



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 9, 2020 – 10:35 PM BST

PDB ID : 4TZ3
Title : Ensemble refinement of the E502A variant of sacteLam55A from Streptomyces sp. SirexAA-E in complex with laminaritetraose
Authors : Bianchetti, C.M.; Takasuka, T.E.; Yik, E.J.; Bergeman, L.F.; Fox, B.G.
Deposited on : 2014-07-09
Resolution : 1.90 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

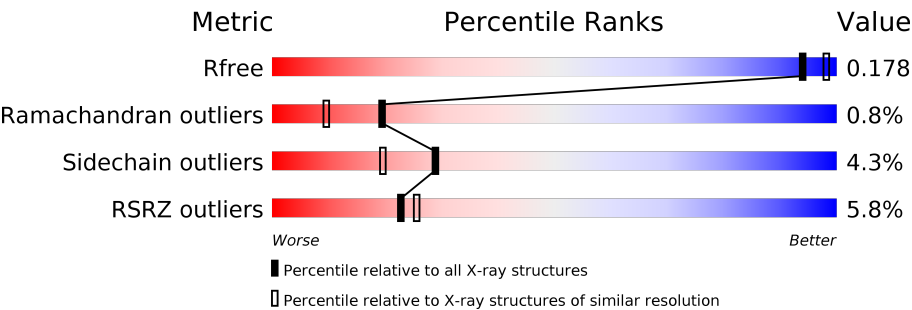
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
<i>R_{free}</i>	130704	6207 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 on 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	1-A	549	<div><div>6%</div><div>95%</div><div>5%</div></div>
1	10-A	549	<div><div>6%</div><div>95%</div><div>5%</div></div>
1	11-A	549	<div><div>6%</div><div>95%</div><div>5%</div></div>
1	12-A	549	<div><div>6%</div><div>95%</div><div>5%</div></div>
1	13-A	549	<div><div>6%</div><div>95%</div><div>5%</div></div>
1	14-A	549	<div><div>6%</div><div>94%</div><div>5%</div></div>
1	15-A	549	<div><div>6%</div><div>94%</div><div>5%</div></div>

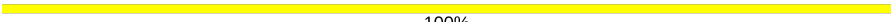
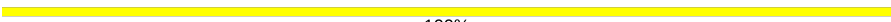
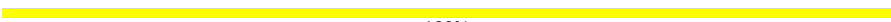
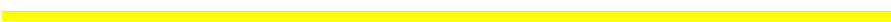








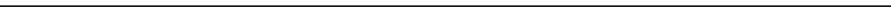


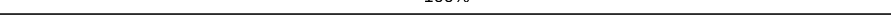
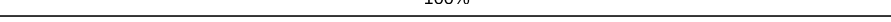
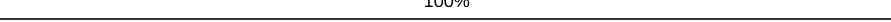
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Mol	Chain	Length	Quality of chain
1	16-A	549	
1	17-A	549	
1	18-A	549	
1	19-A	549	
1	2-A	549	
1	20-A	549	
1	21-A	549	
1	22-A	549	
1	23-A	549	
1	24-A	549	
1	25-A	549	
1	3-A	549	
1	4-A	549	
1	5-A	549	
1	6-A	549	
1	7-A	549	
1	8-A	549	
1	9-A	549	
2	1-B	4	
2	10-B	4	
2	11-B	4	
2	12-B	4	
2	13-B	4	
2	14-B	4	
2	15-B	4	

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Mol	Chain	Length	Quality of chain
2	16-B	4	 100%
2	17-B	4	 100%
2	18-B	4	 100%
2	19-B	4	 100%
2	2-B	4	 100%
2	20-B	4	 100%
2	21-B	4	 100%
2	22-B	4	 100%
2	23-B	4	 100%
2	24-B	4	 100%
2	25-B	4	 100%
2	3-B	4	 100%
2	4-B	4	 100%
2	5-B	4	 100%
2	6-B	4	 100%
2	7-B	4	 100%
2	8-B	4	 100%
2	9-B	4	 100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 210296 atoms, of which 98125 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 Putative secreted protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	1-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	2-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	3-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	4-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	5-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	6-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	7-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	8-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	9-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	10-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	11-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	12-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	13-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	14-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	15-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	16-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			

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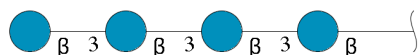
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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	17-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	18-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	19-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	20-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	21-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	22-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	23-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	24-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			
1	25-A	549	Total	C	H	N	O	S	0	0	0
			8074	2629	3907	708	825	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	502	ALA	GLU	engineered mutation	UNP G2NFJ9

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



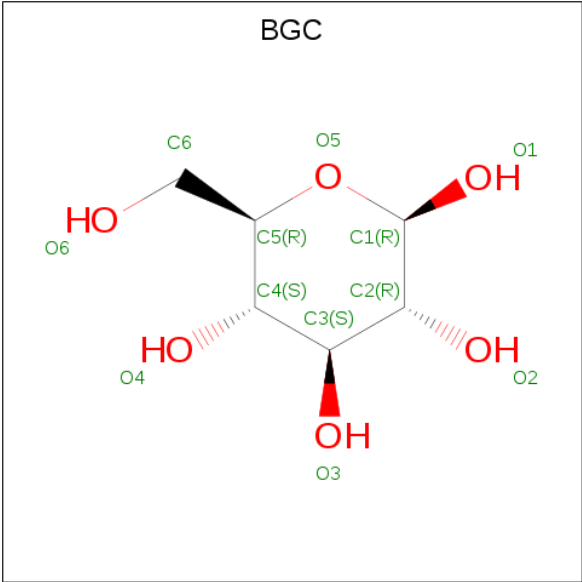
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	1-B	4	Total	C	O	0	0	0
			45	24	21			
2	2-B	4	Total	C	O	0	0	0
			45	24	21			
2	3-B	4	Total	C	O	0	0	0
			45	24	21			
2	4-B	4	Total	C	O	0	0	0
			45	24	21			
2	5-B	4	Total	C	O	0	0	0
			45	24	21			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	6-B	4	Total 45	C 24	O 21	0	0	0
2	7-B	4	Total 45	C 24	O 21	0	0	0
2	8-B	4	Total 45	C 24	O 21	0	0	0
2	9-B	4	Total 45	C 24	O 21	0	0	0
2	10-B	4	Total 45	C 24	O 21	0	0	0
2	11-B	4	Total 45	C 24	O 21	0	0	0
2	12-B	4	Total 45	C 24	O 21	0	0	0
2	13-B	4	Total 45	C 24	O 21	0	0	0
2	14-B	4	Total 45	C 24	O 21	0	0	0
2	15-B	4	Total 45	C 24	O 21	0	0	0
2	16-B	4	Total 45	C 24	O 21	0	0	0
2	17-B	4	Total 45	C 24	O 21	0	0	0
2	18-B	4	Total 45	C 24	O 21	0	0	0
2	19-B	4	Total 45	C 24	O 21	0	0	0
2	20-B	4	Total 45	C 24	O 21	0	0	0
2	21-B	4	Total 45	C 24	O 21	0	0	0
2	22-B	4	Total 45	C 24	O 21	0	0	0
2	23-B	4	Total 45	C 24	O 21	0	0	0
2	24-B	4	Total 45	C 24	O 21	0	0	0
2	25-B	4	Total 45	C 24	O 21	0	0	0

- Molecule 3 is beta-D-glucopyranose (three-letter code: BGC) (formula: C₆H₁₂O₆).



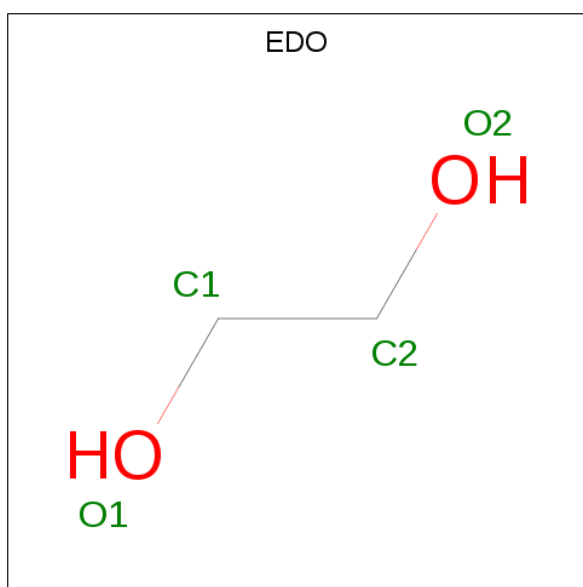
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	1-A	1	Total	C	O	0	0
			11	6	5		
3	2-A	1	Total	C	O	0	0
			11	6	5		
3	3-A	1	Total	C	O	0	0
			11	6	5		
3	4-A	1	Total	C	O	0	0
			11	6	5		
3	5-A	1	Total	C	O	0	0
			11	6	5		
3	6-A	1	Total	C	O	0	0
			11	6	5		
3	7-A	1	Total	C	O	0	0
			11	6	5		
3	8-A	1	Total	C	O	0	0
			11	6	5		
3	9-A	1	Total	C	O	0	0
			11	6	5		
3	10-A	1	Total	C	O	0	0
			11	6	5		
3	11-A	1	Total	C	O	0	0
			11	6	5		
3	12-A	1	Total	C	O	0	0
			11	6	5		
3	13-A	1	Total	C	O	0	0
			11	6	5		
3	14-A	1	Total	C	O	0	0
			11	6	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	15-A	1	Total	C	O	0	0
			11	6	5		
3	16-A	1	Total	C	O	0	0
			11	6	5		
3	17-A	1	Total	C	O	0	0
			11	6	5		
3	18-A	1	Total	C	O	0	0
			11	6	5		
3	19-A	1	Total	C	O	0	0
			11	6	5		
3	20-A	1	Total	C	O	0	0
			11	6	5		
3	21-A	1	Total	C	O	0	0
			11	6	5		
3	22-A	1	Total	C	O	0	0
			11	6	5		
3	23-A	1	Total	C	O	0	0
			11	6	5		
3	24-A	1	Total	C	O	0	0
			11	6	5		
3	25-A	1	Total	C	O	0	0
			11	6	5		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	1-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	2-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	3-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	4-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	5-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	6-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	7-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	8-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	9-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	10-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	11-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	12-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	13-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	14-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	15-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	16-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	17-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	18-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	19-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	20-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	21-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	22-A	1	Total	C	H	O	0	0
			10	2	6	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	23-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	24-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	25-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	1-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	2-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	3-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	4-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	5-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	6-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	7-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	8-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	9-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	10-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	11-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	12-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	13-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	14-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	15-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	16-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	17-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	18-A	1	Total	C	H	O	0	0
			10	2	6	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	19-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	20-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	21-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	22-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	23-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	24-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	25-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	1-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	2-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	3-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	4-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	5-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	6-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	7-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	8-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	9-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	10-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	11-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	12-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	13-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	14-A	1	Total	C	H	O	0	0
			10	2	6	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	15-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	16-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	17-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	18-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	19-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	20-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	21-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	22-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	23-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	24-A	1	Total	C	H	O	0	0
			10	2	6	2		
4	25-A	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	1-A	253	Total	O	0	0
			253	253		
5	2-A	248	Total	O	0	0
			248	248		
5	3-A	259	Total	O	0	0
			259	259		
5	4-A	249	Total	O	0	0
			249	249		
5	5-A	246	Total	O	0	0
			246	246		
5	6-A	245	Total	O	0	0
			245	245		
5	7-A	258	Total	O	0	0
			258	258		
5	8-A	264	Total	O	0	0
			264	264		

