3:C:601:EDO:H21	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:238:TYR:HA	3:C:601:EDO:H22	1.97	0.47
1:C:455:ASN:OD1	1:C:455:ASN:C	2.58	0.47
1:D:268:PHE:HA	1:D:269:TYR:HB3	1.97	0.47
1:A:25:ALA:O	1:A:267:PRO:HD2	2.15	0.46
1:B:502:VAL:O	1:B:508:THR:HA	2.15	0.46
1:A:25:ALA:HA	1:A:64:THR:O	2.16	0.46
1:C:455:ASN:O	1:C:457:HIS:HD2	1.99	0.46
1:B:432:ALA:HB2	1:B:487:ILE:HD11	1.97	0.46
1:C:193:HIS:HE1	3:C:601:EDO:C1	2.28	0.46
1:A:442:GLU:O	1:A:443:ASN:HB2	2.16	0.45
1:A:455:ASN:O	1:A:457:HIS:HD2	1.99	0.45
1:A:187:ALA:HA	1:A:221:ILE:O	2.16	0.45
1:A:223:THR:O	1:A:224:ALA:HB2	2.16	0.45
1:B:390:TYR:CE2	1:B:396:ILE:CG1	3.00	0.45
1:C:176:ASP:OD1	1:C:457:HIS:CE1	2.66	0.45
1:B:192:ASP:HB2	1:B:203:PHE:CG	2.52	0.45
1:C:227:TYR:HD1	3:C:601:EDO:H21	1.77	0.45
1:A:331:HIS:ND1	1:A:336:ILE:HD12	2.32	0.44
1:D:228:SER:H	3:D:603:EDO:C1	2.30	0.44
1:C:131:ASN:ND2	1:C:169:SER:OG	2.51	0.44
1:C:150:MET:HE2	1:C:150:MET:HA	1.99	0.44
1:D:25:ALA:O	1:D:267:PRO:HD2	2.18	0.44
1:A:176:ASP:OD1	1:A:457:HIS:CE1	2.66	0.44
1:A:181:SER:O	1:A:218:LYS:NZ	2.50	0.44
1:B:223:THR:HA	1:B:266:CYS:HB2	1.98	0.44
1:D:259:ASP:OD2	1:D:327:ASN:HB3	2.18	0.44
1:C:495:ASN:HD21	1:C:514:GLN:NE2	2.15	0.44
1:C:170:ALA:O	1:C:188:TYR:HA	2.18	0.44
1:B:71:TYR:HB3	1:B:75:GLN:HB2	2.00	0.43
1:C:510:LYS:NZ	7:C:713:HOH:O	2.51	0.43
1:C:25:ALA:HA	1:C:64:THR:O	2.17	0.43
1:A:423:GLU:HG2	1:A:424:LYS:HG2	1.99	0.43
1:A:193:HIS:HE1	3:A:601:EDO:O1	2.02	0.43
1:D:237:GLN:OE1	5:D:604:GOL:C1	2.67	0.43
1:D:329:SER:OG	1:D:331:HIS:HE1	2.02	0.43
1:A:76:LEU:HD11	1:A:124:ILE:HD11	2.01	0.43
1:B:171:ASN:HA	1:B:189:ASN:HB2	2.00	0.43
1:A:228:SER:H	3:A:601:EDO:C2	2.31	0.43
1:D:442:GLU:O	1:D:443:ASN:CB	2.67	0.42
1:A:283:HIS:CE1	1:A:291:TRP:CH2	3.08	0.42
1:C:100:ASP:OD1	1:C:101:PRO:HD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:369:MET:HE3	1:B:369:MET:HB2	1.96	0.42
1:D:473:GLN:HG2	1:D:476:LYS:HE2	2.01	0.42
1:A:449:ASP:HB2	7:A:940:HOH:O	2.18	0.42
1:D:416:LEU:HD23	1:D:416:LEU:C	2.44	0.42
1:A:131:ASN:OD1	1:A:169:SER:OG	2.21	0.42
1:A:283:HIS:CE1	1:A:291:TRP:CZ2	3.07	0.42
1:B:448:ASP:O	7:B:701:HOH:O	2.22	0.42
1:C:128:LEU:HA	1:C:167:THR:O	2.20	0.42
1:C:189:ASN:ND2	1:C:223:THR:OG1	2.53	0.42
1:C:227:TYR:HA	3:C:601:EDO:C2	2.50	0.42
1:D:427:ASN:HD22	1:D:519:ARG:HH11	1.64	0.42
1:A:332:THR:HA	1:A:407:SER:O	2.19	0.41
1:D:192:ASP:HB3	1:D:227:TYR:OH	2.20	0.41
1:B:227:TYR:CD1	3:B:604:EDO:H12	2.55	0.41
1:D:151:ASN:HD22	1:D:151:ASN:HA	1.74	0.41
1:A:434:ILE:HD11	1:A:496:ALA:HB1	2.01	0.41
1:A:498:ALA:O	1:A:512:HIS:HA	2.20	0.41
1:B:227:TYR:HD1	3:B:604:EDO:H12	1.86	0.41
1:C:426:LEU:HB2	1:C:519:ARG:HA	2.02	0.41
1:A:426:LEU:HB2	1:A:519:ARG:HA	2.02	0.41
1:C:200:THR:O	1:C:499:GLY:HA3	2.21	0.41
1:C:268:PHE:CE2	2:G:1:BGC:H5	2.56	0.41
1:C:348:LYS:HB3	1:C:390:TYR:HB2	2.03	0.41
1:A:391:ASP:C	1:A:391:ASP:OD1	2.64	0.41
1:B:24:LYS:HE3	7:B:827:HOH:O	2.21	0.41
1:C:259:ASP:OD2	1:C:327:ASN:HB3	2.20	0.41
1:C:338:LEU:HD23	1:C:371:ASP:HB3	2.02	0.41
1:A:369:MET:HE3	1:A:369:MET:HB2	1.84	0.41
1:B:119:LYS:HA	1:B:161:HIS:CE1	2.56	0.40
1:D:225:PHE:CZ	1:D:265:MET:HE3	2.56	0.40
1:D:228:SER:H	3:D:603:EDO:H11	1.87	0.40
1:D:108:GLU:HG2	1:D:156:ILE:CD1	2.52	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	526/537 (98%)	508 (97%)	16 (3%)	2 (0%)	30	37
1	B	524/537 (98%)	507 (97%)	16 (3%)	1 (0%)	44	54
1	C	525/537 (98%)	505 (96%)	17 (3%)	3 (1%)	22	26
1	D	526/537 (98%)	511 (97%)	14 (3%)	1 (0%)	44	54
All	All	2101/2148 (98%)	2031 (97%)	63 (3%)	7 (0%)	37	46

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	142	THR
1	A	519	ARG
1	B	142	THR
1	C	519	ARG
1	A	142	THR
1	C	279	GLU
1	D	142	THR

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	464/473 (98%)	457 (98%)	7 (2%)	60	75
1	B	462/473 (98%)	457 (99%)	5 (1%)	70	83
1	C	463/473 (98%)	458 (99%)	5 (1%)	70	83
1	D	464/473 (98%)	459 (99%)	5 (1%)	70	83
All	All	1853/1892 (98%)	1831 (99%)	22 (1%)	67	80

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	LEU
1	A	185	VAL
1	A	269	TYR
1	A	364	THR
1	A	369	MET
1	A	441	SER
1	A	503	ARG
1	B	4	LYS
1	B	185	VAL
1	B	269	TYR
1	B	369	MET
1	B	476	LYS
1	C	70	GLN
1	C	185	VAL
1	C	269	TYR
1	C	369	MET
1	C	477	LYS
1	D	40	GLU
1	D	269	TYR
1	D	369	MET
1	D	441	SER
1	D	474	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	A	15	HIS
1	A	158	HIS
1	A	193	HIS
1	A	253	ASN
1	A	301	ASN
1	A	399	ASN
1	A	443	ASN
1	A	457	HIS
1	B	189	ASN
1	B	342	ASN
1	B	403	ASN
1	B	465	GLN
1	B	514	GLN
1	C	131	ASN
1	C	134	GLN
1	C	137	HIS

