



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 28, 2025 – 10:22 AM EDT

PDB ID : 9C9O / pdb_00009c9o
Title : M. tuberculosis PKS13 acyltransferase (AT) domain in complex with SuFEx inhibitor CMX410 - reaction product
Authors : Krieger, I.V.; Tang, S.; Sacchetini, J.C.; TB Structural Genomics Consortium (TBSGC)
Deposited on : 2024-06-14
Resolution : 2.02 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 : 2022.3.0, CSD as543be (2022)
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 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

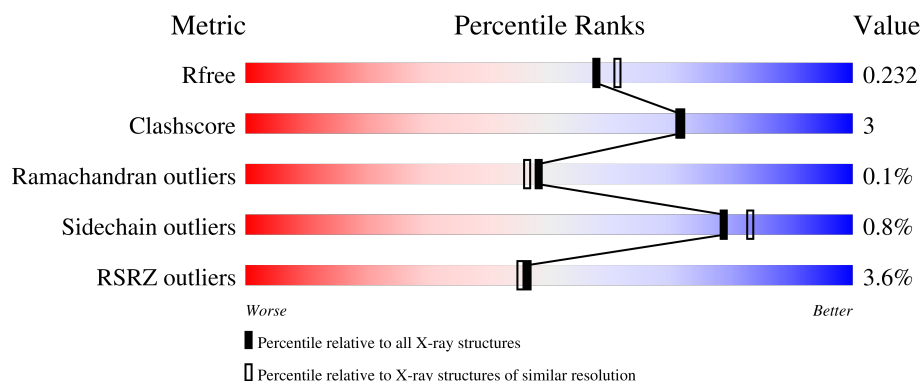
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


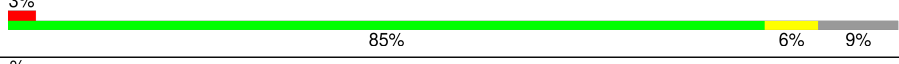
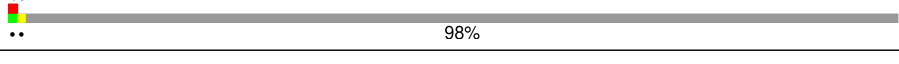

The reported resolution of this entry is 2.02 Å.

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	12358 (2.04-2.00)
Clashscore	180529	13897 (2.04-2.00)
Ramachandran outliers	177936	13770 (2.04-2.00)
Sidechain outliers	177891	13769 (2.04-2.00)
RSRZ outliers	164620	12358 (2.04-2.00)

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	512	
2	B	512	
2	C	512	
3	D	2	

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Mol	Chain	Length	Quality of chain
3	E	2	 50%50%
3	F	2	 50%50%
3	G	2	 50%50%
3	H	2	 100%
3	I	2	 50%50%
3	J	2	 100%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7875 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 Polyketide synthase Pks13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	468	3559	2252	617	677	13	0	3	0

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	552	MET	-	initiating methionine	UNP I6X8D2
A	553	HIS	-	expression tag	UNP I6X8D2
A	554	HIS	-	expression tag	UNP I6X8D2
A	555	HIS	-	expression tag	UNP I6X8D2
A	556	HIS	-	expression tag	UNP I6X8D2
A	557	HIS	-	expression tag	UNP I6X8D2
A	558	HIS	-	expression tag	UNP I6X8D2
A	559	SER	-	expression tag	UNP I6X8D2
A	560	SER	-	expression tag	UNP I6X8D2
A	561	GLY	-	expression tag	UNP I6X8D2
A	562	VAL	-	expression tag	UNP I6X8D2
A	563	ASP	-	expression tag	UNP I6X8D2
A	564	LEU	-	expression tag	UNP I6X8D2
A	565	GLY	-	expression tag	UNP I6X8D2
A	566	THR	-	expression tag	UNP I6X8D2
A	567	GLU	-	expression tag	UNP I6X8D2
A	568	ASN	-	expression tag	UNP I6X8D2
A	569	LEU	-	expression tag	UNP I6X8D2
A	570	TYR	-	expression tag	UNP I6X8D2
A	571	PHE	-	expression tag	UNP I6X8D2
A	572	GLN	-	expression tag	UNP I6X8D2
A	573	SER	-	expression tag	UNP I6X8D2
A	574	ASN	-	expression tag	UNP I6X8D2
A	575	ALA	-	expression tag	UNP I6X8D2
A	801	A1ATO	SER	conflict	UNP I6X8D2

- Molecule 2 is a protein called Polyketide synthase Pks13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	468	Total	C	N	O	S	0	1	0
			3554	2245	619	678	12			
2	C	10	Total	C	N	O		0	0	0
			80	53	11	16				

There are 48 discrepancies between the modelled and reference sequences:

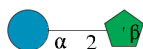
Chain	Residue	Modelled	Actual	Comment	Reference
B	552	MET	-	initiating methionine	UNP I6X8D2
B	553	HIS	-	expression tag	UNP I6X8D2
B	554	HIS	-	expression tag	UNP I6X8D2
B	555	HIS	-	expression tag	UNP I6X8D2
B	556	HIS	-	expression tag	UNP I6X8D2
B	557	HIS	-	expression tag	UNP I6X8D2
B	558	HIS	-	expression tag	UNP I6X8D2
B	559	SER	-	expression tag	UNP I6X8D2
B	560	SER	-	expression tag	UNP I6X8D2
B	561	GLY	-	expression tag	UNP I6X8D2
B	562	VAL	-	expression tag	UNP I6X8D2
B	563	ASP	-	expression tag	UNP I6X8D2
B	564	LEU	-	expression tag	UNP I6X8D2
B	565	GLY	-	expression tag	UNP I6X8D2
B	566	THR	-	expression tag	UNP I6X8D2
B	567	GLU	-	expression tag	UNP I6X8D2
B	568	ASN	-	expression tag	UNP I6X8D2
B	569	LEU	-	expression tag	UNP I6X8D2
B	570	TYR	-	expression tag	UNP I6X8D2
B	571	PHE	-	expression tag	UNP I6X8D2
B	572	GLN	-	expression tag	UNP I6X8D2
B	573	SER	-	expression tag	UNP I6X8D2
B	574	ASN	-	expression tag	UNP I6X8D2
B	575	ALA	-	expression tag	UNP I6X8D2
C	552	MET	-	initiating methionine	UNP I6X8D2
C	553	HIS	-	expression tag	UNP I6X8D2
C	554	HIS	-	expression tag	UNP I6X8D2
C	555	HIS	-	expression tag	UNP I6X8D2
C	556	HIS	-	expression tag	UNP I6X8D2
C	557	HIS	-	expression tag	UNP I6X8D2
C	558	HIS	-	expression tag	UNP I6X8D2
C	559	SER	-	expression tag	UNP I6X8D2
C	560	SER	-	expression tag	UNP I6X8D2
C	561	GLY	-	expression tag	UNP I6X8D2
C	562	VAL	-	expression tag	UNP I6X8D2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	563	ASP	-	expression tag	UNP I6X8D2
C	564	LEU	-	expression tag	UNP I6X8D2
C	565	GLY	-	expression tag	UNP I6X8D2
C	566	THR	-	expression tag	UNP I6X8D2
C	567	GLU	-	expression tag	UNP I6X8D2
C	568	ASN	-	expression tag	UNP I6X8D2
C	569	LEU	-	expression tag	UNP I6X8D2
C	570	TYR	-	expression tag	UNP I6X8D2
C	571	PHE	-	expression tag	UNP I6X8D2
C	572	GLN	-	expression tag	UNP I6X8D2
C	573	SER	-	expression tag	UNP I6X8D2
C	574	ASN	-	expression tag	UNP I6X8D2
C	575	ALA	-	expression tag	UNP I6X8D2

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



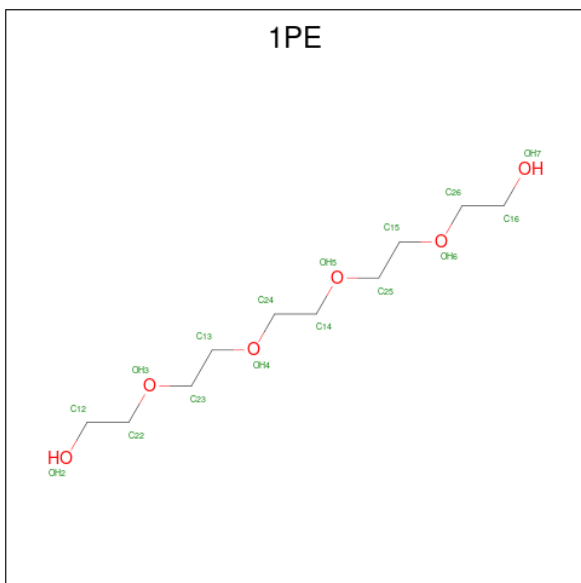
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	D	2	Total	C	O	0	0	0
			23	12	11			
3	E	2	Total	C	O	0	0	0
			23	12	11			
3	F	2	Total	C	O	0	0	0
			23	12	11			
3	G	2	Total	C	O	0	0	0
			23	12	11			
3	H	2	Total	C	O	0	0	0
			23	12	11			
3	I	2	Total	C	O	0	0	0
			23	12	11			
3	J	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



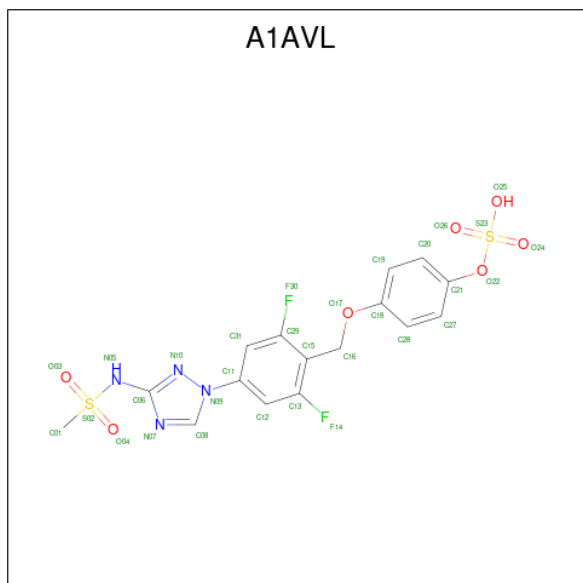
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			13	8	5		
5	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 6 is 4-(2,6-difluoro-4-[3-(methanesulfonamido)-1H-1,2,4-triazol-1-yl]phenyl)methoxy)phenyl hydrogen sulfate (CCD ID: A1AVL) (formula: C₁₆H₁₄F₂N₄O₇S₂) (labeled as "Ligand of Interest" by depositor).