Continued on next page...

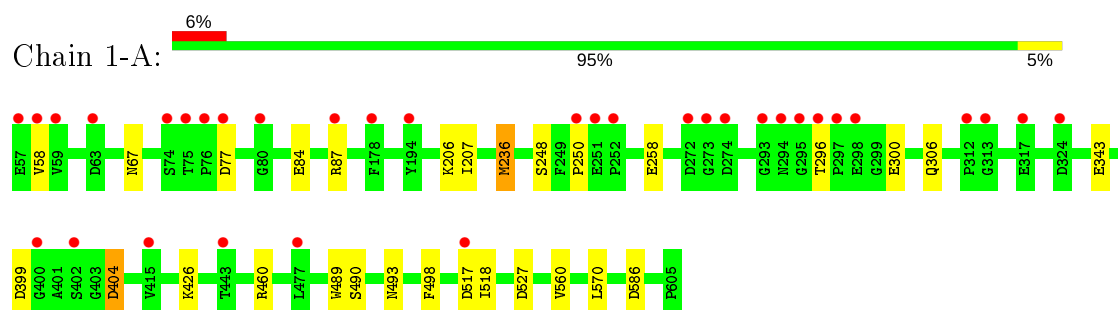
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	9-A	251	Total	O	0	0
			251	251		
5	10-A	248	Total	O	0	0
			248	248		
5	11-A	264	Total	O	0	0
			264	264		
5	12-A	246	Total	O	0	0
			246	246		
5	13-A	247	Total	O	0	0
			247	247		
5	14-A	242	Total	O	0	0
			242	242		
5	15-A	262	Total	O	0	0
			262	262		
5	16-A	252	Total	O	0	0
			252	252		
5	17-A	243	Total	O	0	0
			243	243		
5	18-A	255	Total	O	0	0
			255	255		
5	19-A	274	Total	O	0	0
			274	274		
5	20-A	242	Total	O	0	0
			242	242		
5	21-A	240	Total	O	0	0
			240	240		
5	22-A	250	Total	O	0	0
			250	250		
5	23-A	259	Total	O	0	0
			259	259		
5	24-A	250	Total	O	0	0
			250	250		
5	25-A	249	Total	O	0	0
			249	249		

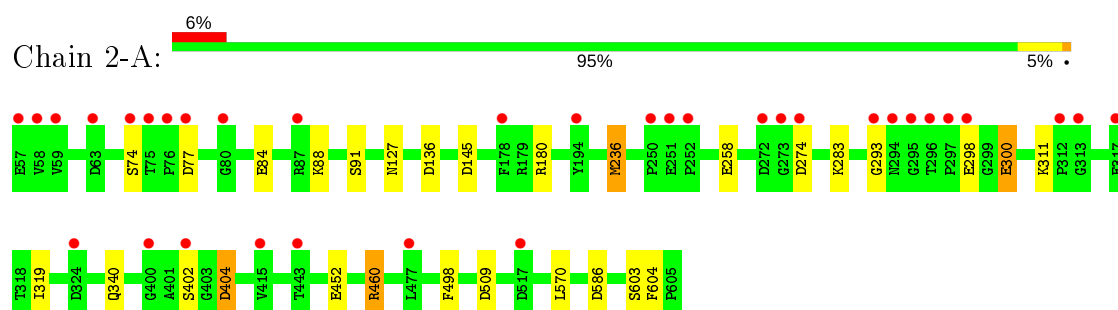
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