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Mol	Chain	Res	Type
1	C	151	ASN
1	C	189	ASN
1	C	193	HIS
1	C	331	HIS
1	C	403	ASN
1	C	457	HIS
1	C	483	ASN
1	C	514	GLN
1	D	15	HIS
1	D	151	ASN
1	D	189	ASN
1	D	331	HIS
1	D	427	ASN
1	D	443	ASN
1	D	457	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

12 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	BGC	E	1	2	12,12,12	0.48	0	17,17,17	0.58	0
2	BGC	E	2	2	11,11,12	1.54	2 (18%)	15,15,17	0.87	1 (6%)
2	BGC	E	3	2	11,11,12	0.65	0	15,15,17	1.10	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BGC	F	1	2	12,12,12	0.64	0	17,17,17	0.58	0
2	BGC	F	2	2	11,11,12	0.92	1 (9%)	15,15,17	0.82	1 (6%)
2	BGC	F	3	2	11,11,12	0.60	0	15,15,17	1.28	1 (6%)
2	BGC	G	1	2	12,12,12	0.75	0	17,17,17	0.59	0
2	BGC	G	2	2	11,11,12	0.68	0	15,15,17	0.93	1 (6%)
2	BGC	G	3	2	11,11,12	0.80	0	15,15,17	1.33	3 (20%)
2	BGC	H	1	2	12,12,12	0.65	0	17,17,17	0.56	0
2	BGC	H	2	2	11,11,12	0.83	0	15,15,17	0.65	0
2	BGC	H	3	2	11,11,12	0.78	0	15,15,17	1.27	2 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	E	1	2	-	0/2/22/22	0/1/1/1
2	BGC	E	2	2	-	0/2/19/22	0/1/1/1
2	BGC	E	3	2	-	2/2/19/22	0/1/1/1
2	BGC	F	1	2	-	0/2/22/22	0/1/1/1
2	BGC	F	2	2	-	0/2/19/22	0/1/1/1
2	BGC	F	3	2	-	2/2/19/22	0/1/1/1
2	BGC	G	1	2	-	0/2/22/22	0/1/1/1
2	BGC	G	2	2	-	0/2/19/22	0/1/1/1
2	BGC	G	3	2	-	2/2/19/22	0/1/1/1
2	BGC	H	1	2	-	0/2/22/22	0/1/1/1
2	BGC	H	2	2	-	0/2/19/22	0/1/1/1
2	BGC	H	3	2	-	2/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	2	BGC	C2-C3	4.32	1.58	1.52
2	E	2	BGC	O5-C5	2.10	1.47	1.43
2	F	2	BGC	O3-C3	2.03	1.47	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	3	BGC	C1-O5-C5	3.80	117.34	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	BGC	C1-C2-C3	3.12	113.50	109.67
2	G	3	BGC	O5-C5-C6	2.99	111.88	107.20
2	E	2	BGC	O2-C2-C3	2.48	115.11	110.14
2	F	2	BGC	C1-C2-C3	2.34	112.54	109.67
2	H	3	BGC	O5-C5-C6	2.30	110.81	107.20
2	G	3	BGC	O5-C5-C4	-2.22	105.42	110.83
2	H	3	BGC	O5-C5-C4	-2.19	105.49	110.83
2	G	3	BGC	O2-C2-C3	2.15	114.44	110.14

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	3	BGC	O5-C5-C6-O6
2	H	3	BGC	O5-C5-C6-O6
2	H	3	BGC	C4-C5-C6-O6
2	G	3	BGC	C4-C5-C6-O6
2	F	3	BGC	C4-C5-C6-O6
2	F	3	BGC	O5-C5-C6-O6
2	E	3	BGC	O5-C5-C6-O6
2	E	3	BGC	C4-C5-C6-O6

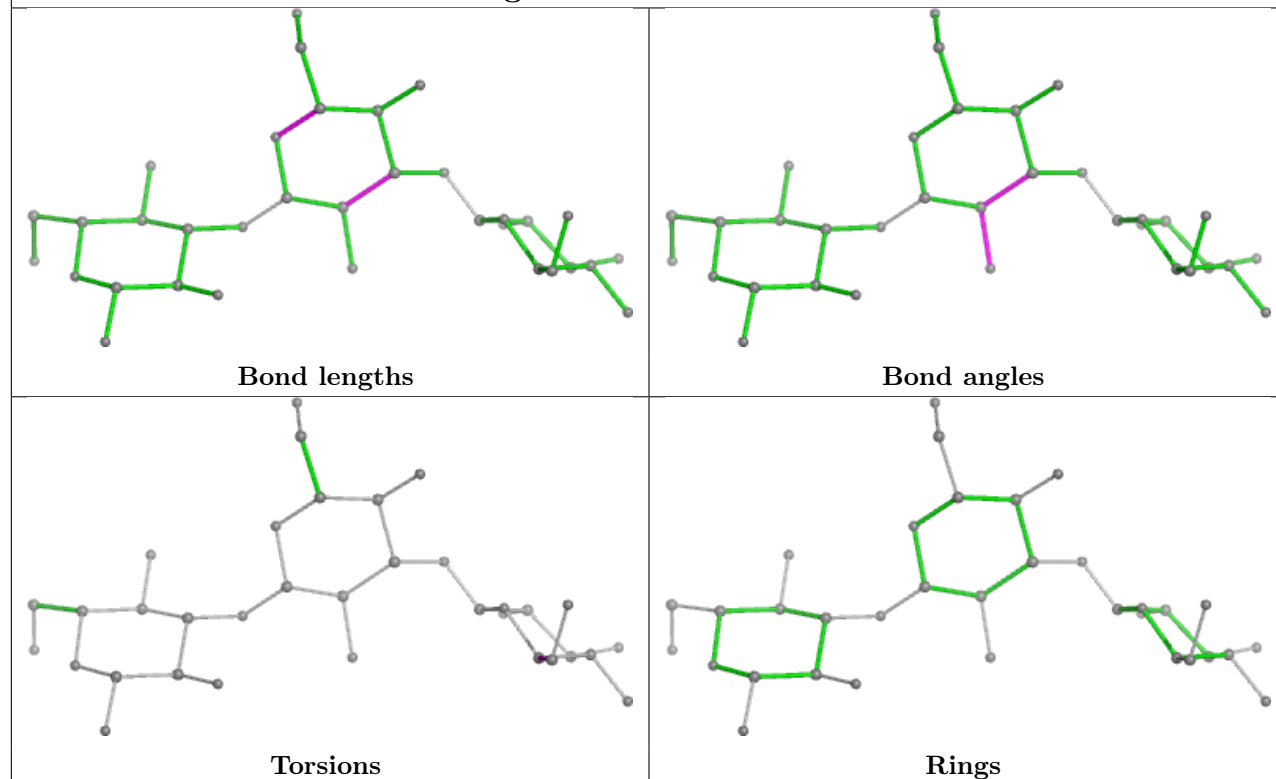
There are no ring outliers.