- Molecule 1: Polyketide synthase Pks13



[illegible]

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



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



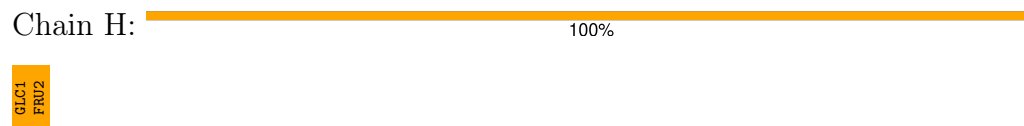
- Molecule 3: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose



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



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



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

Chain I:  50% 50%

GLC1
FRU2

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

Chain J:  100%

GLC1
FRU2

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	106.67Å 106.67Å 258.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	52.24 – 2.02 52.24 – 2.02	Depositor EDS
% Data completeness (in resolution range)	80.5 (52.24-2.02) 80.6 (52.24-2.02)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.01Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.194 , 0.231 0.196 , 0.232	Depositor DCC
R_{free} test set	4938 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	31.3	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7875	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.69% 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: A1AVL, FRU, SO4, 1PE, GLC, A1ATO

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	0.30	0/3635	0.46	0/4933
2	B	0.33	0/3631	0.50	0/4932
2	C	0.35	0/81	0.39	0/108
All	All	0.31	0/7347	0.48	0/9973

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	3559	0	3528	19	0
2	B	3554	0	3516	21	0
2	C	80	0	69	4	0
3	D	23	0	20	0	0
3	E	23	0	20	1	0
3	F	23	0	20	1	0
3	G	23	0	20	0	0
3	H	23	0	20	2	0
3	I	23	0	20	2	0
3	J	23	0	20	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	10	0	0	0	0
4	B	10	0	0	0	0
5	A	26	0	34	1	0
6	A	31	0	0	0	0
7	A	212	0	0	2	0
7	B	230	0	0	7	0
7	C	2	0	0	0	0
All	All	7875	0	7287	47	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 3.

All (47) 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)
3:I:2:FRU:C2	3:I:2:FRU:O5	1.65	1.30
3:E:2:FRU:C2	3:E:2:FRU:O5	1.64	1.26
3:F:2:FRU:O5	3:F:2:FRU:C2	1.64	1.19
3:H:2:FRU:C2	3:H:2:FRU:O5	1.65	1.17
2:B:1019:GLN:HE22	2:C:579:GLU:HA	1.58	0.68
2:B:889:LEU:HD22	2:B:899:ALA:HB1	1.76	0.66
1:A:900:ARG:NH2	7:A:1201:HOH:O	2.28	0.66
2:B:889:LEU:HD11	2:B:901:LYS:HB2	1.77	0.65
2:B:974:ARG:NH1	7:B:1202:HOH:O	2.28	0.62
1:A:632:PHE:HB3	1:A:636:ARG:HH22	1.65	0.61
1:A:762:ASP:HB3	1:A:765:GLN:HG3	1.85	0.58
3:H:1:GLC:H5	3:H:2:FRU:H61	1.86	0.58
1:A:974:ARG:NH2	1:A:1003:ASP:OD2	2.36	0.56
2:B:826:ARG:NH2	7:B:1204:HOH:O	2.35	0.56
1:A:722:GLN:HE22	5:A:1104:1PE:H232	1.72	0.54
2:B:1019:GLN:NE2	2:C:579:GLU:HA	2.21	0.54
2:B:957:GLU:HG2	7:B:1384:HOH:O	2.08	0.54
2:B:1061:ARG:HD3	7:B:1345:HOH:O	2.06	0.54
1:A:844:LEU:HD13	1:A:901:LYS:HD3	1.89	0.53
2:B:921:THR:OG1	2:B:960:LYS:HD2	2.09	0.53
2:B:1047:ARG:HD2	7:B:1304:HOH:O	2.09	0.52
1:A:677:VAL:HG22	1:A:701:VAL:HG22	1.92	0.52
2:B:823:ILE:O	2:B:827:SER:OG	2.20	0.51
1:A:1029:MET:O	1:A:1039:LEU:HD22	2.10	0.51
1:A:632:PHE:H	1:A:636:ARG:NH1	2.09	0.50
1:A:793:LYS:HD2	7:A:1258:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1019:GLN:HE22	2:C:579:GLU:CA	2.26	0.49
2:B:946:ILE:HD12	2:B:953:ILE:HD13	1.94	0.48
1:A:849:GLU:HB3	1:A:897:LYS:HB3	1.95	0.48
2:B:829:LEU:HD21	2:B:919:GLU:HG2	1.96	0.47
2:B:869:VAL:HB	2:B:877:VAL:HB	1.97	0.46
7:B:1412:HOH:O	3:I:2:FRU:H62	2.17	0.45
1:A:830:MET:HE3	1:A:908:SER:HB3	1.98	0.44
1:A:947:LYS:HB2	1:A:947:LYS:NZ	2.33	0.44
1:A:1042:ARG:HE	1:A:1042:ARG:HB2	1.47	0.44
1:A:724:ARG:HD3	1:A:763:ASP:O	2.17	0.44
2:B:932:PRO:HG2	7:B:1206:HOH:O	2.17	0.43
2:B:644:LEU:HD21	2:B:665:LEU:HD11	2.00	0.43
2:B:944:ARG:HB3	2:B:944:ARG:NH1	2.34	0.43
1:A:1029:MET:HG2	1:A:1039:LEU:HD11	2.01	0.42
1:A:676:VAL:HG11	1:A:1030:ALA:HB1	2.01	0.42
2:B:942:GLU:HA	2:B:968:TYR:CD1	2.54	0.42
2:C:578:ASP:OD2	2:C:582:ASN:HB2	2.20	0.41
1:A:681:ASP:OD1	1:A:684:GLU:HG3	2.20	0.41
1:A:1008:ASP:N	1:A:1008:ASP:OD1	2.55	0.40
2:B:769:ILE:HD11	2:B:904:THR:HA	2.04	0.40
2:B:597:LEU:HD21	2:B:817:ARG:NH2	2.36	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	466/512 (91%)	459 (98%)	7 (2%)	0	100	100
2	B	467/512 (91%)	457 (98%)	9 (2%)	1 (0%)	44	41
2	C	8/512 (2%)	8 (100%)	0	0	100	100
All	All	941/1536 (61%)	924 (98%)	16 (2%)	1 (0%)	48	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	955	ASP

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	366/405 (90%)	362 (99%)	4 (1%)	70	75
2	B	366/406 (90%)	363 (99%)	3 (1%)	79	83
2	C	8/406 (2%)	8 (100%)	0	100	100
All	All	740/1217 (61%)	733 (99%)	7 (1%)	79	80

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	756	VAL
1	A	826[A]	ARG
1	A	826[B]	ARG
1	A	889	LEU
2	B	802	LEU
2	B	857	GLU
2	B	1039	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	680	HIS
1	A	722	GLN
1	A	780	GLN
1	A	925	GLN
1	A	1031	GLN
2	B	780	GLN
2	B	1010	GLN
2	B	1019	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

14 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$
3	GLC	D	1	3	11,11,12	0.59	0	15,15,17	0.66	0
3	FRU	D	2	3	11,12,12	4.46	6 (54%)	10,18,18	0.94	1 (10%)
3	GLC	E	1	3	11,11,12	0.57	0	15,15,17	1.23	2 (13%)
3	FRU	E	2	3	11,12,12	4.53	5 (45%)	10,18,18	1.16	0
3	GLC	F	1	3	11,11,12	0.60	0	15,15,17	0.91	0
3	FRU	F	2	3	11,12,12	4.71	5 (45%)	10,18,18	1.15	1 (10%)
3	GLC	G	1	3	11,11,12	0.49	0	15,15,17	0.85	0
3	FRU	G	2	3	11,12,12	4.47	7 (63%)	10,18,18	1.22	1 (10%)
3	GLC	H	1	3	11,11,12	0.52	0	15,15,17	1.30	3 (20%)
3	FRU	H	2	3	11,12,12	4.88	5 (45%)	10,18,18	0.71	0
3	GLC	I	1	3	11,11,12	0.51	0	15,15,17	1.05	1 (6%)
3	FRU	I	2	3	11,12,12	4.92	6 (54%)	10,18,18	0.91	1 (10%)
3	GLC	J	1	3	11,11,12	0.61	0	15,15,17	1.20	1 (6%)
3	FRU	J	2	3	11,12,12	4.36	6 (54%)	10,18,18	1.24	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
3	GLC	D	1	3	-	0/2/19/22	0/1/1/1
3	FRU	D	2	3	-	2/5/24/24	0/1/1/1
3	GLC	E	1	3	-	0/2/19/22	0/1/1/1
3	FRU	E	2	3	-	0/5/24/24	0/1/1/1
3	GLC	F	1	3	-	0/2/19/22	0/1/1/1
3	FRU	F	2	3	-	0/5/24/24	0/1/1/1
3	GLC	G	1	3	-	2/2/19/22	0/1/1/1
3	FRU	G	2	3	-	2/5/24/24	0/1/1/1
3	GLC	H	1	3	-	0/2/19/22	0/1/1/1
3	FRU	H	2	3	-	2/5/24/24	0/1/1/1
3	GLC	I	1	3	-	0/2/19/22	0/1/1/1
3	FRU	I	2	3	-	0/5/24/24	0/1/1/1
3	GLC	J	1	3	-	0/2/19/22	0/1/1/1
3	FRU	J	2	3	-	0/5/24/24	0/1/1/1