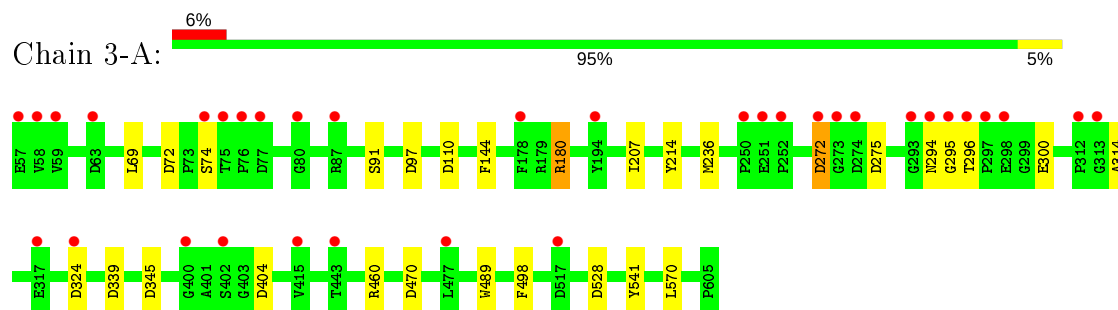
- Molecule 1: Putative secreted protein



- Molecule 1: Putative secreted protein

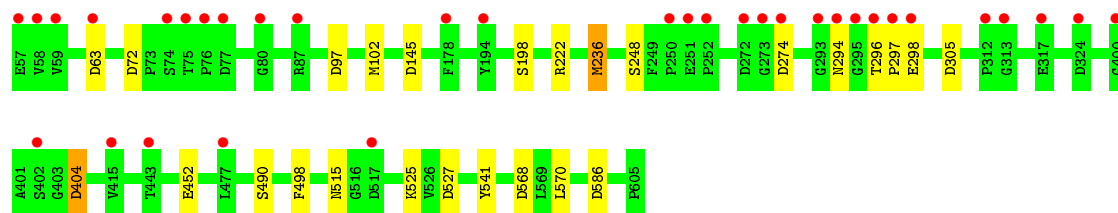


- Molecule 1: Putative secreted protein

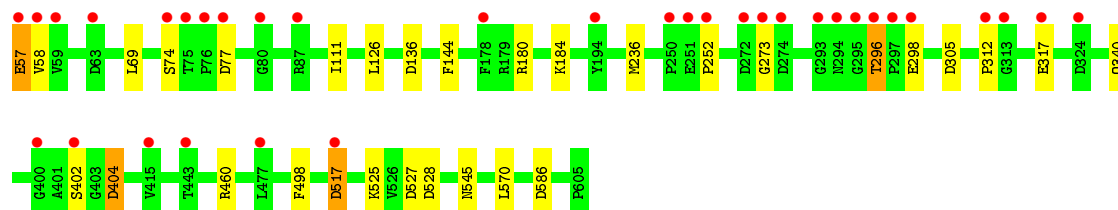


- Molecule 1: Putative secreted protein

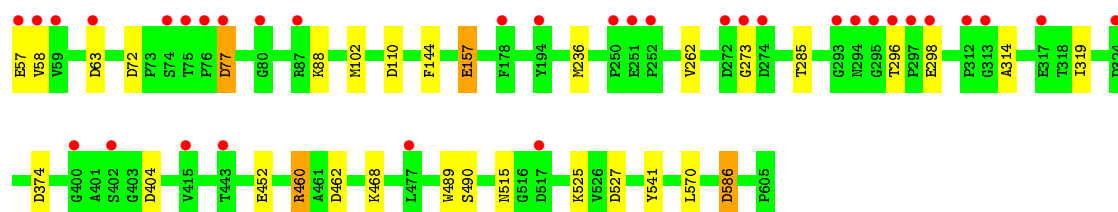




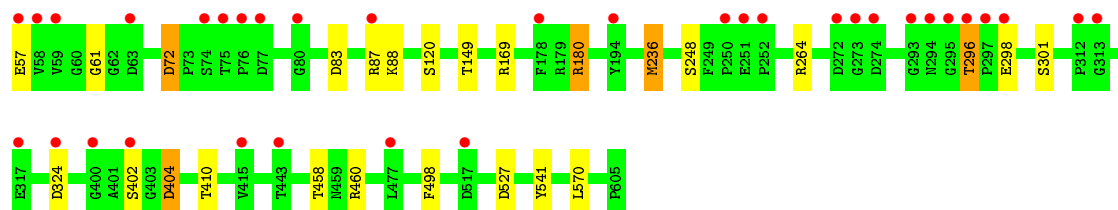
- Molecule 1: Putative secreted protein



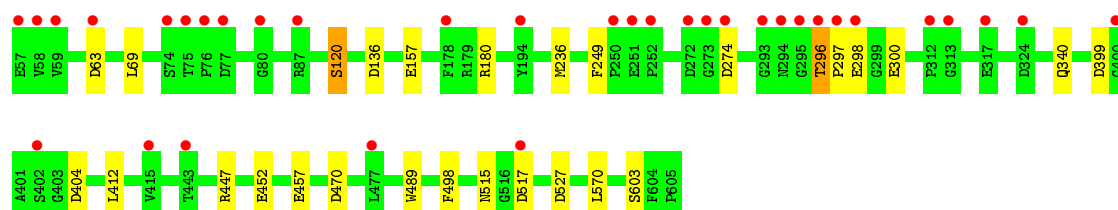
- Molecule 1: Putative secreted protein



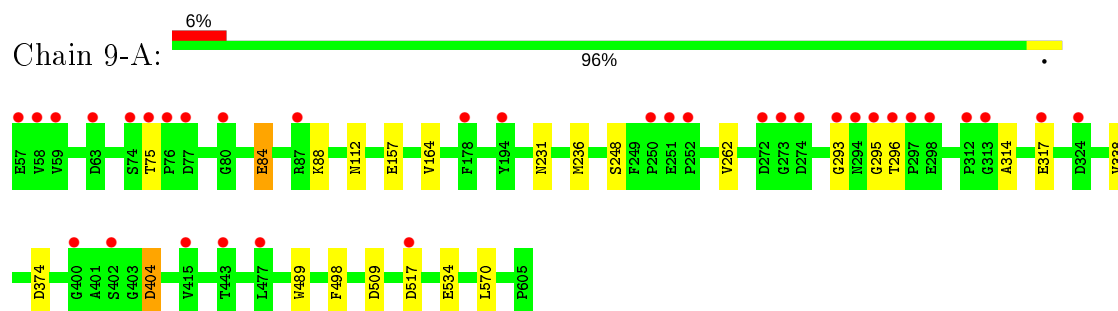
- Molecule 1: Putative secreted protein



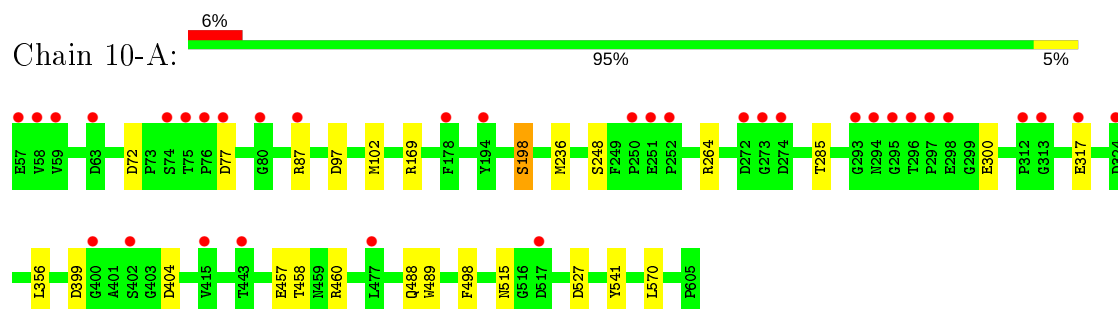
- Molecule 1: Putative secreted protein



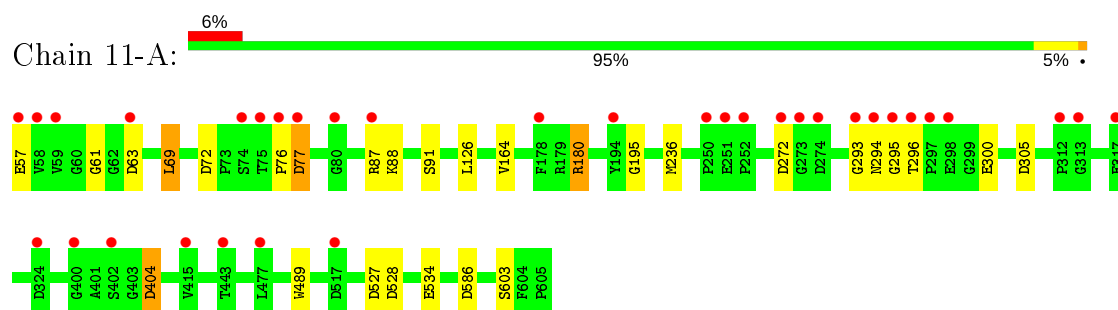
- Molecule 1: Putative secreted protein



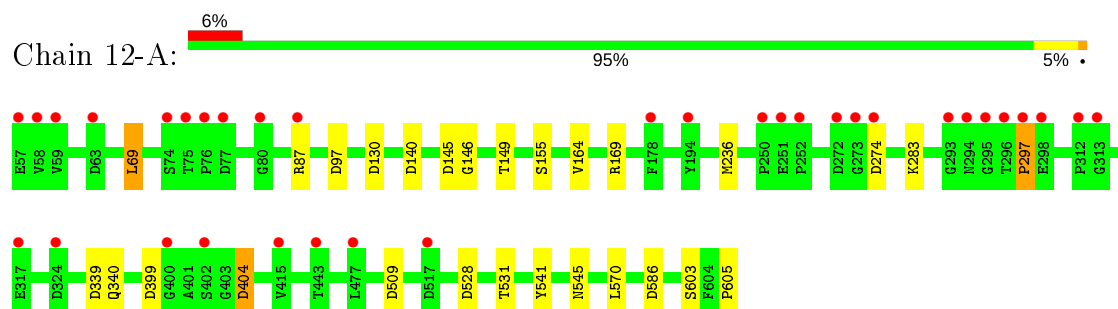
- Molecule 1: Putative secreted protein



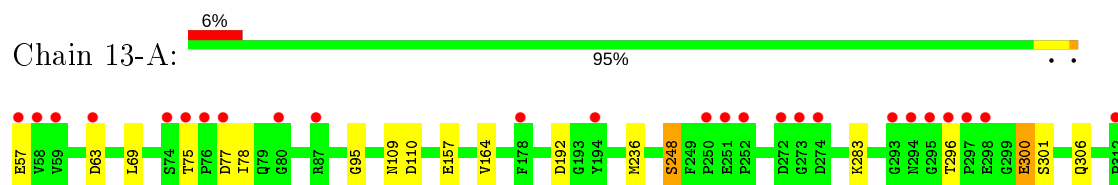
- Molecule 1: Putative secreted protein

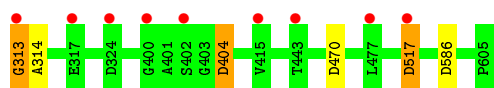


- Molecule 1: Putative secreted protein

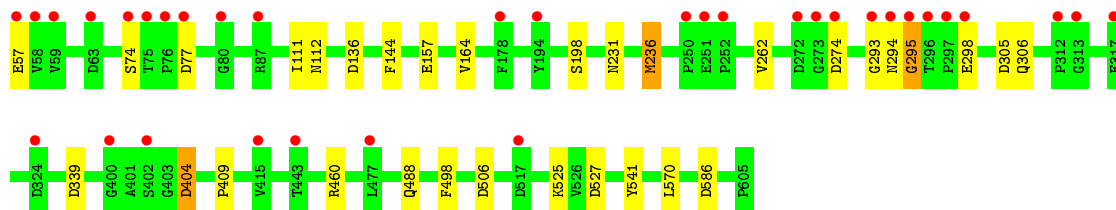


- Molecule 1: Putative secreted protein

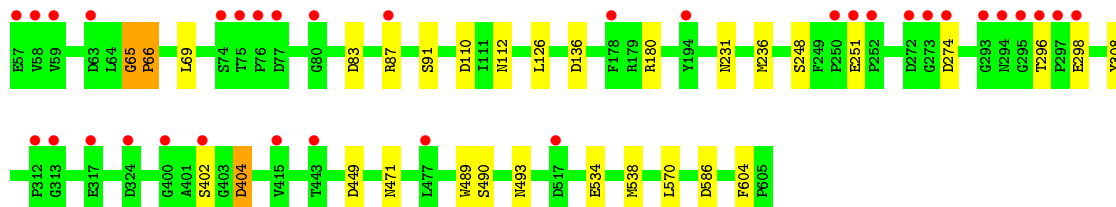




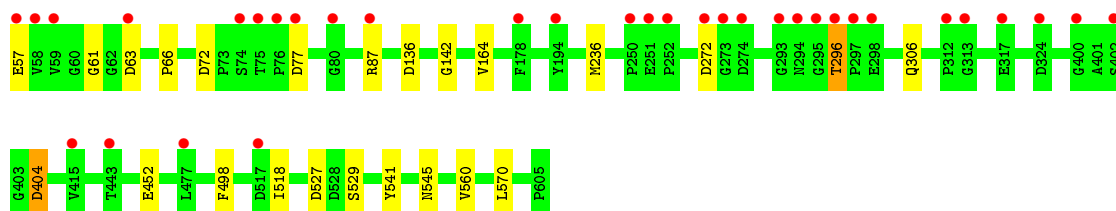
- Molecule 1: Putative secreted protein



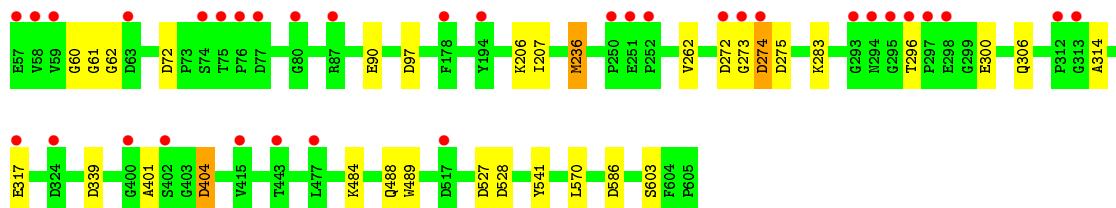
- Molecule 1: Putative secreted protein



- Molecule 1: Putative secreted protein

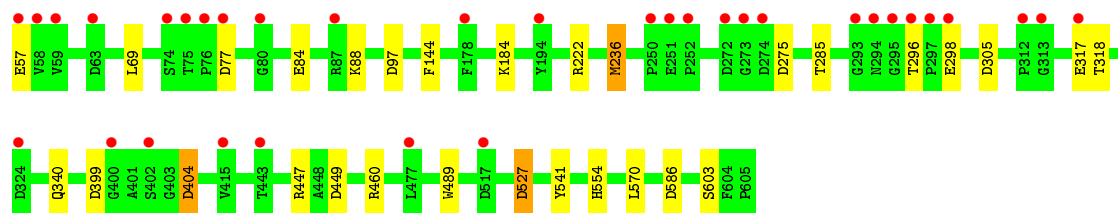