1 monomer is involved in 1 short contact:

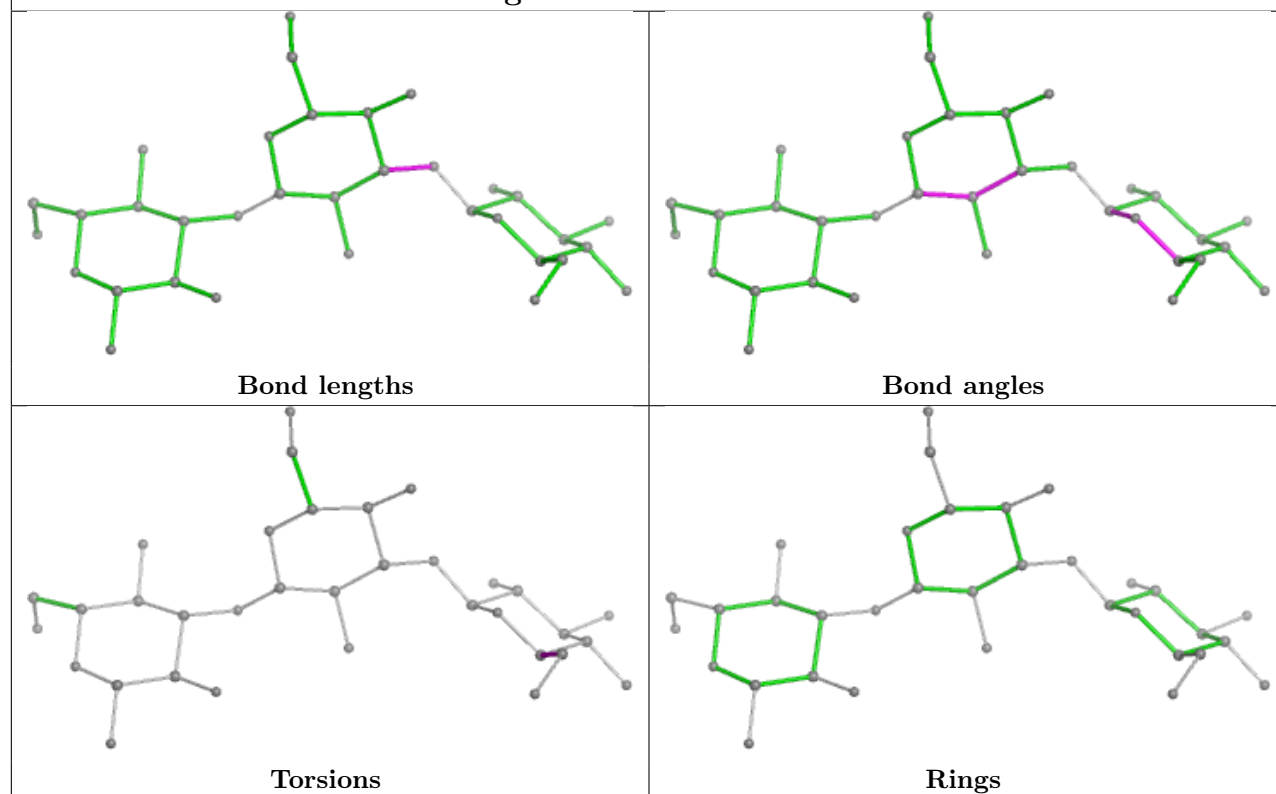
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	1	BGC	1	0

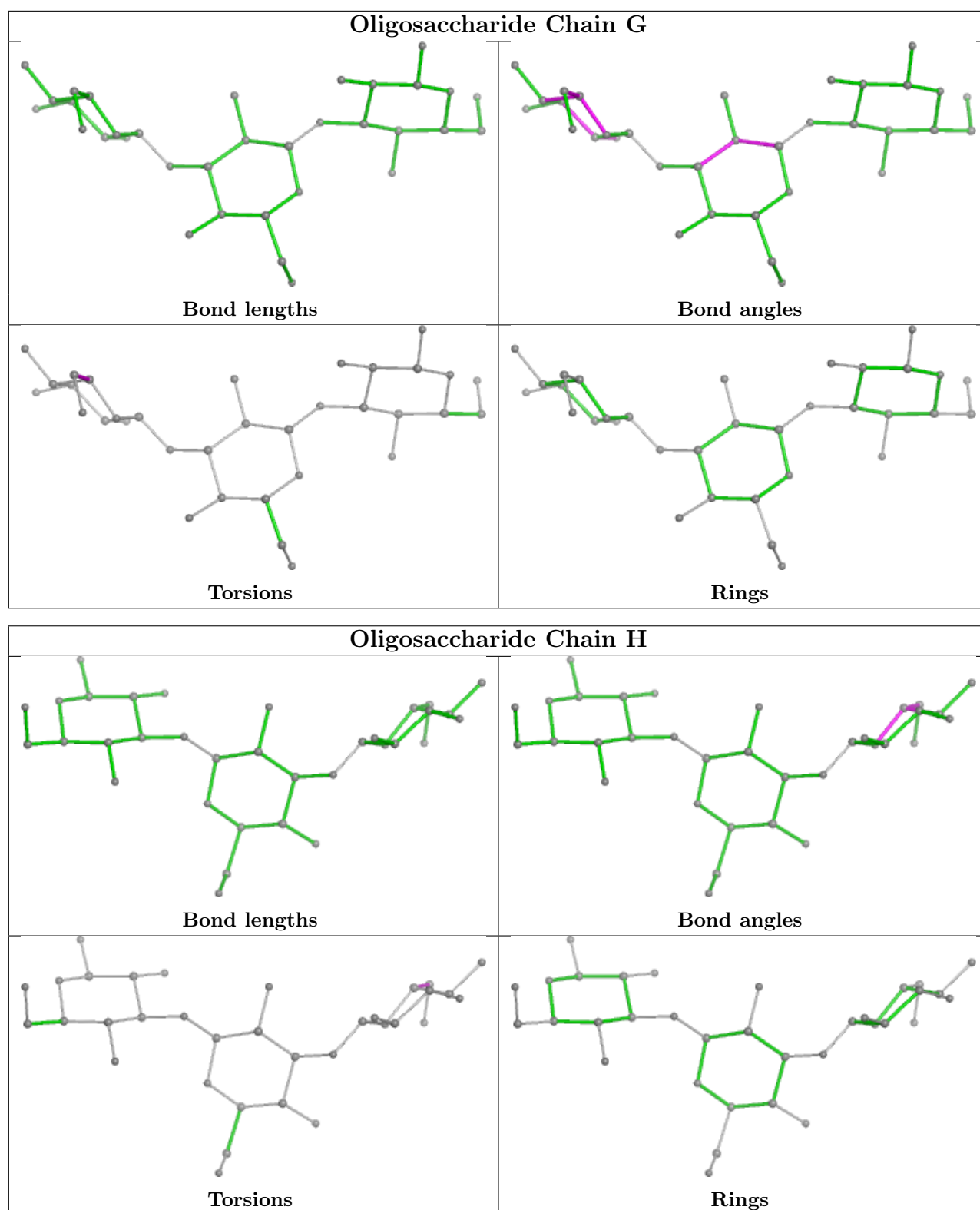
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain E



Oligosaccharide Chain F





5.6 Ligand geometry [i](#)

19 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	BGC	B	602	-	12,12,12	1.24	0	17,17,17	0.81	0
5	GOL	D	605	-	5,5,5	0.20	0	5,5,5	0.33	0
5	GOL	D	606	-	5,5,5	0.09	0	5,5,5	0.32	0
3	EDO	C	601	-	3,3,3	0.22	0	2,2,2	0.41	0
5	GOL	D	602	-	5,5,5	0.10	0	5,5,5	0.24	0
5	GOL	D	607	-	5,5,5	0.09	0	5,5,5	0.23	0
3	EDO	B	601	-	3,3,3	0.76	0	2,2,2	0.45	0
3	EDO	D	603	-	3,3,3	0.22	0	2,2,2	0.14	0
4	BGC	A	602	-	12,12,12	1.15	1 (8%)	17,17,17	1.10	1 (5%)
5	GOL	B	605	-	5,5,5	0.28	0	5,5,5	0.66	0
3	EDO	B	604	-	3,3,3	0.50	0	2,2,2	0.28	0
6	PGE	C	602	-	9,9,9	0.48	0	8,8,8	0.44	0
5	GOL	A	603	-	5,5,5	0.18	0	5,5,5	0.40	0
3	EDO	A	601	-	3,3,3	0.17	0	2,2,2	0.35	0
5	GOL	B	606	-	5,5,5	0.26	0	5,5,5	0.52	0
5	GOL	B	603	-	5,5,5	0.32	0	5,5,5	0.59	0
5	GOL	C	603	-	5,5,5	0.12	0	5,5,5	0.28	0
5	GOL	D	604	-	5,5,5	0.30	0	5,5,5	0.62	0
4	BGC	D	601	-	12,12,12	0.18	0	17,17,17	0.35	0