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	2	FRU	O5-C2	14.00	1.65	1.43
3	H	2	FRU	O5-C2	13.91	1.65	1.43
3	F	2	FRU	O5-C2	13.41	1.64	1.43
3	E	2	FRU	O5-C2	12.93	1.64	1.43
3	G	2	FRU	O5-C2	12.48	1.63	1.43
3	D	2	FRU	O5-C2	12.33	1.63	1.43
3	J	2	FRU	O5-C2	12.02	1.62	1.43
3	I	2	FRU	O5-C5	-4.71	1.33	1.43
3	G	2	FRU	O5-C5	-4.62	1.33	1.43
3	J	2	FRU	O5-C5	-4.47	1.34	1.43
3	D	2	FRU	O5-C5	-4.42	1.34	1.43
3	H	2	FRU	O5-C5	-4.39	1.34	1.43
3	F	2	FRU	O5-C5	-4.34	1.34	1.43
3	E	2	FRU	O5-C5	-4.34	1.34	1.43
3	H	2	FRU	C1-C2	4.21	1.61	1.52
3	I	2	FRU	C1-C2	4.15	1.61	1.52
3	F	2	FRU	C1-C2	4.12	1.61	1.52
3	D	2	FRU	O4-C4	-4.06	1.32	1.43
3	J	2	FRU	C1-C2	3.98	1.61	1.52
3	H	2	FRU	O4-C4	-3.82	1.33	1.43
3	E	2	FRU	C1-C2	3.77	1.60	1.52
3	D	2	FRU	C1-C2	3.69	1.60	1.52
3	E	2	FRU	O4-C4	-3.63	1.34	1.43
3	F	2	FRU	O4-C4	-3.63	1.34	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	2	FRU	O4-C4	-3.63	1.34	1.43
3	G	2	FRU	O4-C4	-3.62	1.34	1.43
3	I	2	FRU	O4-C4	-3.52	1.34	1.43
3	G	2	FRU	C1-C2	3.31	1.59	1.52
3	I	2	FRU	C4-C5	3.11	1.60	1.53
3	H	2	FRU	C4-C5	2.83	1.60	1.53
3	G	2	FRU	C4-C5	2.74	1.59	1.53
3	D	2	FRU	C4-C5	2.61	1.59	1.53
3	F	2	FRU	C4-C5	2.59	1.59	1.53
3	J	2	FRU	C4-C5	2.46	1.59	1.53
3	E	2	FRU	C4-C5	2.36	1.59	1.53
3	G	2	FRU	O3-C3	2.30	1.47	1.42
3	D	2	FRU	O2-C2	-2.25	1.36	1.40
3	I	2	FRU	O3-C3	2.24	1.47	1.42
3	J	2	FRU	O2-C2	-2.16	1.37	1.40
3	G	2	FRU	O2-C2	-2.01	1.37	1.40

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	1	GLC	C1-O5-C5	3.30	116.61	112.19
3	I	1	GLC	C1-O5-C5	2.94	116.12	112.19
3	E	1	GLC	C1-O5-C5	2.86	116.02	112.19
3	H	1	GLC	C1-O5-C5	2.73	115.84	112.19
3	G	2	FRU	O1-C1-C2	-2.59	105.95	111.67
3	H	1	GLC	C6-C5-C4	-2.53	106.80	113.02
3	H	1	GLC	O5-C5-C6	2.49	112.52	107.66
3	F	2	FRU	C6-C5-C4	-2.39	109.46	115.10
3	E	1	GLC	C6-C5-C4	-2.35	107.24	113.02
3	D	2	FRU	O1-C1-C2	-2.10	107.03	111.67
3	I	2	FRU	O2-C2-O5	2.03	113.22	109.33

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	G	2	FRU	O1-C1-C2-O2
3	D	2	FRU	C4-C5-C6-O6
3	D	2	FRU	O5-C5-C6-O6
3	G	1	GLC	O5-C5-C6-O6
3	H	2	FRU	C4-C5-C6-O6
3	H	2	FRU	O5-C5-C6-O6

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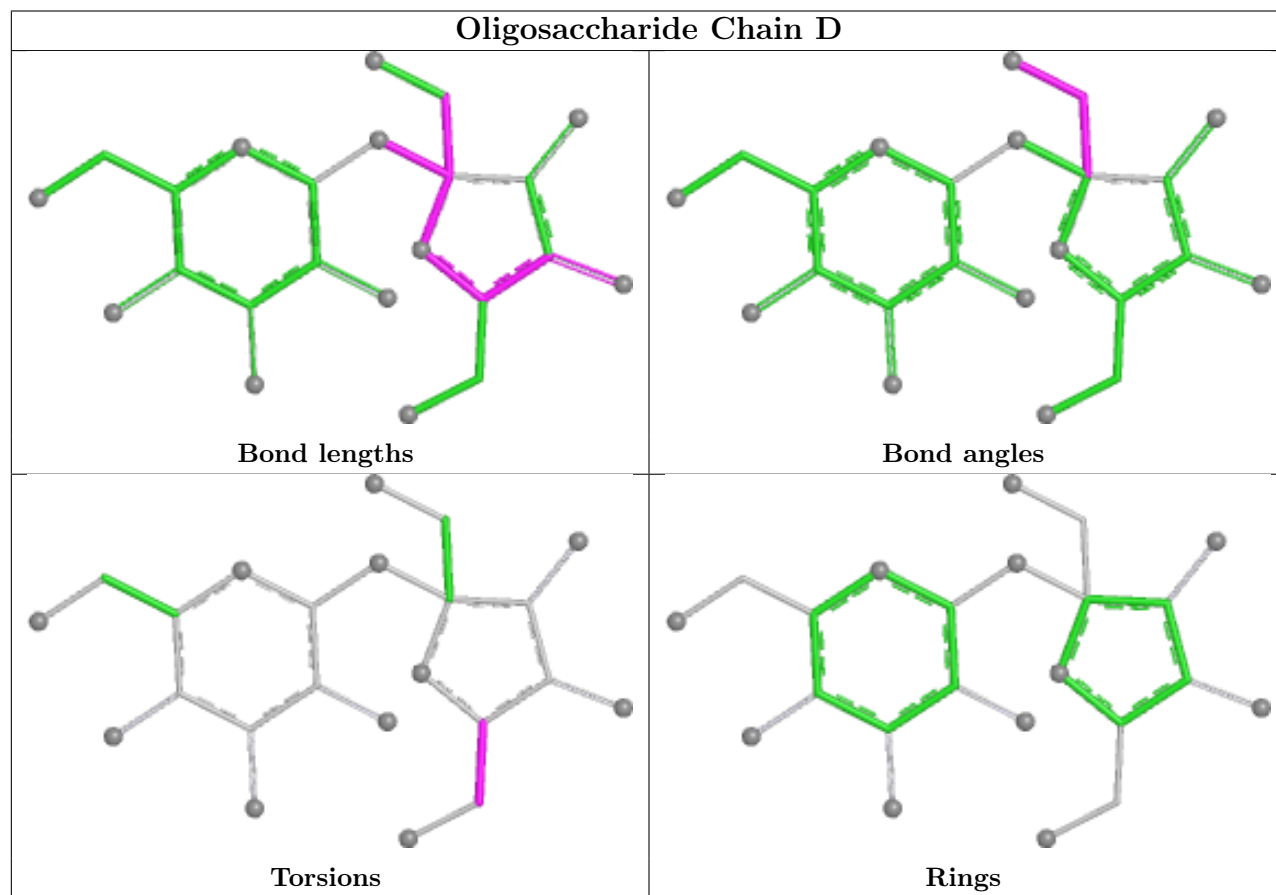
Mol	Chain	Res	Type	Atoms
3	G	2	FRU	O1-C1-C2-O5
3	G	1	GLC	C4-C5-C6-O6

There are no ring outliers.

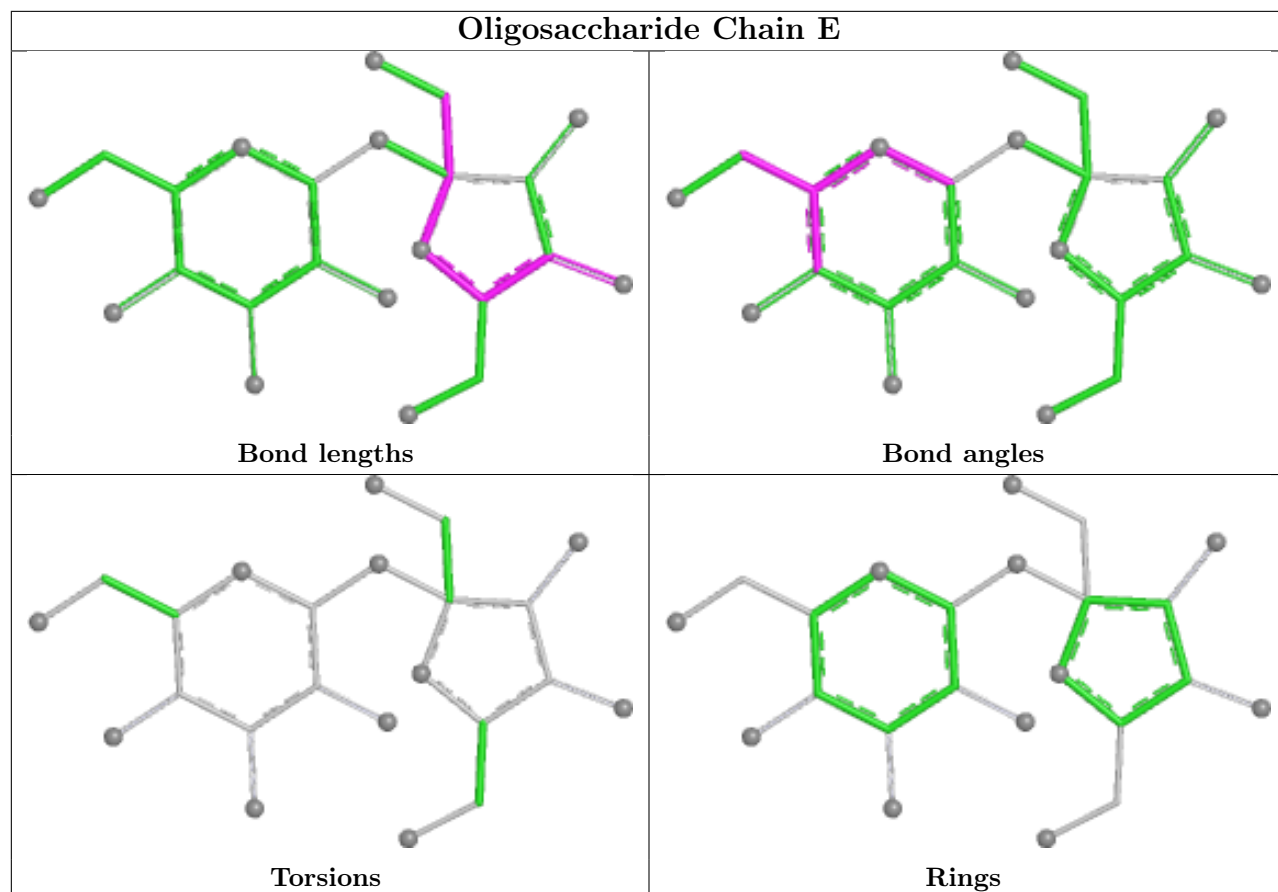
5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	1	GLC	1	0
3	H	2	FRU	2	0
3	I	2	FRU	2	0
3	E	2	FRU	1	0
3	F	2	FRU	1	0