- Molecule 1: Putative secreted protein

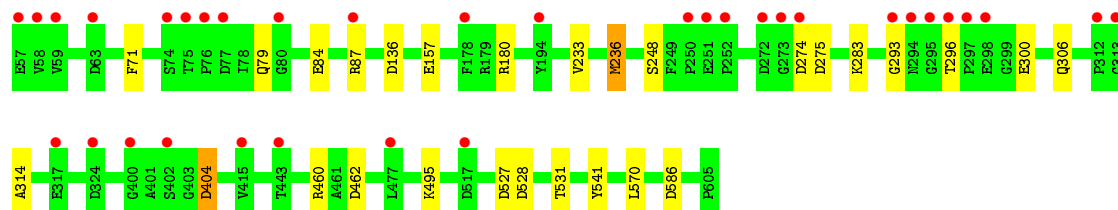


- Molecule 1: Putative secreted protein

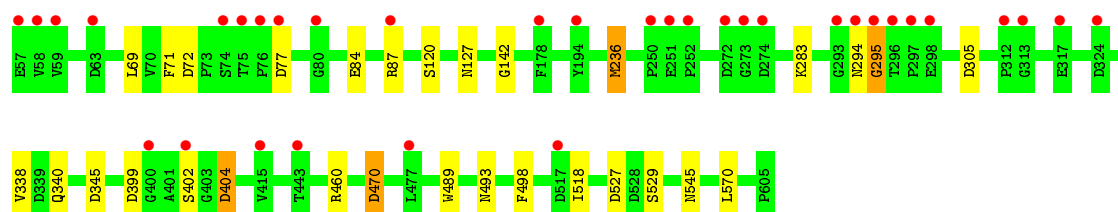




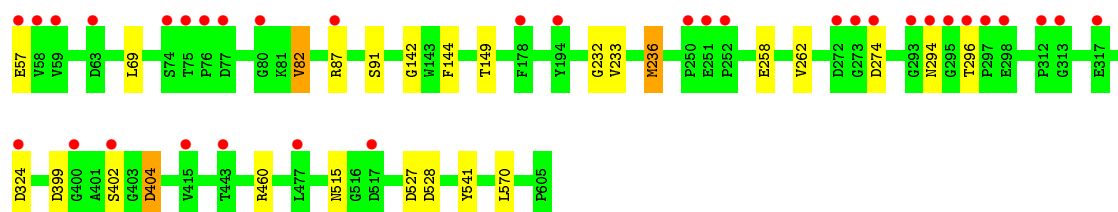
- Molecule 1: Putative secreted protein



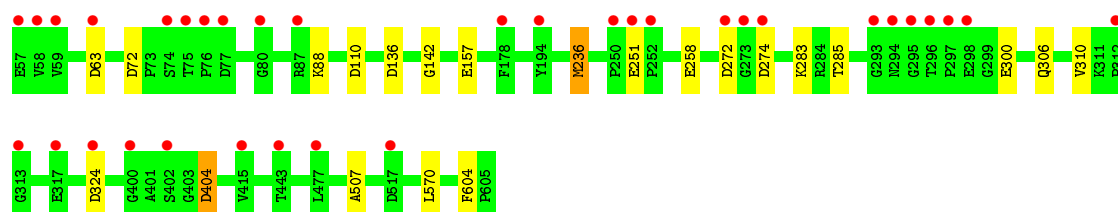
- Molecule 1: Putative secreted protein



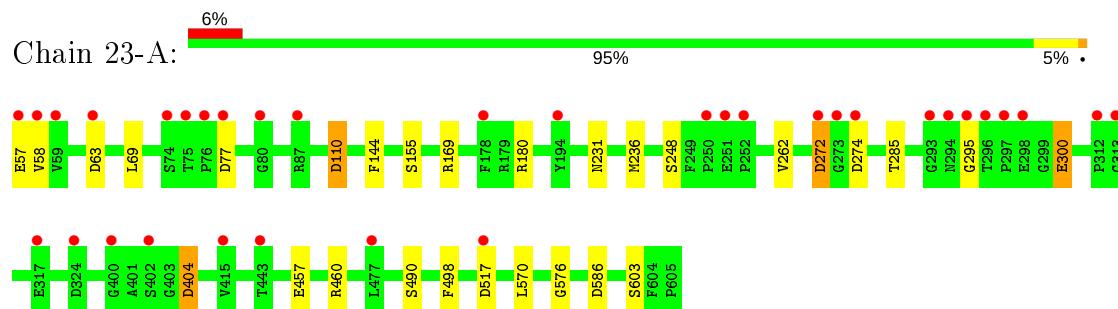
- Molecule 1: Putative secreted protein



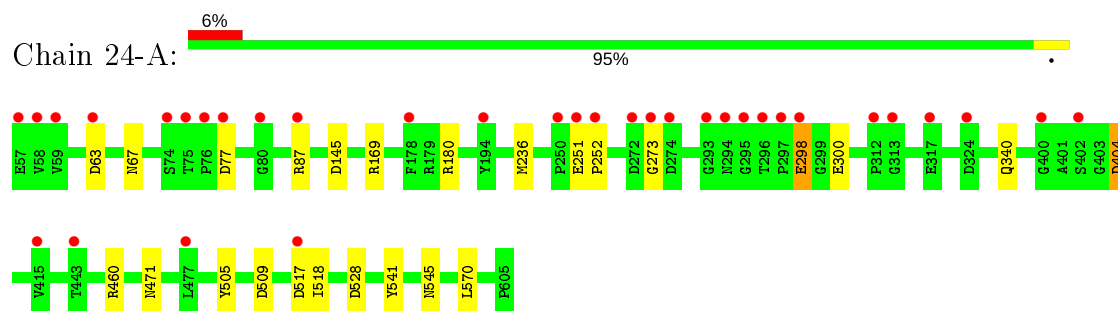
- Molecule 1: Putative secreted protein



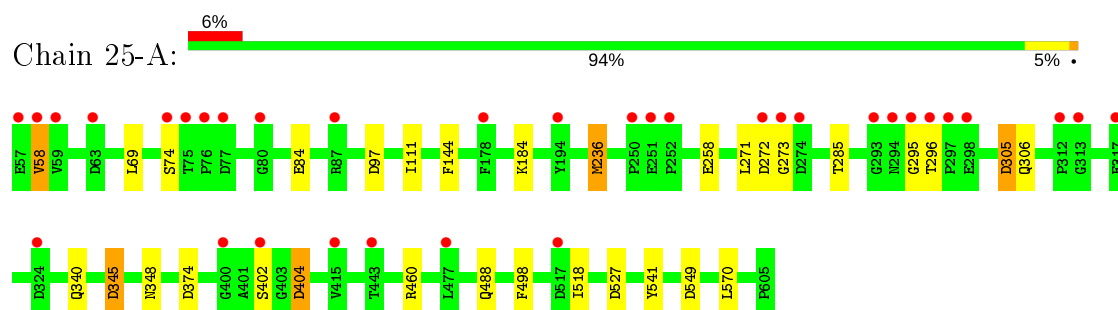
- Molecule 1: Putative secreted protein



- Molecule 1: Putative secreted protein



- Molecule 1: Putative secreted protein



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



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




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

Chain 3-B:  100%


BGC1
BGC2
BGC3
BGC4

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

Chain 4-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 5-B:  100%


BGC1
BGC2
BGC3
BGC4

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

Chain 6-B:  100%


BGC1
BGC2
BGC3
BGC4

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

Chain 7-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 8-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 9-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 10-B:  100%