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	BGC	B	602	-	-	1/2/22/22	0/1/1/1
5	GOL	D	605	-	-	4/4/4/4	-
5	GOL	D	606	-	-	3/4/4/4	-
3	EDO	C	601	-	-	0/1/1/1	-
5	GOL	D	602	-	-	4/4/4/4	-
5	GOL	D	607	-	-	0/4/4/4	-
3	EDO	B	601	-	-	0/1/1/1	-
3	EDO	D	603	-	-	0/1/1/1	-
4	BGC	A	602	-	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	B	605	-	-	2/4/4/4	-
3	EDO	B	604	-	-	1/1/1/1	-
6	PGE	C	602	-	-	4/7/7/7	-
5	GOL	A	603	-	-	2/4/4/4	-
3	EDO	A	601	-	-	0/1/1/1	-
5	GOL	B	606	-	-	0/4/4/4	-
5	GOL	B	603	-	-	2/4/4/4	-
5	GOL	C	603	-	-	2/4/4/4	-
5	GOL	D	604	-	-	4/4/4/4	-
4	BGC	D	601	-	-	0/2/22/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	602	BGC	O5-C5	2.66	1.50	1.44

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	602	BGC	O2-C2-C3	2.53	116.21	110.35

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	603	GOL	O1-C1-C2-C3
5	D	602	GOL	O1-C1-C2-C3
5	D	604	GOL	O1-C1-C2-C3
5	D	604	GOL	C1-C2-C3-O3
5	D	605	GOL	O1-C1-C2-O2
5	D	605	GOL	O1-C1-C2-C3
5	D	605	GOL	C1-C2-C3-O3
5	D	606	GOL	C1-C2-C3-O3
5	C	603	GOL	O1-C1-C2-O2
6	C	602	PGE	O1-C1-C2-O2
5	A	603	GOL	O1-C1-C2-C3
5	B	603	GOL	O1-C1-C2-C3
5	B	605	GOL	O1-C1-C2-C3
5	D	602	GOL	C1-C2-C3-O3
5	D	606	GOL	O1-C1-C2-C3

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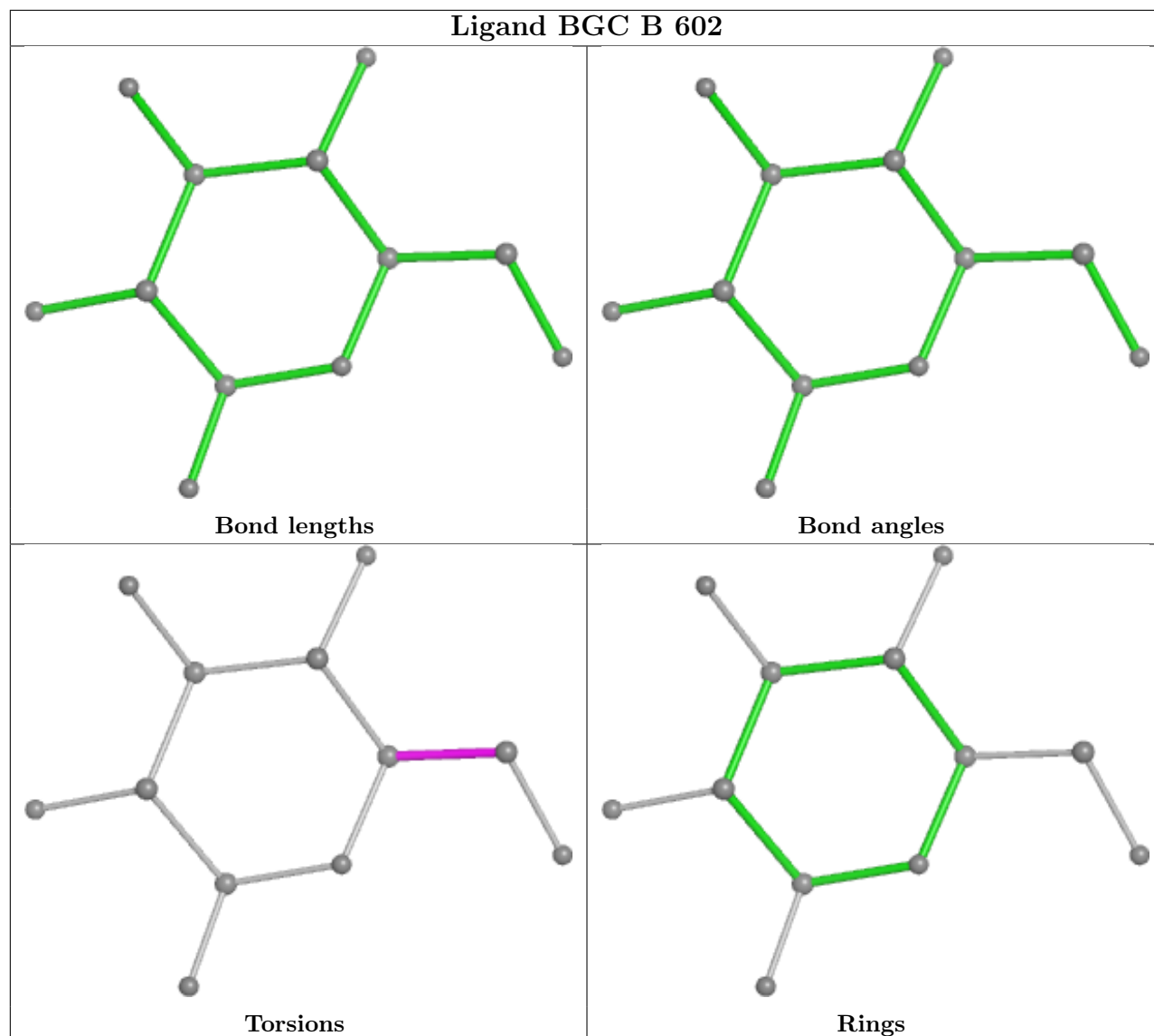
Mol	Chain	Res	Type	Atoms
5	B	603	GOL	O1-C1-C2-O2
5	B	605	GOL	O1-C1-C2-O2
5	D	602	GOL	O1-C1-C2-O2
5	D	604	GOL	O1-C1-C2-O2
5	D	604	GOL	O2-C2-C3-O3
6	C	602	PGE	O3-C5-C6-O4
5	D	606	GOL	O2-C2-C3-O3
5	D	602	GOL	O2-C2-C3-O3
6	C	602	PGE	C3-C4-O3-C5
5	A	603	GOL	O1-C1-C2-O2
6	C	602	PGE	O2-C3-C4-O3
5	D	605	GOL	O2-C2-C3-O3
3	B	604	EDO	O1-C1-C2-O2
4	B	602	BGC	O5-C5-C6-O6