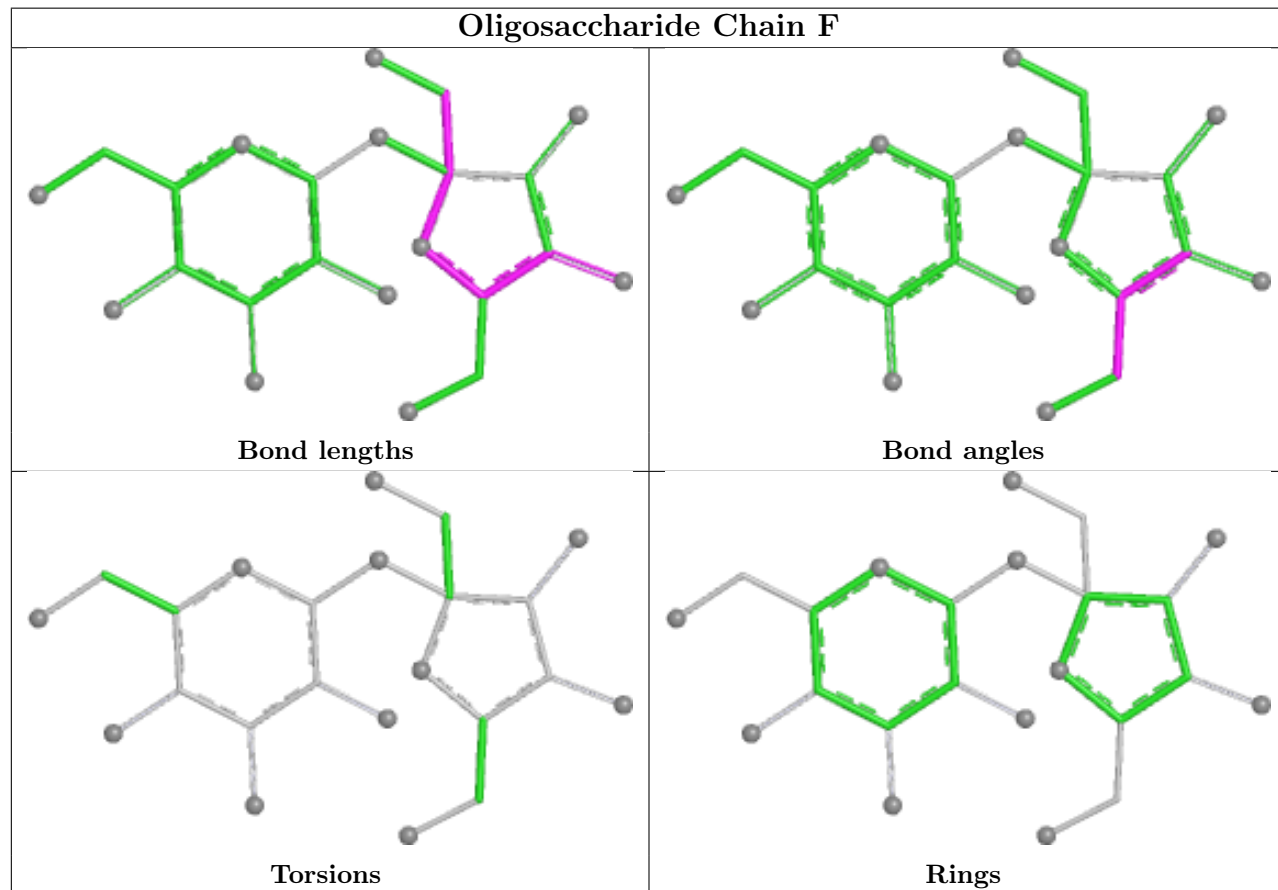
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