B6C1
B6C2
B6C3
B6C4

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

Chain 11-B:  100%

B6C1
B6C2
B6C3
B6C4

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

Chain 12-B:  100%

B6C1
B6C2
B6C3
B6C4

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

Chain 13-B:  100%

B6C1
B6C2
B6C3
B6C4

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

Chain 14-B:  100%

B6C1
B6C2
B6C3
B6C4

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

Chain 15-B:  100%

B6C1
B6C2
B6C3
B6C4

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

Chain 16-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 17-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 18-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 19-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 20-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 21-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 22-B:  100%


BGC1
BGC2
BGC3
BGC4

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

Chain 23-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 24-B:  100%

BGC1
BGC2
BGC3
BGC4

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

Chain 25-B:  100%

BGC1
BGC2
BGC3
BGC4

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.28Å 100.21Å 54.22Å 90.00° 99.46° 90.00°	Depositor
Resolution (Å)	28.33 – 1.90 28.33 – 1.90	Depositor EDS
% Data completeness (in resolution range)	93.2 (28.33-1.90) 90.3 (28.33-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.97 (at 1.91Å)	Xtriage
Refinement program	PHENIX (phenix.ensemble_refinement: 1.9_1692)	Depositor
R, R_{free}	0.113 , 0.151 0.143 , 0.178	Depositor DCC
R_{free} test set	1989 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	19.3	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 114.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	210296	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.90% 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: BGC, 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	1-A	0.85	5/4280 (0.1%)	0.93	11/5848 (0.2%)
1	2-A	0.85	4/4280 (0.1%)	0.91	7/5848 (0.1%)
1	3-A	0.81	1/4280 (0.0%)	0.91	7/5848 (0.1%)
1	4-A	0.83	3/4280 (0.1%)	0.91	7/5848 (0.1%)
1	5-A	0.80	2/4280 (0.0%)	0.89	7/5848 (0.1%)
1	6-A	0.86	5/4280 (0.1%)	0.91	7/5848 (0.1%)
1	7-A	0.83	4/4280 (0.1%)	0.91	7/5848 (0.1%)
1	8-A	0.83	4/4280 (0.1%)	0.86	4/5848 (0.1%)
1	9-A	0.83	5/4280 (0.1%)	0.88	4/5848 (0.1%)
1	10-A	0.85	4/4280 (0.1%)	0.91	6/5848 (0.1%)
1	11-A	0.80	2/4280 (0.0%)	0.90	6/5848 (0.1%)
1	12-A	0.83	2/4280 (0.0%)	0.92	8/5848 (0.1%)
1	13-A	0.83	2/4280 (0.0%)	0.92	7/5848 (0.1%)
1	14-A	0.84	6/4280 (0.1%)	0.92	6/5848 (0.1%)
1	15-A	0.81	1/4280 (0.0%)	0.92	5/5848 (0.1%)
1	16-A	0.82	2/4280 (0.0%)	0.90	5/5848 (0.1%)
1	17-A	0.82	4/4280 (0.1%)	0.91	8/5848 (0.1%)
1	18-A	0.82	4/4280 (0.1%)	0.94	13/5848 (0.2%)
1	19-A	0.87	4/4280 (0.1%)	0.94	9/5848 (0.2%)
1	20-A	0.85	5/4280 (0.1%)	0.92	9/5848 (0.2%)
1	21-A	0.81	2/4280 (0.0%)	0.89	6/5848 (0.1%)
1	22-A	0.82	3/4280 (0.1%)	0.89	3/5848 (0.1%)
1	23-A	0.80	1/4280 (0.0%)	0.90	12/5848 (0.2%)
1	24-A	0.83	3/4280 (0.1%)	0.97	15/5848 (0.3%)
1	25-A	0.86	4/4280 (0.1%)	0.96	13/5848 (0.2%)
All	All	0.83	82/107000 (0.1%)	0.91	192/146200 (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	3-A	0	1
1	14-A	0	1
1	17-A	0	1
1	19-A	0	1
1	24-A	0	1
All	All	0	5

The worst 5 of 82 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	1-A	236	MET	CB-CG	8.44	1.78	1.51
1	10-A	198	SER	CB-OG	-8.39	1.31	1.42
1	19-A	404	ASP	CB-CG	8.26	1.69	1.51
1	4-A	404	ASP	CB-CG	8.12	1.68	1.51
1	1-A	236	MET	CG-SD	8.07	2.02	1.81

The worst 5 of 192 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	4-A	236	MET	CG-SD-CE	-14.71	76.66	100.20
1	18-A	236	MET	CG-SD-CE	-14.68	76.71	100.20
1	1-A	236	MET	CG-SD-CE	-14.32	77.28	100.20
1	24-A	87	ARG	NE-CZ-NH2	-13.75	113.42	120.30
1	14-A	236	MET	CG-SD-CE	-12.86	79.63	100.20

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	14-A	295	GLY	Peptide
1	17-A	62	GLY	Peptide
1	19-A	293	GLY	Peptide
1	24-A	251	GLU	Peptide
1	3-A	295	GLY	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1-A	4167	3907	3919	0	0
1	2-A	4167	3907	3919	0	0
1	3-A	4167	3907	3919	0	0
1	4-A	4167	3907	3919	0	0
1	5-A	4167	3907	3919	0	0
1	6-A	4167	3907	3919	0	0
1	7-A	4167	3907	3919	0	0
1	8-A	4167	3907	3919	0	0
1	9-A	4167	3907	3919	0	0
1	10-A	4167	3907	3919	0	0
1	11-A	4167	3907	3919	0	0
1	12-A	4167	3907	3919	0	0
1	13-A	4167	3907	3919	0	0
1	14-A	4167	3907	3919	0	0
1	15-A	4167	3907	3919	0	0
1	16-A	4167	3907	3919	0	0
1	17-A	4167	3907	3919	0	0
1	18-A	4167	3907	3919	0	0
1	19-A	4167	3907	3919	0	0
1	20-A	4167	3907	3919	0	0
1	21-A	4167	3907	3919	0	0
1	22-A	4167	3907	3919	0	0
1	23-A	4167	3907	3919	0	0
1	24-A	4167	3907	3919	0	0
1	25-A	4167	3907	3919	0	0
2	1-B	45	0	37	0	0
2	2-B	45	0	39	0	0
2	3-B	45	0	39	0	0
2	4-B	45	0	39	0	0
2	5-B	45	0	36	0	0
2	6-B	45	0	39	0	0
2	7-B	45	0	38	0	0
2	8-B	45	0	38	0	0
2	9-B	45	0	38	0	0
2	10-B	45	0	38	0	0
2	11-B	45	0	39	0	0
2	12-B	45	0	39	0	0
2	13-B	45	0	39	0	0
2	14-B	45	0	39	0	0
2	15-B	45	0	38	0	0
2	16-B	45	0	39	0	0
2	17-B	45	0	39	0	0
2	18-B	45	0	38	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	19-B	45	0	39	0	0
2	20-B	45	0	38	0	0
2	21-B	45	0	37	0	0
2	22-B	45	0	39	0	0
2	23-B	45	0	38	0	0
2	24-B	45	0	38	0	0
2	25-B	45	0	39	0	0
3	1-A	11	0	10	0	0
3	2-A	11	0	10	0	0
3	3-A	11	0	10	0	0
3	4-A	11	0	10	0	0
3	5-A	11	0	10	0	0
3	6-A	11	0	10	0	0
3	7-A	11	0	10	0	0
3	8-A	11	0	10	0	0
3	9-A	11	0	10	0	0
3	10-A	11	0	10	0	0
3	11-A	11	0	10	0	0
3	12-A	11	0	10	0	0
3	13-A	11	0	10	0	0
3	14-A	11	0	10	0	0
3	15-A	11	0	10	0	0
3	16-A	11	0	10	0	0
3	17-A	11	0	10	0	0
3	18-A	11	0	10	0	0
3	19-A	11	0	10	0	0
3	20-A	11	0	10	0	0
3	21-A	11	0	10	0	0
3	22-A	11	0	10	0	0
3	23-A	11	0	10	0	0
3	24-A	11	0	10	0	0
3	25-A	11	0	10	0	0
4	1-A	12	18	18	0	0
4	2-A	12	18	18	0	0
4	3-A	12	18	18	0	0
4	4-A	12	18	18	0	0
4	5-A	12	18	18	0	0
4	6-A	12	18	18	0	0
4	7-A	12	18	18	0	0
4	8-A	12	18	18	0	0
4	9-A	12	18	18	0	0
4	10-A	12	18	18	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	11-A	12	18	18	0	0
4	12-A	12	18	18	0	0
4	13-A	12	18	18	0	0
4	14-A	12	18	18	0	0
4	15-A	12	18	18	0	0
4	16-A	12	18	18	0	0
4	17-A	12	18	18	0	0
4	18-A	12	18	18	0	0
4	19-A	12	18	18	0	0
4	20-A	12	18	18	0	0
4	21-A	12	18	18	0	0
4	22-A	12	18	18	0	0
4	23-A	12	18	18	0	0
4	24-A	12	18	18	0	0
4	25-A	12	18	18	0	0
5	1-A	253	0	0	0	0
5	2-A	248	0	0	0	0
5	3-A	259	0	0	0	0
5	4-A	249	0	0	0	0
5	5-A	246	0	0	0	0
5	6-A	245	0	0	0	0
5	7-A	258	0	0	0	0
5	8-A	264	0	0	0	0
5	9-A	251	0	0	0	0
5	10-A	248	0	0	0	0
5	11-A	264	0	0	0	0
5	12-A	246	0	0	0	0
5	13-A	247	0	0	0	0
5	14-A	242	0	0	0	0
5	15-A	262	0	0	0	0
5	16-A	252	0	0	0	0
5	17-A	243	0	0	0	0
5	18-A	255	0	0	0	0
5	19-A	274	0	0	0	0
5	20-A	242	0	0	0	0
5	21-A	240	0	0	0	0
5	22-A	250	0	0	0	0
5	23-A	259	0	0	0	0
5	24-A	250	0	0	0	0
5	25-A	249	0	0	0	0
All	All	112171	98125	99634	0	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). Clashscore could not be calculated for this entry.