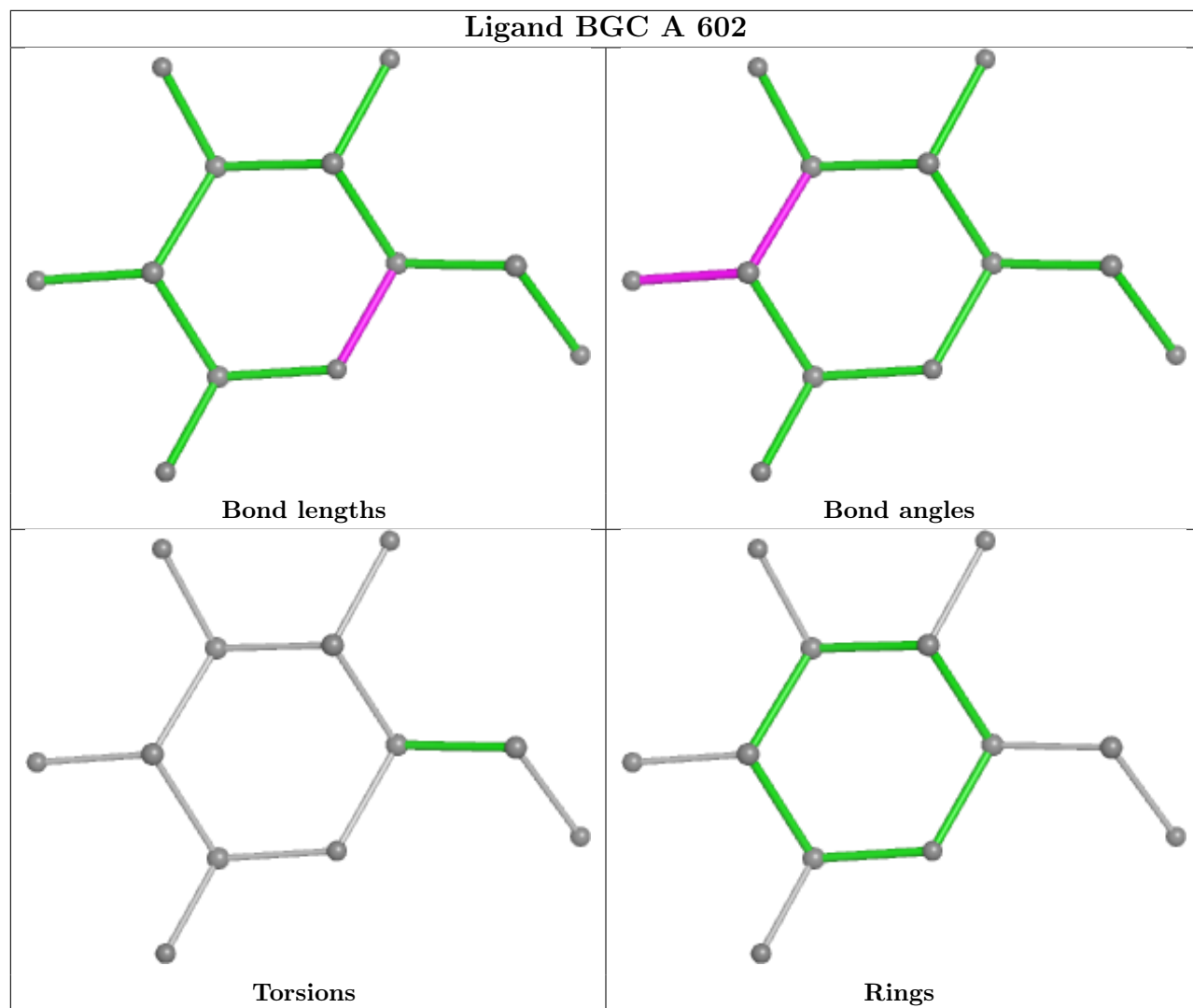
There are no ring outliers.

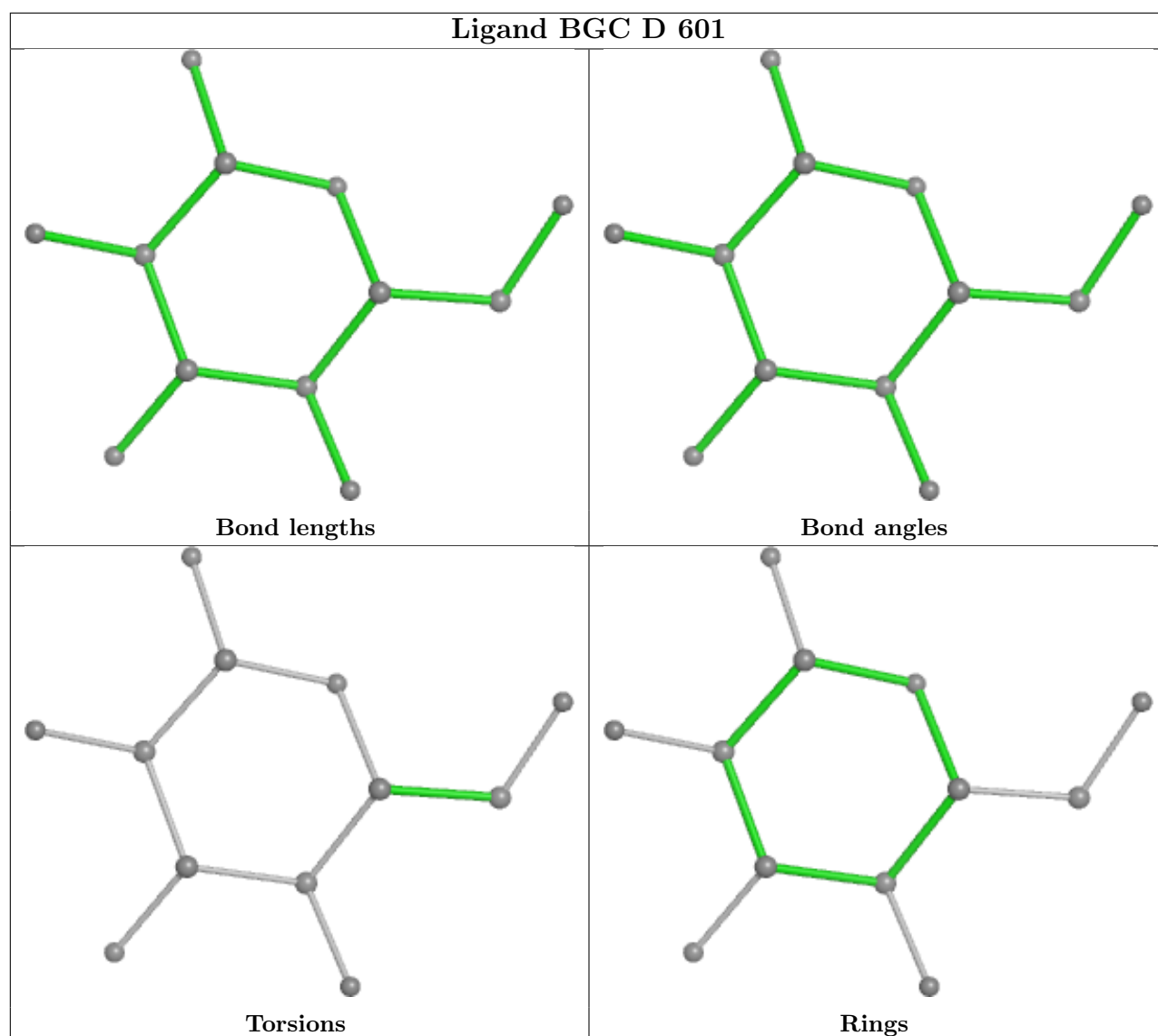
6 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	601	EDO	9	0
3	D	603	EDO	4	0
5	B	605	GOL	1	0
3	B	604	EDO	5	0
3	A	601	EDO	7	0
5	D	604	GOL	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 [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	528/537 (98%)	-0.49	2 (0%) 89 89	24, 32, 50, 72	0
1	B	526/537 (97%)	-0.46	2 (0%) 89 89	24, 34, 53, 76	0
1	C	527/537 (98%)	-0.24	3 (0%) 85 86	25, 39, 65, 86	0
1	D	527/537 (98%)	-0.46	6 (1%) 77 78	19, 33, 52, 80	1 (0%)
All	All	2108/2148 (98%)	-0.41	13 (0%) 85 86	19, 34, 57, 86	1 (0%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	530	LEU	3.7
1	C	105	LYS	3.1
1	D	43	ARG	2.8
1	C	109	ILE	2.6
1	B	235	ASN	2.5
1	D	41	ASP	2.2
1	D	344	LYS	2.2
1	D	40	GLU	2.2
1	A	520	GLY	2.2
1	C	474	LEU	2.2
1	D	529	LYS	2.1
1	D	3	SER	2.0
1	B	4	LYS	2.0