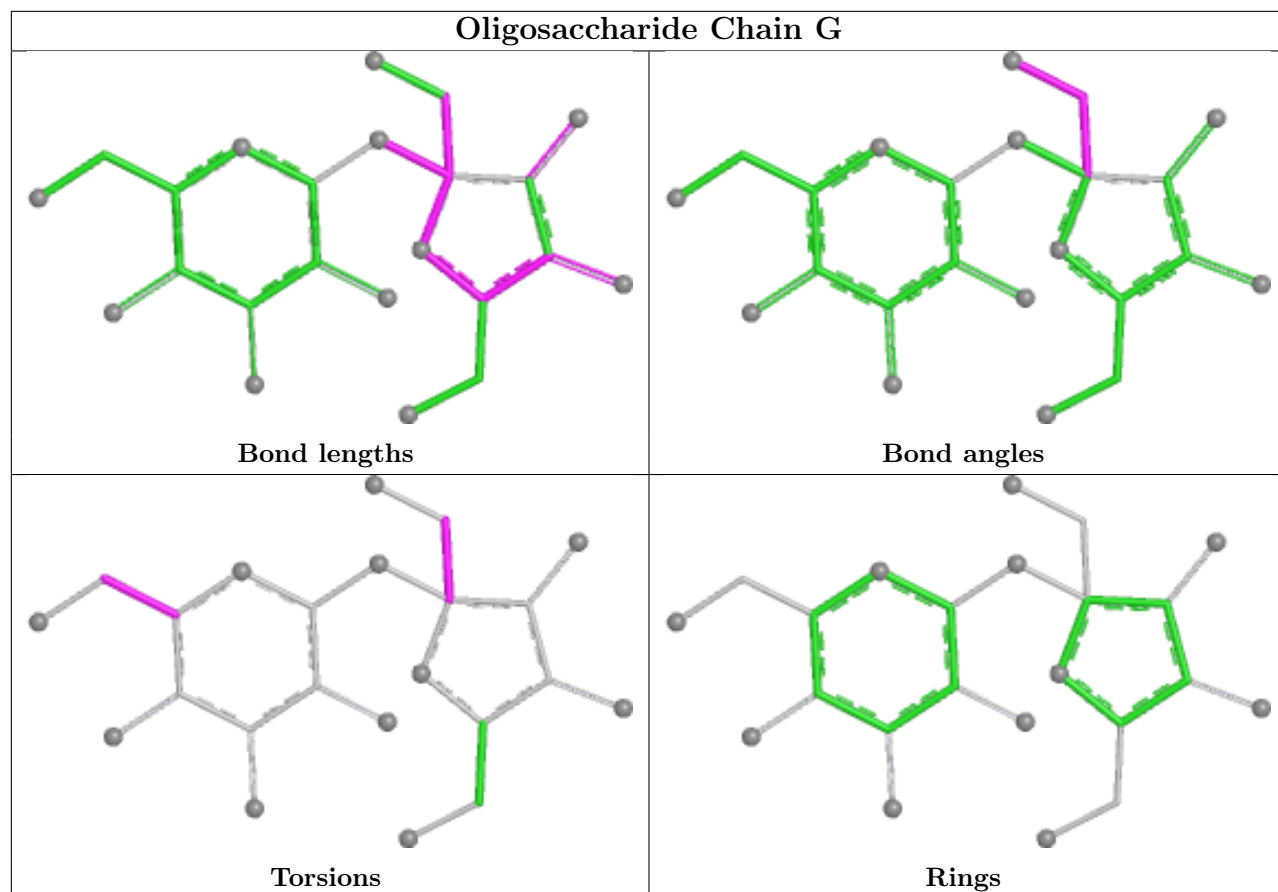
Oligosaccharide Chain E



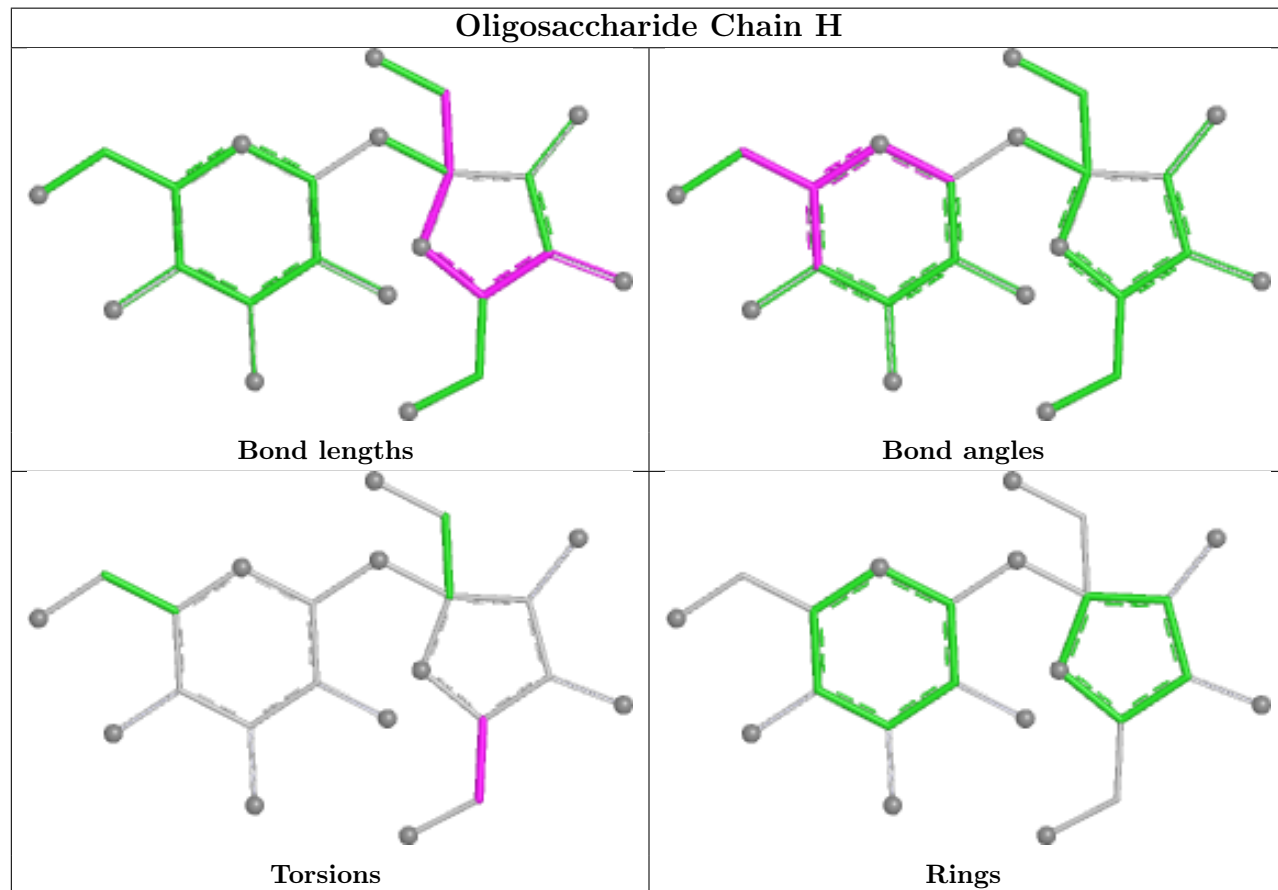
Oligosaccharide Chain F

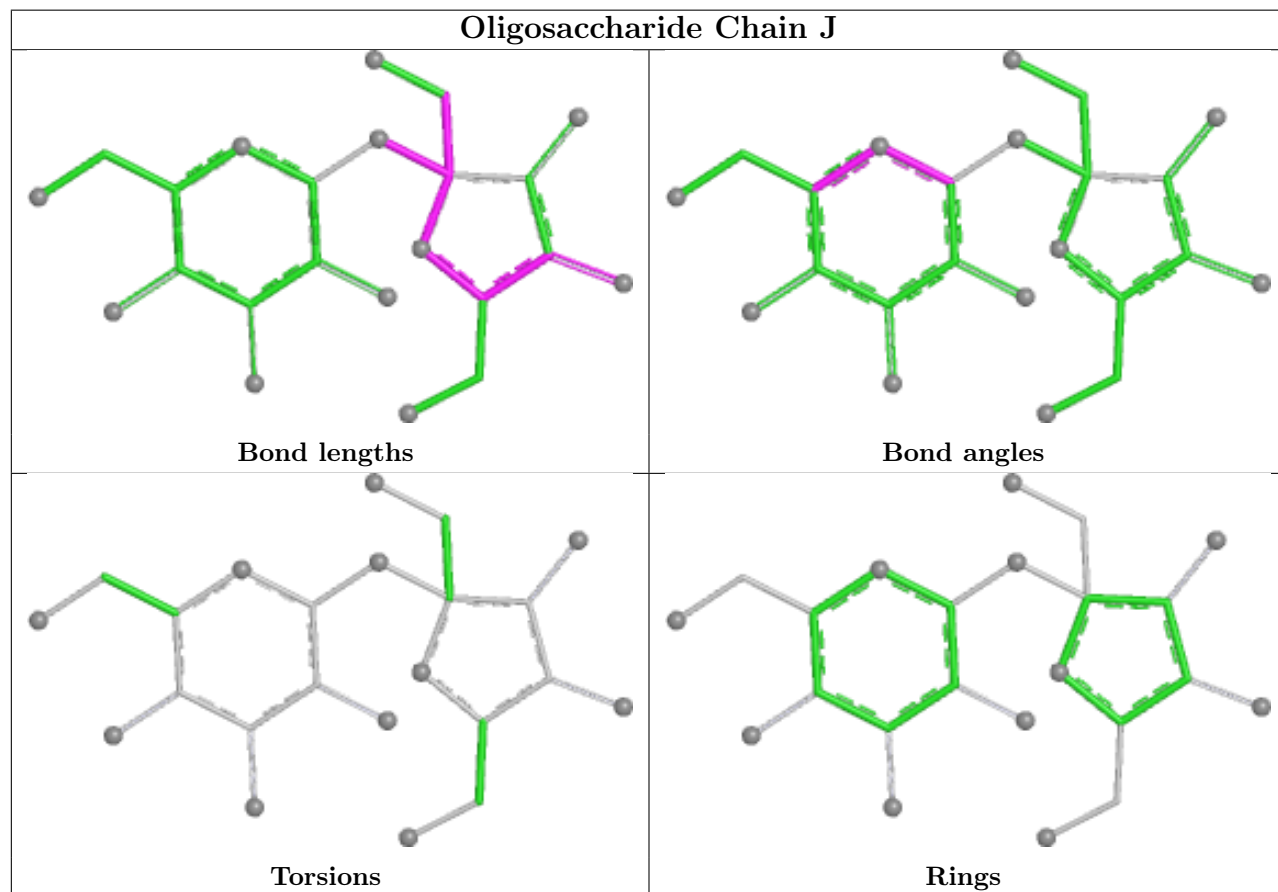
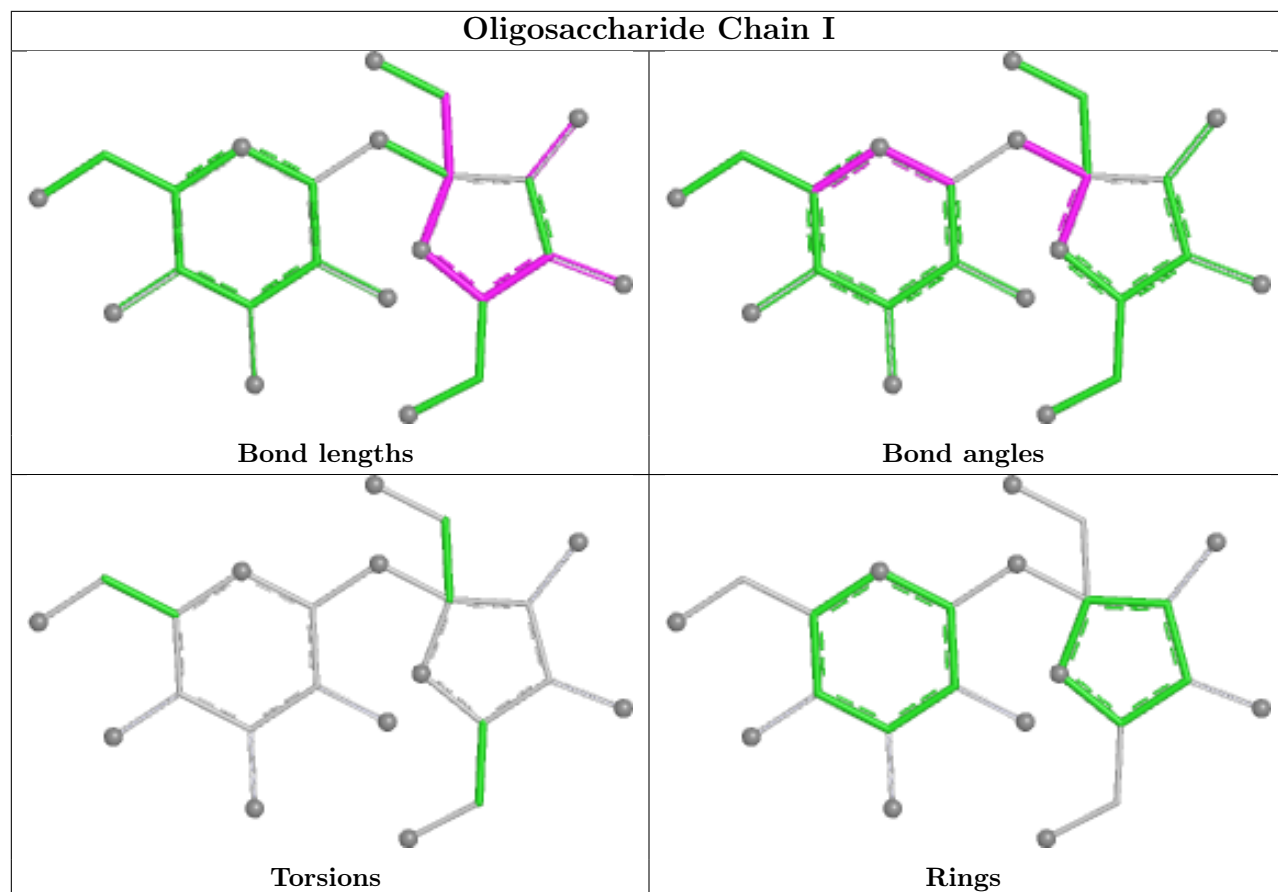


Oligosaccharide Chain G



Oligosaccharide Chain H





5.6 Ligand geometry

7 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	$\# Z > 2$
5	1PE	A	1103	-	12,12,15	0.54	0	11,11,14	0.35	0
4	SO4	B	1102	-	4,4,4	0.27	0	6,6,6	0.15	0
4	SO4	A	1102	-	4,4,4	0.27	0	6,6,6	0.18	0
6	A1AVL	A	1105	-	31,33,33	2.12	8 (25%)	38,49,49	3.10	15 (39%)
4	SO4	A	1101	-	4,4,4	0.28	0	6,6,6	0.13	0
4	SO4	B	1101	-	4,4,4	0.24	0	6,6,6	0.17	0
5	1PE	A	1104	-	12,12,15	0.53	0	11,11,14	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
5	1PE	A	1103	-	-	5/10/10/13	-
6	A1AVL	A	1105	-	-	5/17/19/19	0/3/3/3
5	1PE	A	1104	-	-	4/10/10/13	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1105	A1AVL	C06-N05	6.32	1.46	1.39
6	A	1105	A1AVL	S02-N05	5.42	1.70	1.63
6	A	1105	A1AVL	C11-N09	-4.32	1.32	1.44
6	A	1105	A1AVL	C01-S02	3.10	1.82	1.75
6	A	1105	A1AVL	O22-S23	2.80	1.63	1.58
6	A	1105	A1AVL	O03-S02	2.48	1.47	1.43
6	A	1105	A1AVL	O04-S02	2.42	1.47	1.43
6	A	1105	A1AVL	O24-S23	2.16	1.54	1.45

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1105	A1AVL	O04-S02-O03	-12.62	101.95	118.87
6	A	1105	A1AVL	C12-C13-C15	-6.13	119.89	124.45
6	A	1105	A1AVL	C31-C29-C15	-5.85	120.10	124.45
6	A	1105	A1AVL	C16-C15-C13	-4.18	119.40	122.51
6	A	1105	A1AVL	C29-C15-C13	4.16	120.15	114.54
6	A	1105	A1AVL	C06-N05-S02	-3.68	120.63	125.83
6	A	1105	A1AVL	O22-C21-C27	3.48	125.52	118.70
6	A	1105	A1AVL	C01-S02-N05	2.90	109.83	106.56
6	A	1105	A1AVL	C16-C15-C29	-2.78	120.45	122.51
6	A	1105	A1AVL	O25-S23-O24	2.54	117.44	108.56
6	A	1105	A1AVL	F14-C13-C15	2.49	120.75	117.64
6	A	1105	A1AVL	O22-C21-C20	-2.46	113.87	118.70
6	A	1105	A1AVL	O17-C16-C15	-2.43	101.58	107.66
6	A	1105	A1AVL	O04-S02-C01	2.42	112.09	108.26
6	A	1105	A1AVL	O03-S02-N05	2.03	111.16	107.10

There are no chirality outliers.