There are no clashes within the asymmetric unit.

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	1-A	547/549 (100%)	508 (93%)	34 (6%)	5 (1%)	17	7
1	2-A	547/549 (100%)	503 (92%)	40 (7%)	4 (1%)	22	12
1	3-A	547/549 (100%)	508 (93%)	34 (6%)	5 (1%)	17	7
1	4-A	547/549 (100%)	512 (94%)	30 (6%)	5 (1%)	17	7
1	5-A	547/549 (100%)	510 (93%)	32 (6%)	5 (1%)	17	7
1	6-A	547/549 (100%)	503 (92%)	40 (7%)	4 (1%)	22	12
1	7-A	547/549 (100%)	506 (92%)	38 (7%)	3 (0%)	29	18
1	8-A	547/549 (100%)	509 (93%)	35 (6%)	3 (0%)	29	18
1	9-A	547/549 (100%)	502 (92%)	40 (7%)	5 (1%)	17	7
1	10-A	547/549 (100%)	501 (92%)	45 (8%)	1 (0%)	47	38
1	11-A	547/549 (100%)	511 (93%)	30 (6%)	6 (1%)	14	5
1	12-A	547/549 (100%)	511 (93%)	32 (6%)	4 (1%)	22	12
1	13-A	547/549 (100%)	505 (92%)	37 (7%)	5 (1%)	17	7
1	14-A	547/549 (100%)	509 (93%)	34 (6%)	4 (1%)	22	12
1	15-A	547/549 (100%)	511 (93%)	29 (5%)	7 (1%)	12	4
1	16-A	547/549 (100%)	499 (91%)	43 (8%)	5 (1%)	17	7
1	17-A	547/549 (100%)	502 (92%)	38 (7%)	7 (1%)	12	4
1	18-A	547/549 (100%)	510 (93%)	35 (6%)	2 (0%)	34	24
1	19-A	547/549 (100%)	503 (92%)	40 (7%)	4 (1%)	22	12
1	20-A	547/549 (100%)	505 (92%)	35 (6%)	7 (1%)	12	4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	21-A	547/549 (100%)	513 (94%)	31 (6%)	3 (0%)	29	18
1	22-A	547/549 (100%)	507 (93%)	38 (7%)	2 (0%)	34	24
1	23-A	547/549 (100%)	506 (92%)	34 (6%)	7 (1%)	12	4
1	24-A	547/549 (100%)	501 (92%)	43 (8%)	3 (0%)	29	18
1	25-A	547/549 (100%)	501 (92%)	43 (8%)	3 (0%)	29	18
All	All	13675/13725 (100%)	12656 (92%)	910 (7%)	109 (1%)	19	9

5 of 109 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	2-A	509	ASP
1	3-A	314	ALA
1	4-A	297	PRO
1	6-A	77	ASP
1	6-A	319	ILE

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	1-A	436/436 (100%)	419 (96%)	17 (4%)	32	23
1	2-A	436/436 (100%)	417 (96%)	19 (4%)	28	19
1	3-A	436/436 (100%)	416 (95%)	20 (5%)	27	17
1	4-A	436/436 (100%)	417 (96%)	19 (4%)	28	19
1	5-A	436/436 (100%)	413 (95%)	23 (5%)	22	13
1	6-A	436/436 (100%)	413 (95%)	23 (5%)	22	13
1	7-A	436/436 (100%)	418 (96%)	18 (4%)	30	21
1	8-A	436/436 (100%)	417 (96%)	19 (4%)	28	19
1	9-A	436/436 (100%)	422 (97%)	14 (3%)	39	30
1	10-A	436/436 (100%)	419 (96%)	17 (4%)	32	23
1	11-A	436/436 (100%)	415 (95%)	21 (5%)	25	16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	12-A	436/436 (100%)	416 (95%)	20 (5%)	27	17
1	13-A	436/436 (100%)	420 (96%)	16 (4%)	34	25
1	14-A	436/436 (100%)	412 (94%)	24 (6%)	21	12
1	15-A	436/436 (100%)	415 (95%)	21 (5%)	25	16
1	16-A	436/436 (100%)	420 (96%)	16 (4%)	34	25
1	17-A	436/436 (100%)	419 (96%)	17 (4%)	32	23
1	18-A	436/436 (100%)	414 (95%)	22 (5%)	24	15
1	19-A	436/436 (100%)	419 (96%)	17 (4%)	32	23
1	20-A	436/436 (100%)	419 (96%)	17 (4%)	32	23
1	21-A	436/436 (100%)	416 (95%)	20 (5%)	27	17
1	22-A	436/436 (100%)	420 (96%)	16 (4%)	34	25
1	23-A	436/436 (100%)	419 (96%)	17 (4%)	32	23
1	24-A	436/436 (100%)	424 (97%)	12 (3%)	43	36
1	25-A	436/436 (100%)	411 (94%)	25 (6%)	20	11
All	All	10900/10900 (100%)	10430 (96%)	470 (4%)	29	19

5 of 470 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	12-A	149	THR
1	14-A	488	GLN
1	24-A	67	ASN
1	12-A	283	LYS
1	13-A	248	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 68 such sidechains are listed below:

Mol	Chain	Res	Type
1	11-A	553	GLN
1	13-A	600	GLN
1	24-A	483	ASN
1	12-A	127	ASN
1	12-A	515	ASN

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 ⓘ

100 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	1-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	1-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	1-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	1-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	10-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	10-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	10-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	10-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	11-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	11-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	11-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	11-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	12-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	12-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	12-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	12-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	13-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	13-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BGC	13-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	13-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	14-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	14-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	14-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	14-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	15-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	15-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	15-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	15-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	16-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	16-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	16-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	16-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	17-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
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2	BGC	17-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	17-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	18-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	18-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	18-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	18-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	19-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	19-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	19-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	19-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	2-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	2-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	2-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
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2	BGC	20-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
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Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
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2	BGC	21-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	22-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	22-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
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2	BGC	24-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	24-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	24-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	24-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	25-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	25-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	25-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	25-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	3-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	3-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	3-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
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2	BGC	4-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	4-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	5-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
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2	BGC	5-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	5-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	6-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	6-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	6-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	6-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BGC	7-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	7-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	7-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	7-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	8-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	8-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	8-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	8-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)
2	BGC	9-B	1	2	12,12,12	2.49	4 (33%)	17,17,17	3.27	7 (41%)
2	BGC	9-B	2	2	11,11,12	2.76	4 (36%)	15,15,17	1.76	2 (13%)
2	BGC	9-B	3	2	11,11,12	3.07	5 (45%)	15,15,17	2.04	6 (40%)
2	BGC	9-B	4	2	11,11,12	2.86	5 (45%)	15,15,17	3.06	6 (40%)

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	1-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	1-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	1-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	1-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	10-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	10-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	10-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	10-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	11-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	11-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	11-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	11-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	12-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	12-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	12-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	12-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	13-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	13-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	13-B	3	2	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	13-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	14-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	14-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	14-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	14-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	15-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	15-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	15-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	15-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	16-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	16-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	16-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	16-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	17-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	17-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	17-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	17-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	18-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	18-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	18-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	18-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	19-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	19-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	19-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	19-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	2-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	2-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	2-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	2-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	20-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	20-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	20-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	20-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	21-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	21-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	21-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	21-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	22-B	1	2	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	22-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	22-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	22-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	23-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	23-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	23-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	23-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	24-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	24-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	24-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	24-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	25-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	25-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	25-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	25-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	3-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	3-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	3-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	3-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	4-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	4-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	4-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	4-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	5-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	5-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	5-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	5-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	6-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	6-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	6-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	6-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	7-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	7-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	7-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	7-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	8-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	8-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	8-B	3	2	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	8-B	4	2	-	2/2/19/22	0/1/1/1
2	BGC	9-B	1	2	-	0/2/22/22	0/1/1/1
2	BGC	9-B	2	2	-	0/2/19/22	0/1/1/1
2	BGC	9-B	3	2	-	2/2/19/22	0/1/1/1
2	BGC	9-B	4	2	-	2/2/19/22	0/1/1/1

The worst 5 of 450 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	1-B	3	BGC	C2-C3	-6.20	1.43	1.52
2	4-B	3	BGC	C2-C3	-6.20	1.43	1.52
2	25-B	3	BGC	C2-C3	-6.20	1.43	1.52
2	7-B	3	BGC	C2-C3	-6.20	1.43	1.52
2	9-B	3	BGC	C2-C3	-6.20	1.43	1.52

The worst 5 of 525 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	10-B	4	BGC	C1-O5-C5	8.51	123.72	112.19
2	24-B	4	BGC	C1-O5-C5	8.51	123.72	112.19
2	23-B	4	BGC	C1-O5-C5	8.51	123.72	112.19
2	9-B	4	BGC	C1-O5-C5	8.51	123.72	112.19
2	2-B	4	BGC	C1-O5-C5	8.51	123.72	112.19

There are no chirality outliers.