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

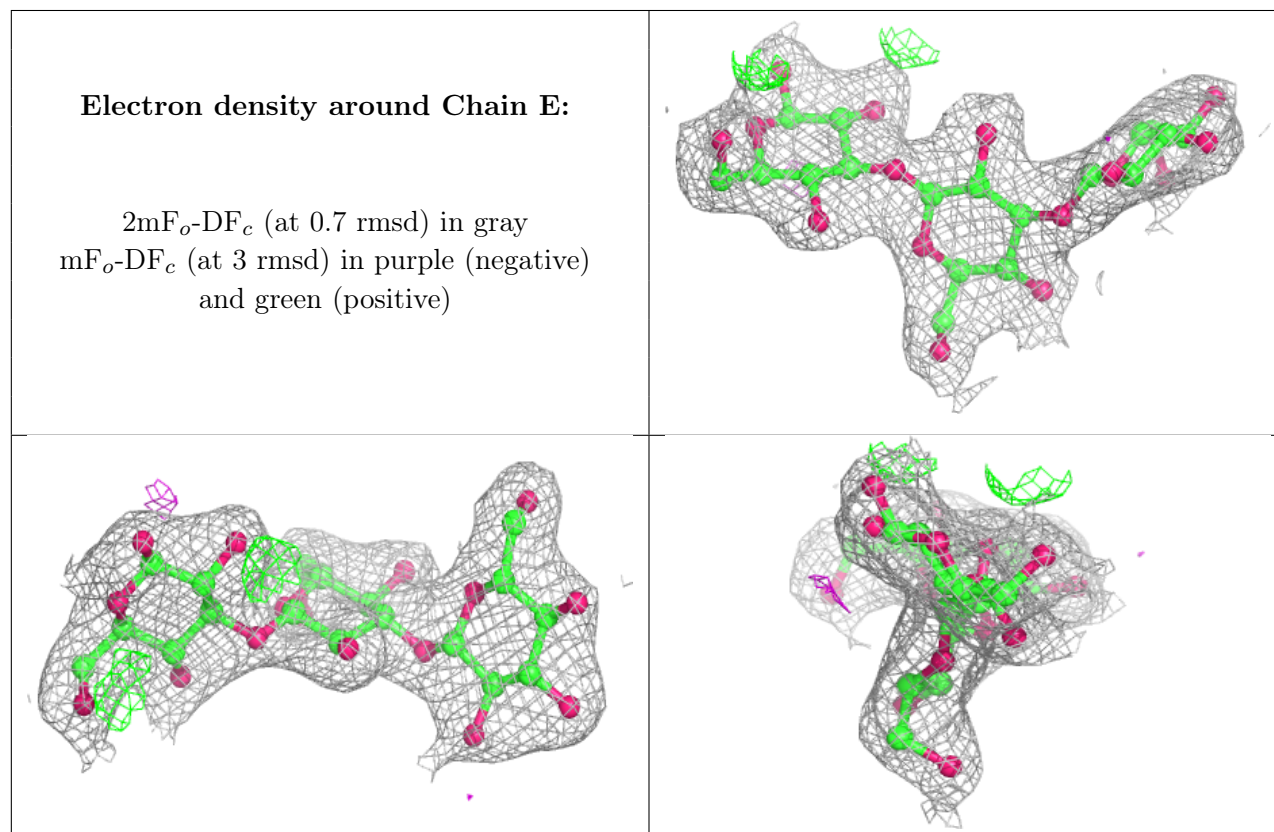
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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.

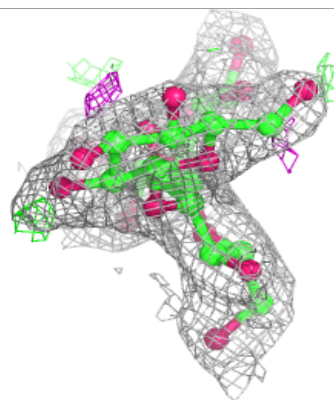
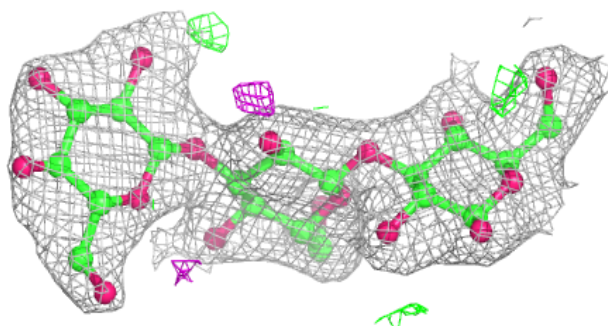
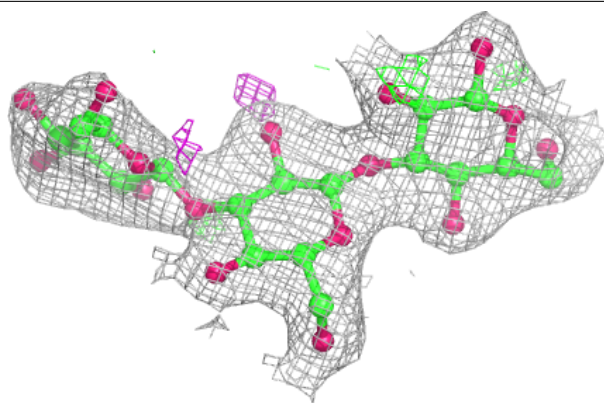
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BGC	G	3	11/12	0.91	0.09	41,48,59,64	0
2	BGC	H	3	11/12	0.91	0.11	40,46,57,60	0
2	BGC	F	3	11/12	0.93	0.08	35,40,50,52	0
2	BGC	G	2	11/12	0.94	0.07	31,34,37,39	0
2	BGC	F	2	11/12	0.94	0.07	24,29,33,33	0
2	BGC	E	3	11/12	0.94	0.08	33,37,42,46	0
2	BGC	E	2	11/12	0.95	0.06	26,28,30,33	0
2	BGC	H	1	12/12	0.96	0.06	22,24,29,29	0
2	BGC	H	2	11/12	0.96	0.06	27,29,31,32	0
2	BGC	G	1	12/12	0.96	0.07	24,26,32,32	0
2	BGC	E	1	12/12	0.97	0.06	22,26,28,30	0
2	BGC	F	1	12/12	0.97	0.06	25,28,30,30	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