All (14) torsion outliers are listed below:

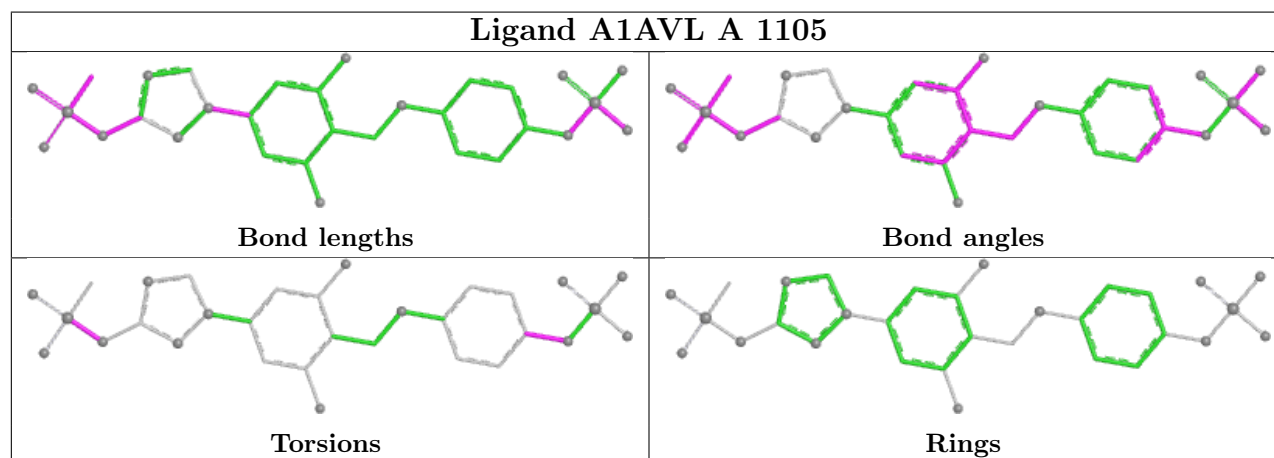
Mol	Chain	Res	Type	Atoms
6	A	1105	A1AVL	C06-N05-S02-C01
6	A	1105	A1AVL	C06-N05-S02-O04
5	A	1104	1PE	OH5-C14-C24-OH4
5	A	1104	1PE	OH2-C12-C22-OH3
5	A	1104	1PE	C23-C13-OH4-C24
5	A	1103	1PE	OH4-C13-C23-OH3
5	A	1103	1PE	OH5-C14-C24-OH4
5	A	1103	1PE	OH7-C16-C26-OH6
6	A	1105	A1AVL	C20-C21-O22-S23
6	A	1105	A1AVL	C27-C21-O22-S23
5	A	1104	1PE	C13-C23-OH3-C22
5	A	1103	1PE	OH6-C15-C25-OH5
6	A	1105	A1AVL	C06-N05-S02-O03
5	A	1103	1PE	C16-C26-OH6-C15

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1104	1PE	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.



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	467/512 (91%)	-0.12	10 (2%) 63 63	20, 32, 49, 75	3 (0%)
2	B	468/512 (91%)	-0.03	17 (3%) 46 45	17, 31, 53, 89	1 (0%)
2	C	10/512 (1%)	3.52	7 (70%) 0 0	60, 76, 87, 100	0
All	All	945/1536 (61%)	-0.04	34 (3%) 46 45	17, 32, 53, 100	4 (0%)

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	584	ILE	5.3
2	C	583	ILE	4.9
2	C	577	PHE	4.8
1	A	841	TYR	4.6
2	C	585	THR	4.5
2	C	576	ARG	4.3
1	A	838	PHE	3.6
2	B	902	PHE	3.6
2	C	580	PHE	3.1
2	B	893	GLU	3.0
2	B	898	PHE	3.0
2	B	621	THR	3.0
1	A	596	GLU	2.9
1	A	857	GLU	2.9
2	B	620	VAL	2.9
2	C	581	GLY	2.9
1	A	620	VAL	2.8
1	A	836[A]	MET	2.7
2	B	858	VAL	2.7
2	B	602	GLU	2.6
2	B	955	ASP	2.6
1	A	898	PHE	2.6
2	B	894	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	1061	ARG	2.5
2	B	861	ASP	2.4
2	B	617	ALA	2.2
2	B	619	GLU	2.2
2	B	900	ARG	2.2
2	B	890	ALA	2.2
1	A	1063	LYS	2.2
1	A	636	ARG	2.2
2	B	616	ALA	2.1
2	B	857	GLU	2.1
2	B	951	GLU	2.1

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](#)

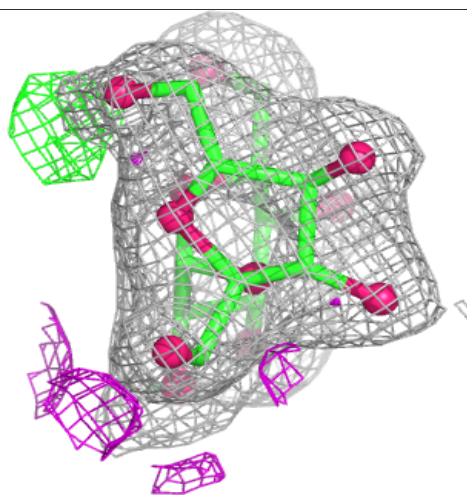
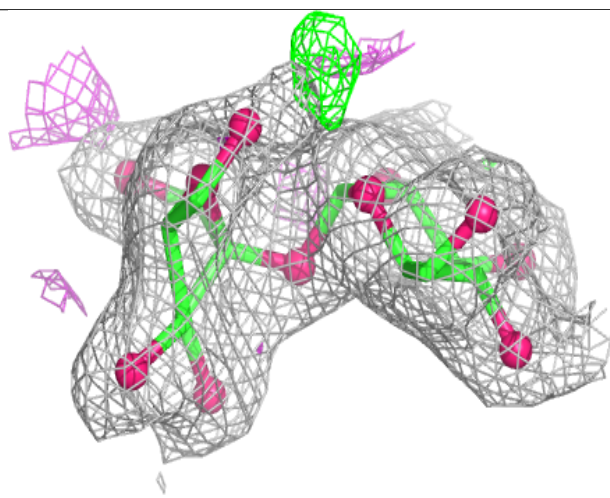
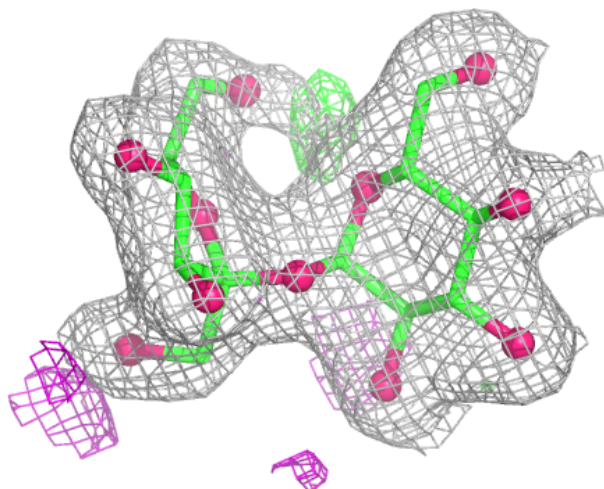
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	FRU	I	2	12/12	0.78	0.16	55,66,73,73	0
3	FRU	H	2	12/12	0.79	0.14	46,57,60,62	0
3	GLC	H	1	11/12	0.84	0.13	37,43,51,55	0
3	GLC	I	1	11/12	0.89	0.11	36,48,62,63	0
3	FRU	D	2	12/12	0.90	0.11	32,37,44,51	0
3	GLC	E	1	11/12	0.90	0.10	31,36,44,55	0
3	FRU	F	2	12/12	0.91	0.09	36,44,48,48	0
3	GLC	G	1	11/12	0.91	0.09	41,43,48,48	0
3	FRU	G	2	12/12	0.93	0.10	34,41,45,46	0
3	GLC	F	1	11/12	0.94	0.08	34,39,42,45	0
3	FRU	E	2	12/12	0.94	0.07	27,32,34,38	0
3	GLC	J	1	11/12	0.94	0.08	26,31,39,49	0
3	GLC	D	1	11/12	0.96	0.07	26,33,37,37	0
3	FRU	J	2	12/12	0.96	0.06	24,28,31,31	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.

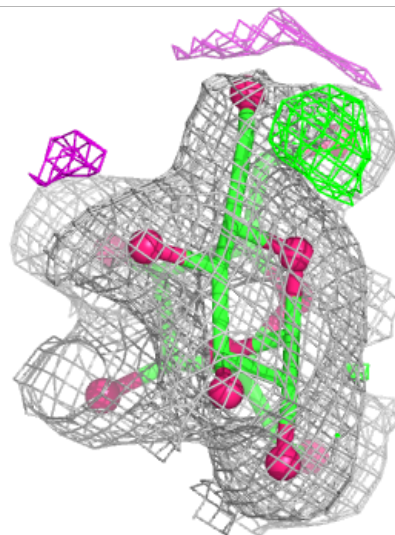
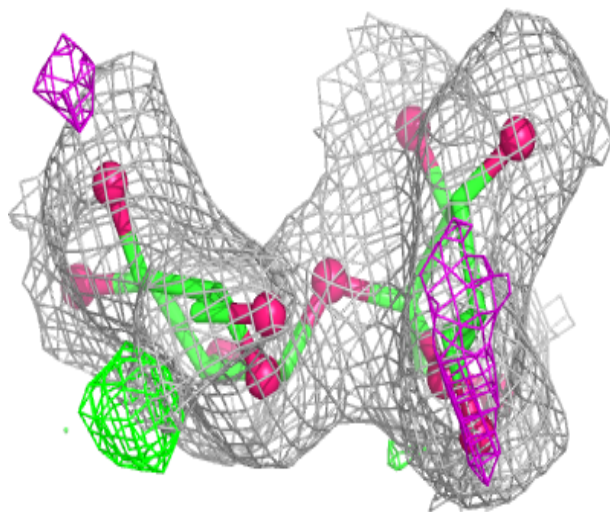
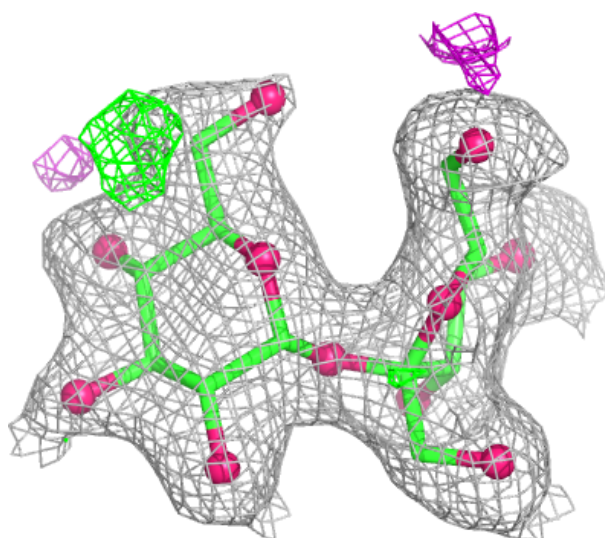
Electron density around Chain D:

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



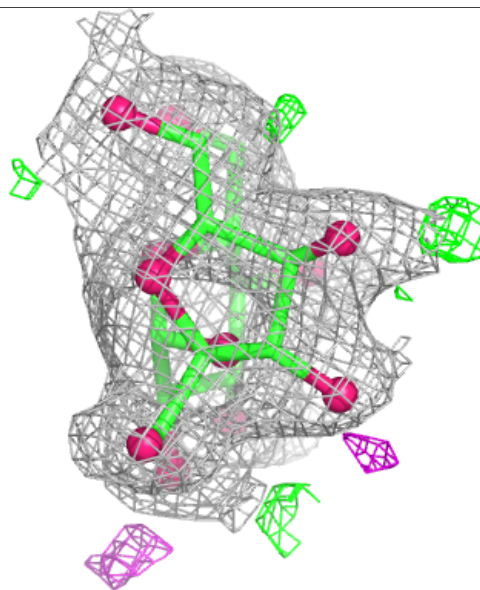
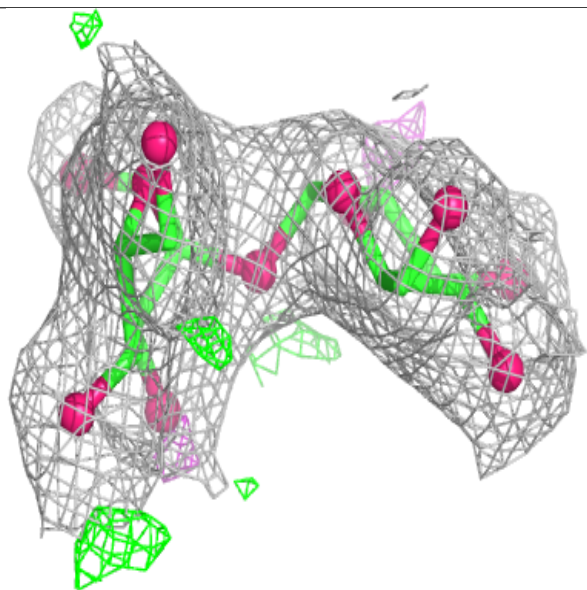
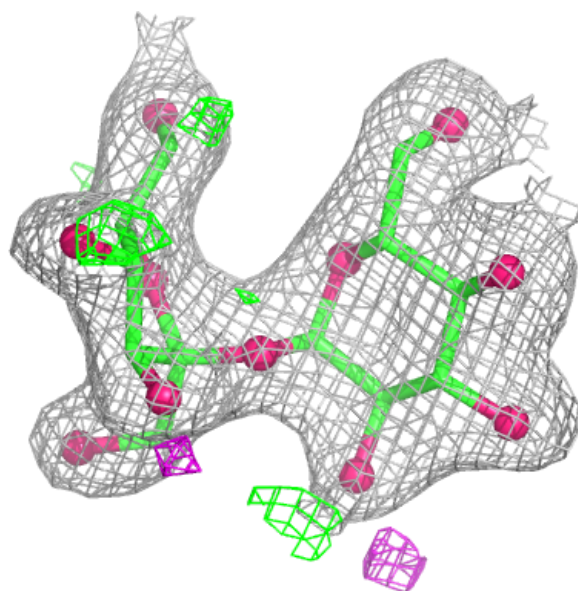
Electron density around Chain E:

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



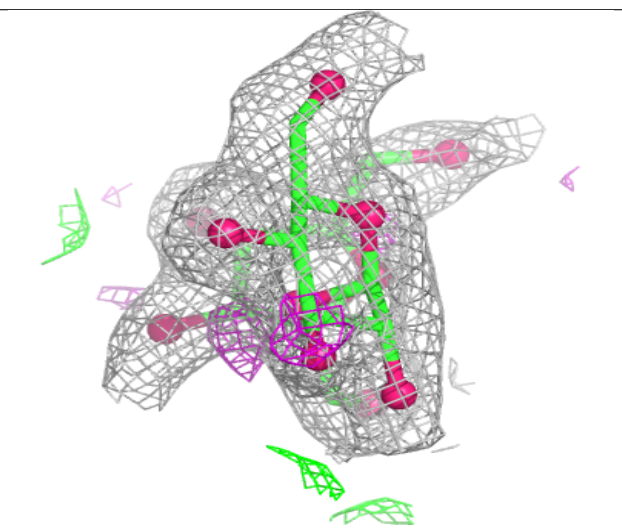
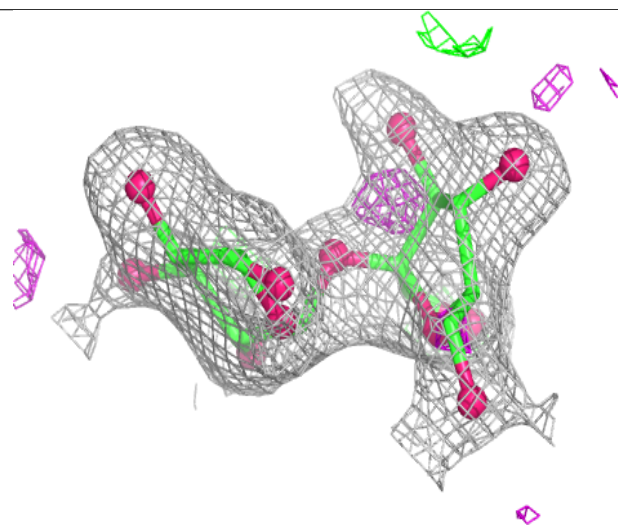
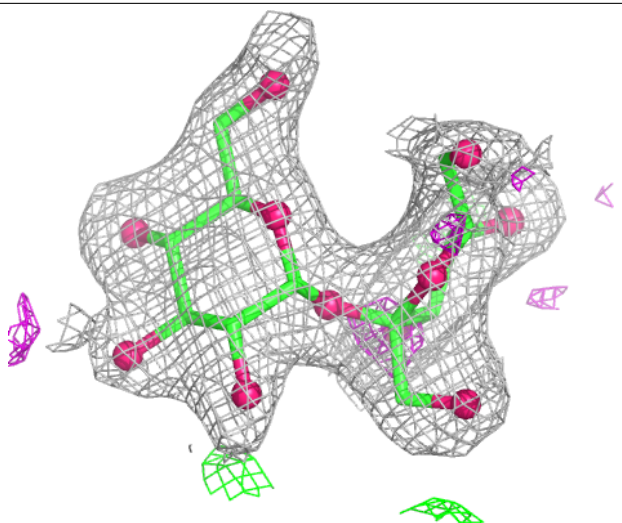
Electron density around Chain F:

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



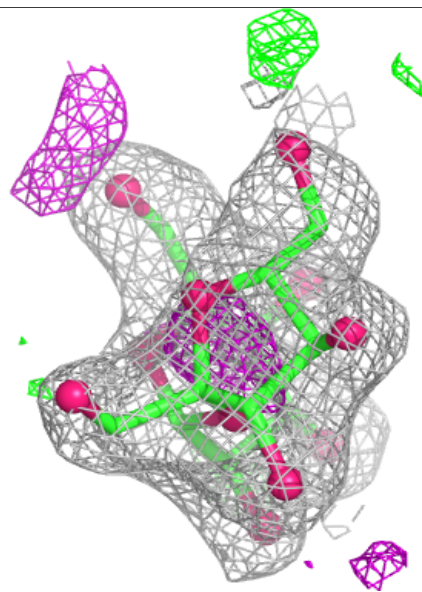
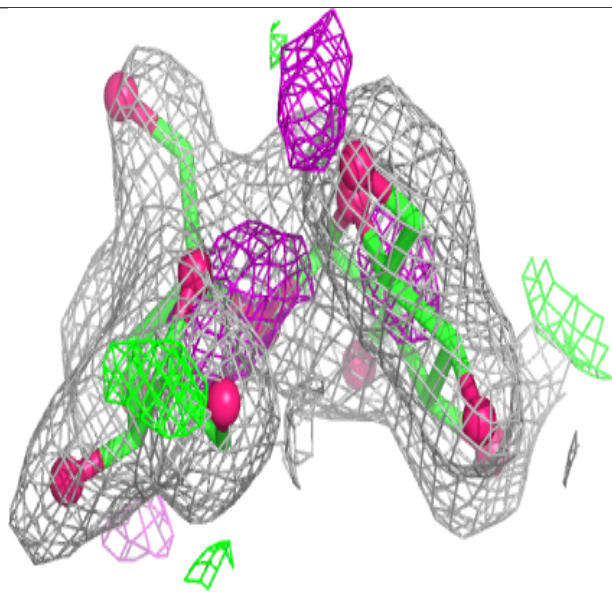
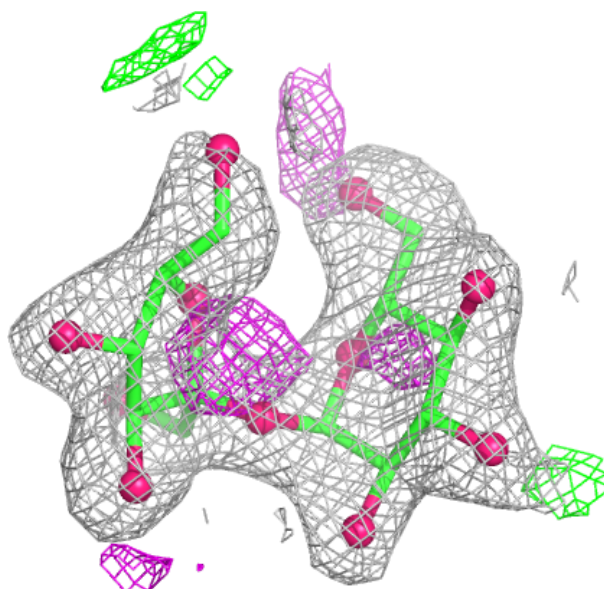
Electron density around Chain G:

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



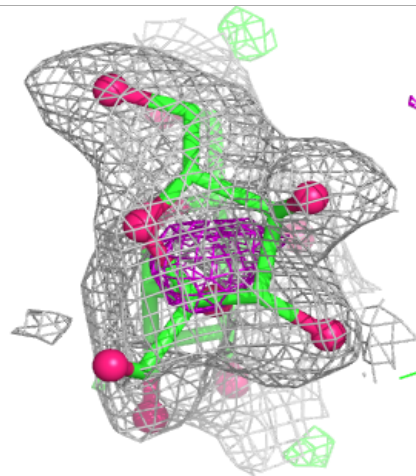
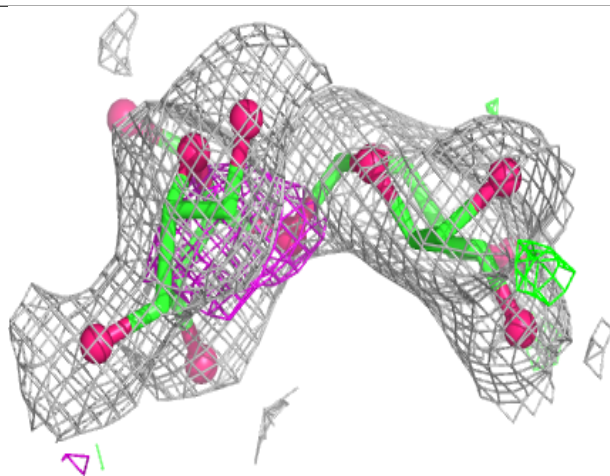
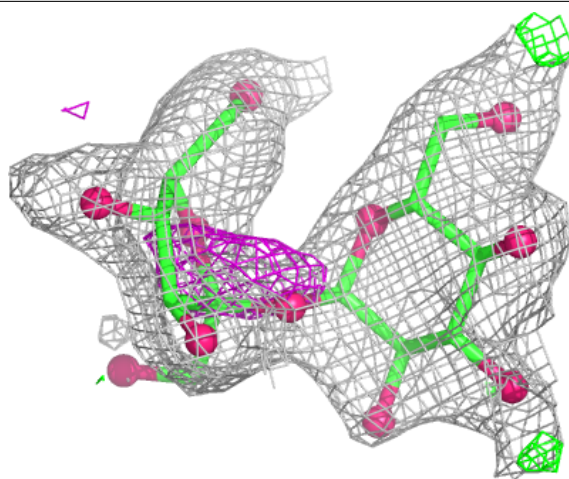
Electron density around Chain H:

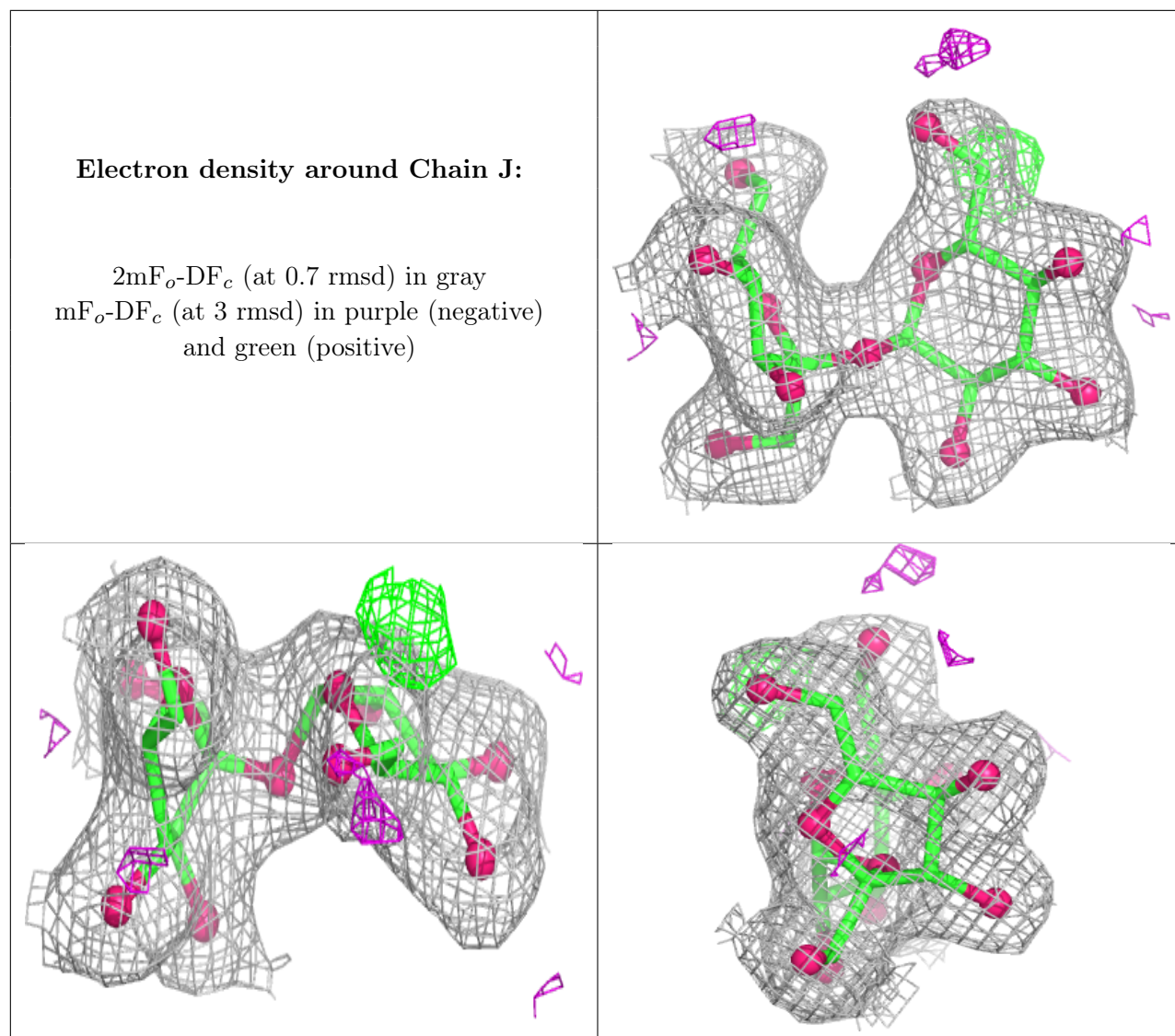
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain I:

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



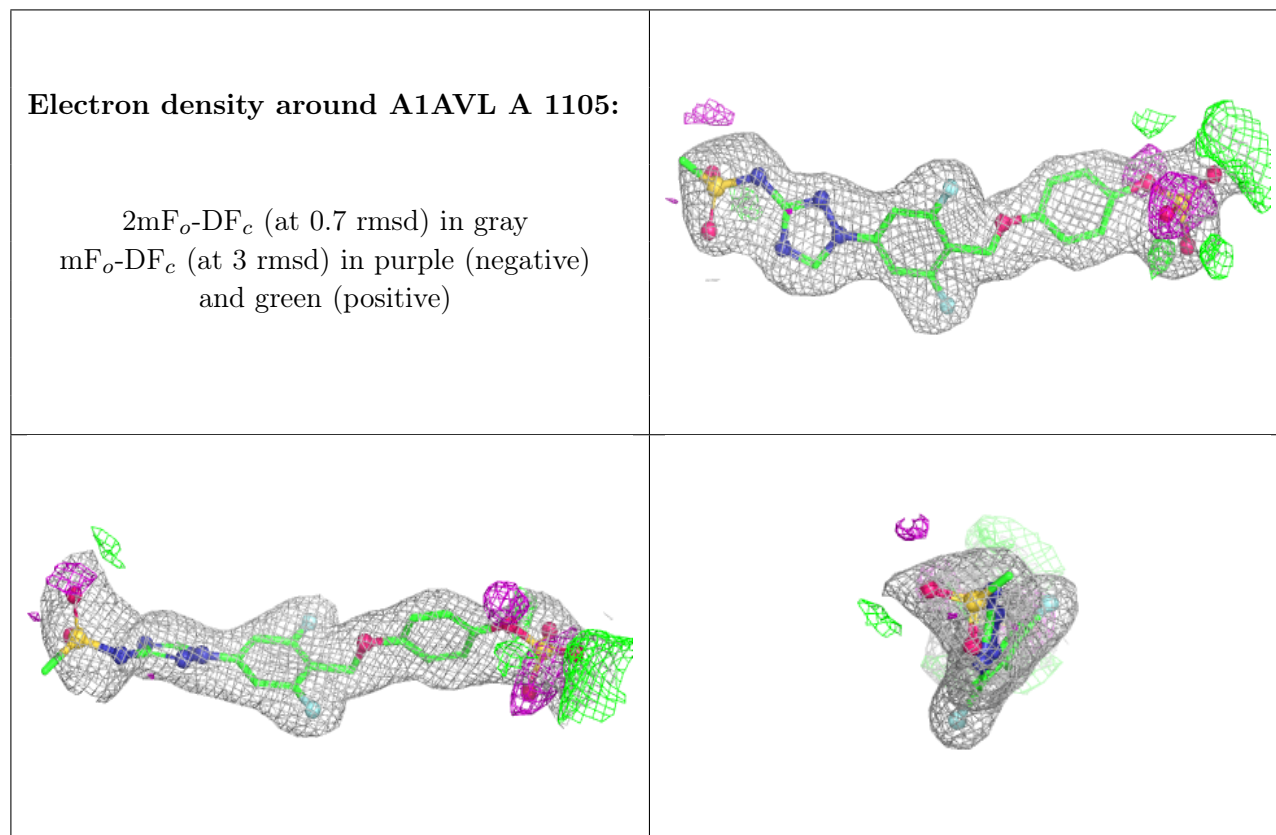


6.4 Ligands ⓘ

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(Å ²)	Q<0.9
4	SO4	B	1102	5/5	0.79	0.10	73,76,84,86	0
4	SO4	A	1101	5/5	0.82	0.12	59,71,76,80	0
5	1PE	A	1104	13/16	0.83	0.16	52,60,66,66	0
6	A1AVL	A	1105	31/31	0.84	0.14	36,44,79,85	0
5	1PE	A	1103	13/16	0.87	0.16	41,49,59,60	0
4	SO4	A	1102	5/5	0.89	0.13	64,65,73,81	0
4	SO4	B	1101	5/5	1.00	0.03	26,27,28,28	5

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.



6.5 Other polymers [i](#)

There are no such residues in this entry.