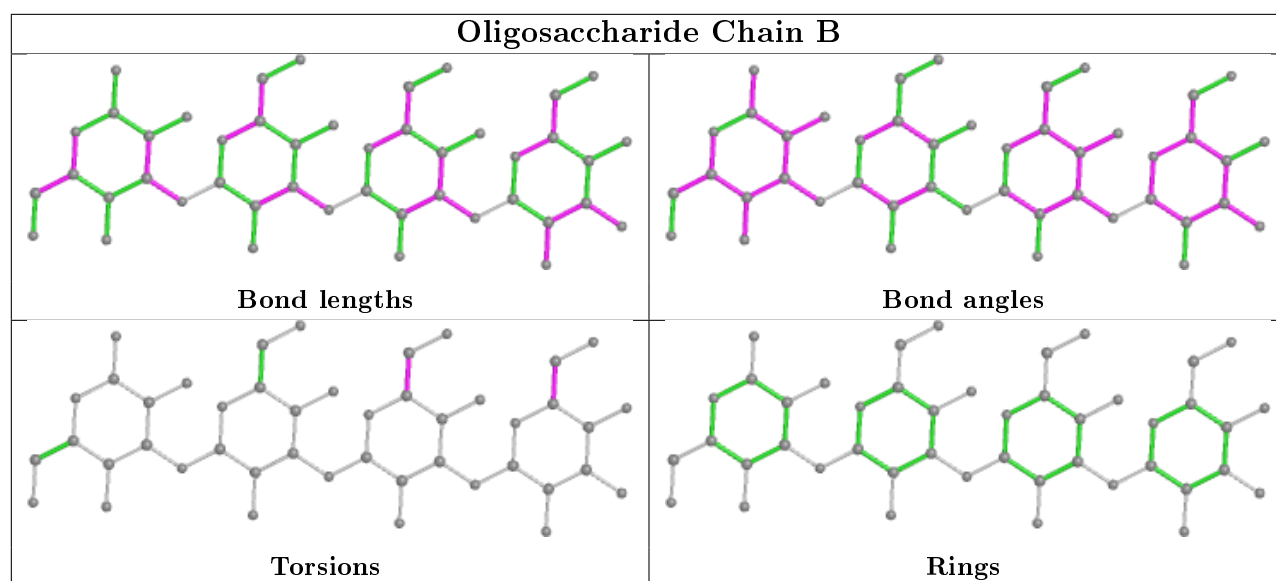
5 of 100 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	10-B	4	BGC	O5-C5-C6-O6
2	24-B	4	BGC	O5-C5-C6-O6
2	23-B	4	BGC	O5-C5-C6-O6
2	9-B	4	BGC	O5-C5-C6-O6
2	2-B	4	BGC	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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



5.6 Ligand geometry [i](#)

100 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	20-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	14-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	3-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	25-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	2-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	1-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	2-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	8-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	3-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	13-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
3	BGC	20-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	2-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
3	BGC	9-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
3	BGC	6-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	9-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	12-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	16-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
3	BGC	23-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	19-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	1-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	12-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	1-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	4-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
3	BGC	21-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	13-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	7-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	23-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	18-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
3	BGC	19-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	4-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	20-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
3	BGC	5-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	11-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	3-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	19-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
3	BGC	12-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
3	BGC	8-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	15-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	13-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	6-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	21-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
3	BGC	18-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	5-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	10-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	22-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	9-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	24-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	15-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
3	BGC	3-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	17-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	9-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	22-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	8-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	7-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	16-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	23-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	18-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	1-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	22-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
3	BGC	4-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
3	BGC	10-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	22-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	16-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	17-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	6-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	19-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	11-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	24-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	24-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	15-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	7-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	12-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
3	BGC	24-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	15-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	8-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	23-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
3	BGC	16-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	17-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	14-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	10-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	11-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	18-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	14-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	11-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	20-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	5-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	7-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
3	BGC	14-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)
4	EDO	13-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	4-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	10-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	21-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	2-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	25-A	702	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	5-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	25-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	17-A	704	-	3,3,3	0.41	0	2,2,2	0.17	0
4	EDO	21-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	6-A	703	-	3,3,3	0.38	0	2,2,2	0.24	0
3	BGC	25-A	701	-	11,11,12	0.74	0	15,15,17	1.27	1 (6%)

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	20-A	703	-	-	0/1/1/1	-
4	EDO	14-A	704	-	-	0/1/1/1	-
4	EDO	3-A	704	-	-	0/1/1/1	-
4	EDO	25-A	703	-	-	0/1/1/1	-
3	BGC	2-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	1-A	703	-	-	0/1/1/1	-
4	EDO	2-A	703	-	-	0/1/1/1	-
4	EDO	8-A	704	-	-	0/1/1/1	-
4	EDO	3-A	702	-	-	0/1/1/1	-
4	EDO	13-A	702	-	-	0/1/1/1	-
3	BGC	20-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	2-A	704	-	-	0/1/1/1	-
3	BGC	9-A	701	-	-	2/2/19/22	0/1/1/1
3	BGC	6-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	9-A	702	-	-	0/1/1/1	-
4	EDO	12-A	704	-	-	0/1/1/1	-
4	EDO	16-A	702	-	-	0/1/1/1	-
3	BGC	23-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	19-A	703	-	-	0/1/1/1	-
4	EDO	1-A	702	-	-	0/1/1/1	-
4	EDO	12-A	703	-	-	0/1/1/1	-
4	EDO	1-A	704	-	-	0/1/1/1	-
4	EDO	4-A	704	-	-	0/1/1/1	-
3	BGC	21-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	13-A	703	-	-	0/1/1/1	-
4	EDO	7-A	702	-	-	0/1/1/1	-
4	EDO	23-A	703	-	-	0/1/1/1	-
4	EDO	18-A	702	-	-	0/1/1/1	-
3	BGC	19-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	4-A	703	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	20-A	702	-	-	0/1/1/1	-
3	BGC	5-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	11-A	703	-	-	0/1/1/1	-
4	EDO	3-A	703	-	-	0/1/1/1	-
4	EDO	19-A	704	-	-	0/1/1/1	-
3	BGC	12-A	701	-	-	2/2/19/22	0/1/1/1
3	BGC	8-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	15-A	703	-	-	0/1/1/1	-
3	BGC	13-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	6-A	704	-	-	0/1/1/1	-
4	EDO	21-A	704	-	-	0/1/1/1	-
3	BGC	18-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	5-A	702	-	-	0/1/1/1	-
4	EDO	10-A	703	-	-	0/1/1/1	-
4	EDO	22-A	702	-	-	0/1/1/1	-
4	EDO	9-A	704	-	-	0/1/1/1	-
4	EDO	24-A	703	-	-	0/1/1/1	-
3	BGC	15-A	701	-	-	2/2/19/22	0/1/1/1
3	BGC	3-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	17-A	703	-	-	0/1/1/1	-
4	EDO	9-A	703	-	-	0/1/1/1	-
3	BGC	22-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	8-A	703	-	-	0/1/1/1	-
4	EDO	7-A	704	-	-	0/1/1/1	-
4	EDO	16-A	704	-	-	0/1/1/1	-
4	EDO	23-A	704	-	-	0/1/1/1	-
4	EDO	18-A	703	-	-	0/1/1/1	-
3	BGC	1-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	22-A	704	-	-	0/1/1/1	-
3	BGC	4-A	701	-	-	2/2/19/22	0/1/1/1
3	BGC	10-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	22-A	703	-	-	0/1/1/1	-
4	EDO	16-A	703	-	-	0/1/1/1	-
3	BGC	17-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	6-A	702	-	-	0/1/1/1	-
4	EDO	19-A	702	-	-	0/1/1/1	-
4	EDO	11-A	704	-	-	0/1/1/1	-
4	EDO	24-A	704	-	-	0/1/1/1	-
4	EDO	24-A	702	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	15-A	702	-	-	0/1/1/1	-
4	EDO	7-A	703	-	-	0/1/1/1	-
4	EDO	12-A	702	-	-	0/1/1/1	-
3	BGC	24-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	15-A	704	-	-	0/1/1/1	-
4	EDO	8-A	702	-	-	0/1/1/1	-
4	EDO	23-A	702	-	-	0/1/1/1	-
3	BGC	16-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	17-A	702	-	-	0/1/1/1	-
4	EDO	14-A	702	-	-	0/1/1/1	-
4	EDO	10-A	702	-	-	0/1/1/1	-
4	EDO	11-A	702	-	-	0/1/1/1	-
4	EDO	18-A	704	-	-	0/1/1/1	-
4	EDO	14-A	703	-	-	0/1/1/1	-
3	BGC	11-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	20-A	704	-	-	0/1/1/1	-
4	EDO	5-A	703	-	-	0/1/1/1	-
3	BGC	7-A	701	-	-	2/2/19/22	0/1/1/1
3	BGC	14-A	701	-	-	2/2/19/22	0/1/1/1
4	EDO	13-A	704	-	-	0/1/1/1	-
4	EDO	4-A	702	-	-	0/1/1/1	-
4	EDO	10-A	704	-	-	0/1/1/1	-
4	EDO	21-A	702	-	-	0/1/1/1	-
4	EDO	2-A	702	-	-	0/1/1/1	-
4	EDO	25-A	702	-	-	0/1/1/1	-
4	EDO	5-A	704	-	-	0/1/1/1	-
4	EDO	25-A	704	-	-	0/1/1/1	-
4	EDO	17-A	704	-	-	0/1/1/1	-
4	EDO	21-A	703	-	-	0/1/1/1	-
4	EDO	6-A	703	-	-	0/1/1/1	-
3	BGC	25-A	701	-	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	2-A	701	BGC	C6-C5-C4	-2.34	107.53	113.00
3	9-A	701	BGC	C6-C5-C4	-2.34	107.53	113.00
3	6-A	701	BGC	C6-C5-C4	-2.34	107.53	113.00
3	23-A	701	BGC	C6-C5-C4	-2.34	107.53	113.00
3	21-A	701	BGC	C6-C5-C4	-2.34	107.53	113.00

There are no chirality outliers.