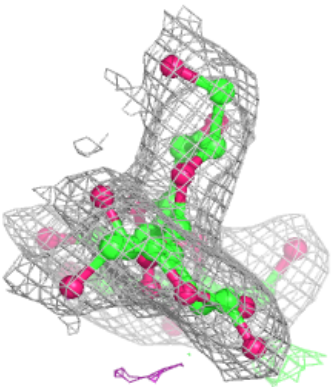
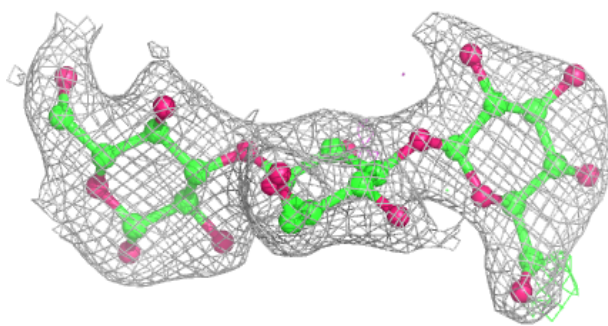
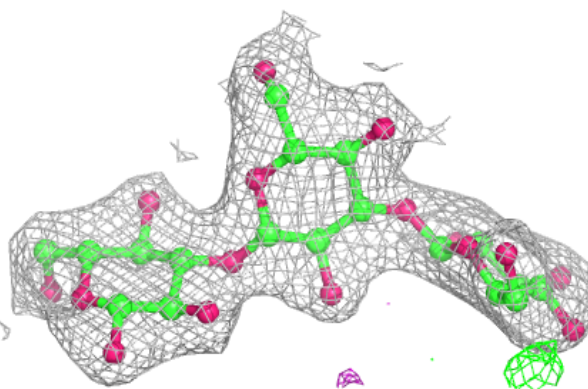


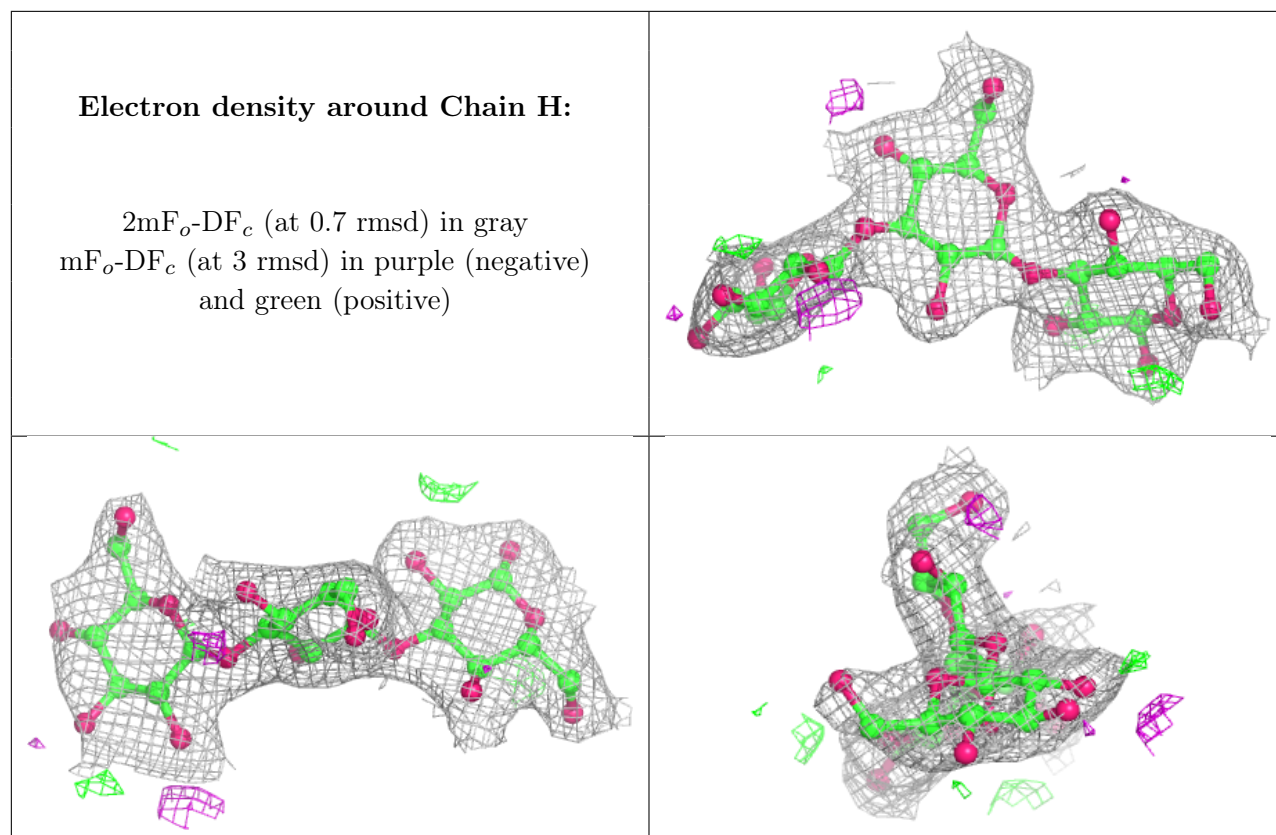
Electron density around Chain F:

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

**Electron density around Chain G:**

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





6.4 Ligands ⓘ

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(\AA^2)	Q<0.9
4	BGC	D	601	12/12	0.75	0.19	41,54,62,63	0
4	BGC	A	602	12/12	0.81	0.14	32,42,47,50	12
4	BGC	B	602	12/12	0.83	0.15	34,45,49,50	12
5	GOL	A	603	6/6	0.83	0.17	46,55,61,61	0
5	GOL	D	602	6/6	0.84	0.16	39,52,59,61	0
5	GOL	D	605	6/6	0.85	0.16	42,53,58,59	0
6	PGE	C	602	10/10	0.85	0.16	52,57,63,66	0
5	GOL	B	605	6/6	0.86	0.16	40,45,47,48	0
5	GOL	D	607	6/6	0.86	0.16	61,63,65,70	0
5	GOL	D	604	6/6	0.86	0.15	44,49,54,56	0
3	EDO	B	601	4/4	0.87	0.14	34,42,44,48	0
5	GOL	B	606	6/6	0.88	0.14	51,60,62,64	0
5	GOL	D	606	6/6	0.88	0.14	57,63,65,66	0
5	GOL	B	603	6/6	0.91	0.12	34,46,47,48	0

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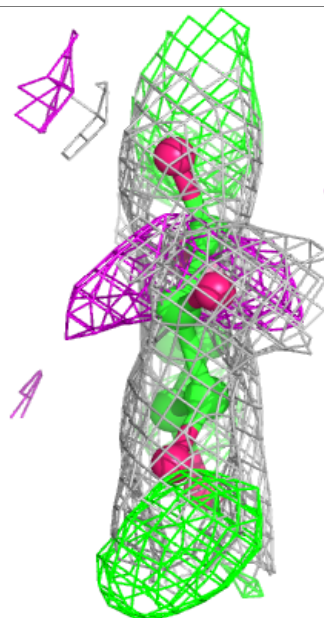
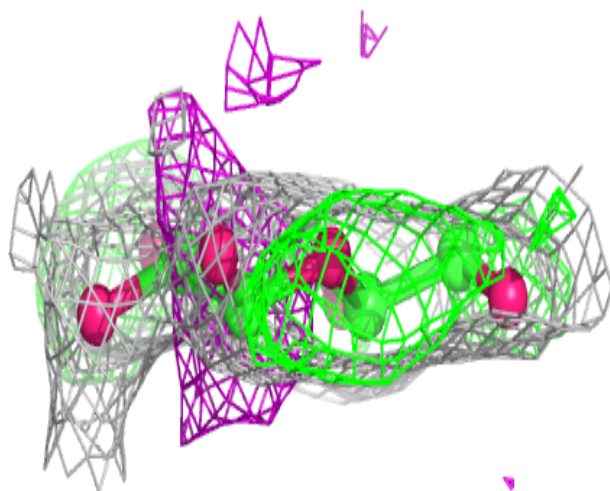
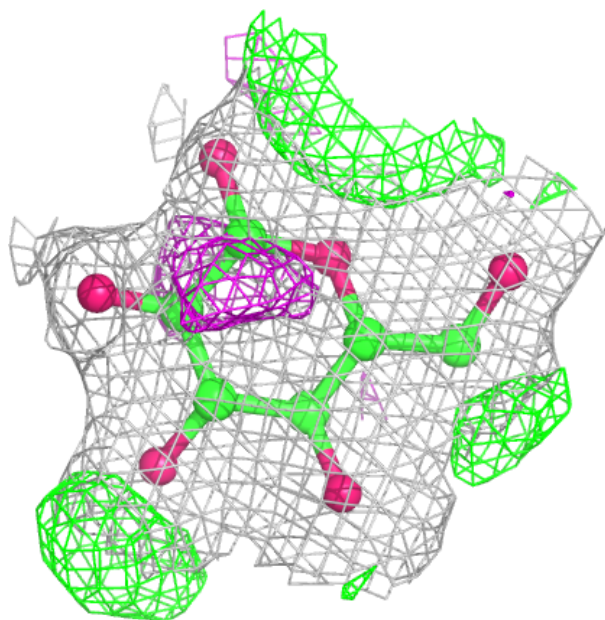
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	EDO	D	603	4/4	0.94	0.18	32,37,37,37	0
3	EDO	A	601	4/4	0.94	0.23	34,34,36,37	0
3	EDO	C	601	4/4	0.95	0.18	33,39,40,43	0
5	GOL	C	603	6/6	0.95	0.09	43,54,58,59	0
3	EDO	B	604	4/4	0.96	0.17	33,37,42,42	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.

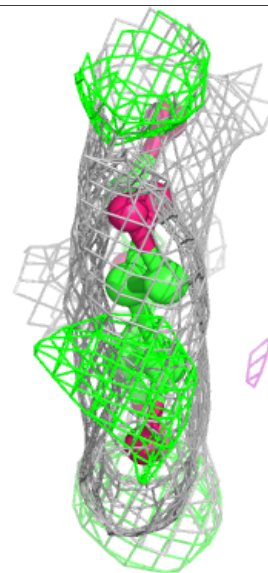
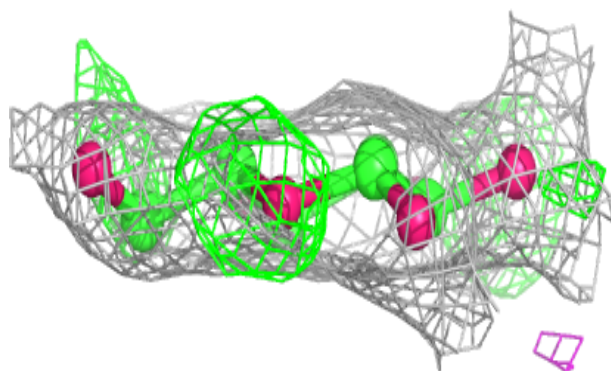
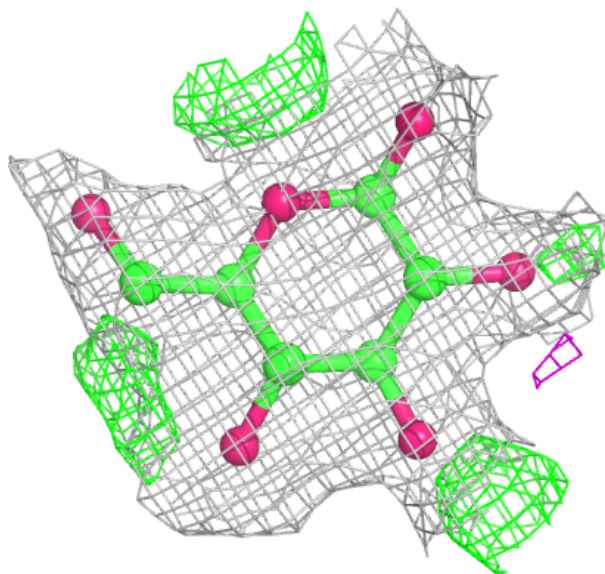
Electron density around BGC D 601:

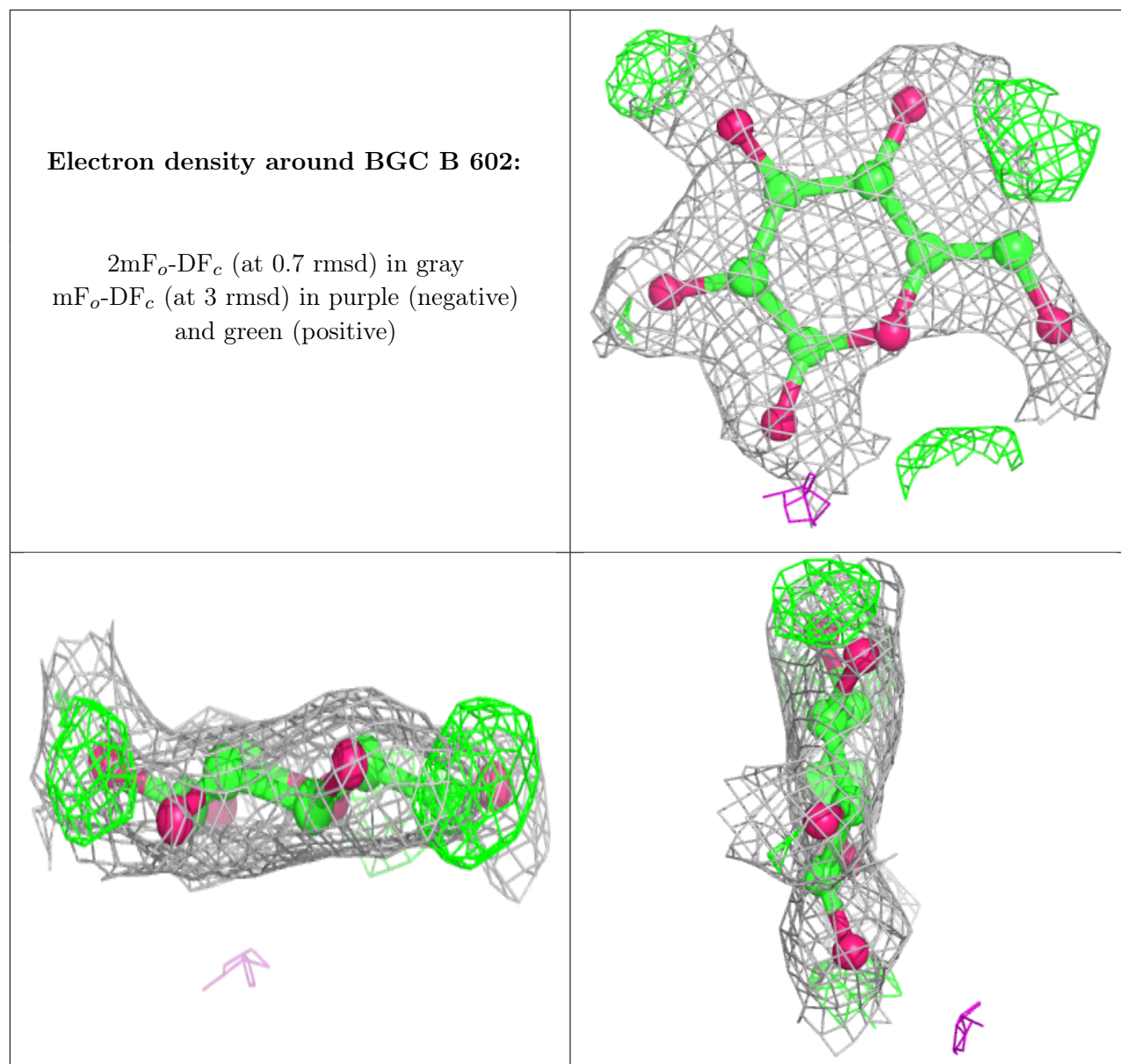
2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around BGC A 602:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.