5 of 50 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	2-A	701	BGC	O5-C5-C6-O6
3	9-A	701	BGC	O5-C5-C6-O6
3	6-A	701	BGC	O5-C5-C6-O6
3	23-A	701	BGC	O5-C5-C6-O6
3	21-A	701	BGC	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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	1-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	2-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	3-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	4-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	5-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	6-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	7-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	8-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	9-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	10-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	11-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	12-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	13-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	14-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	15-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	16-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	17-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	18-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	19-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	20-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	21-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	22-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	23-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)
1	24-A	549/549 (100%)	0.06	34 (6%)	20	23	17, 18, 20, 22	549 (100%)

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	25-A	549/549 (100%)	0.06	34 (6%) 20 23	17, 18, 20, 22	549 (100%)
All	All	13725/13725 (100%)	0.06	850 (6%) 23 23	17, 18, 20, 22	13725 (100%)

The worst 5 of 850 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	1-A	293	GLY	8.6
1	2-A	293	GLY	8.6
1	3-A	293	GLY	8.6
1	4-A	293	GLY	8.6
1	5-A	293	GLY	8.6

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(Å ²)	Q<0.9
2	BGC	22-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	18-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	7-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	3-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	23-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	24-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	19-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	17-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	9-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	24-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	16-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	20-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	11-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	13-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	21-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	18-B	1	12/12	0.96	0.15	18,21,31,34	12

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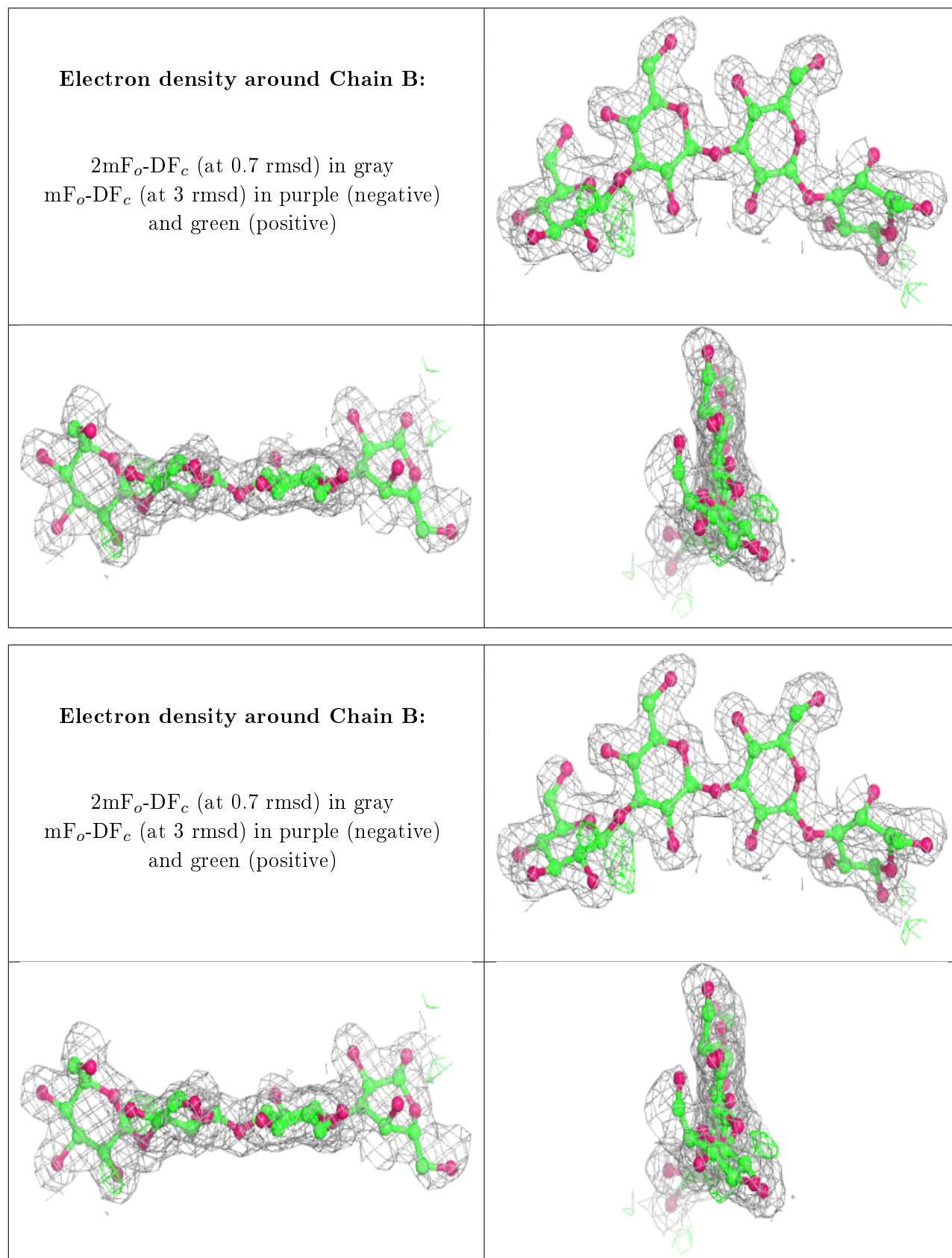
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BGC	20-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	12-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	25-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	10-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	6-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	9-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	8-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	14-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	10-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	16-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	5-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	2-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	1-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	7-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	5-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	1-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	2-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	15-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	22-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	8-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	3-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	6-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	13-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	4-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	17-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	25-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	11-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	4-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	21-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	12-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	23-B	1	12/12	0.96	0.15	18,21,31,34	12
2	BGC	19-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	14-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	15-B	2	11/12	0.96	0.06	16,20,21,23	11
2	BGC	19-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	24-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	6-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	3-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	17-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	16-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	12-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	1-B	3	11/12	0.97	0.07	15,17,22,23	11

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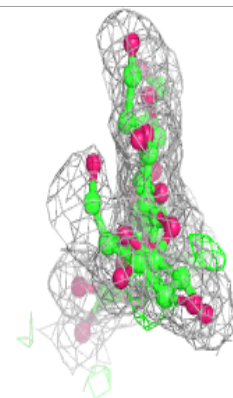
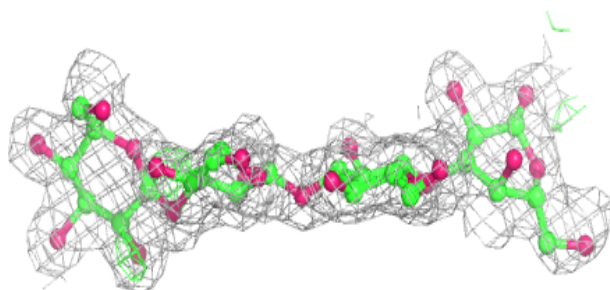
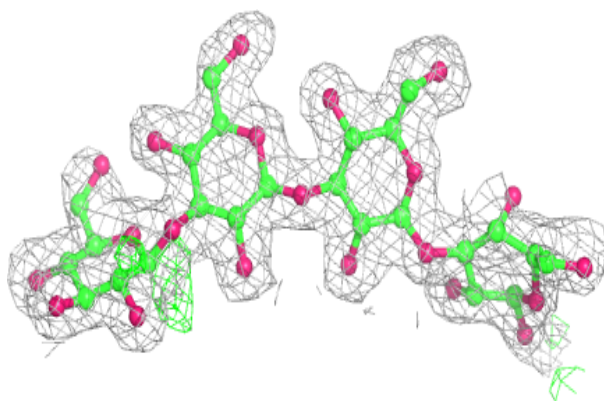
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BGC	1-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	15-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	16-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	9-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	17-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	5-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	25-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	15-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	21-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	14-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	4-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	5-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	12-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	4-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	11-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	8-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	10-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	6-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	20-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	20-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	13-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	22-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	2-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	7-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	23-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	21-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	18-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	22-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	14-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	18-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	11-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	7-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	9-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	25-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	8-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	24-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	2-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	3-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	19-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	23-B	3	11/12	0.97	0.07	15,17,22,23	11
2	BGC	10-B	4	11/12	0.97	0.10	14,17,18,18	11
2	BGC	13-B	3	11/12	0.97	0.07	15,17,22,23	11

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.

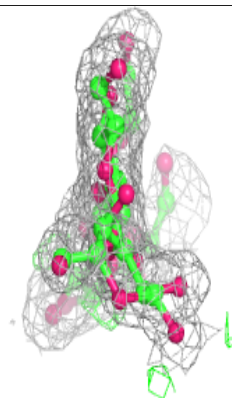
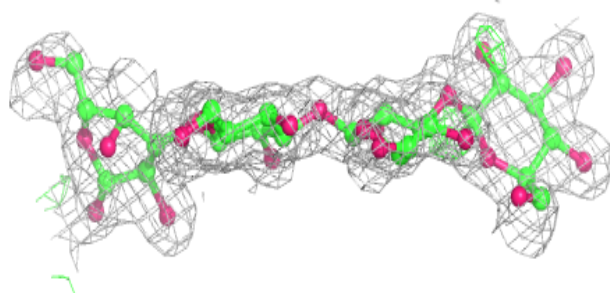
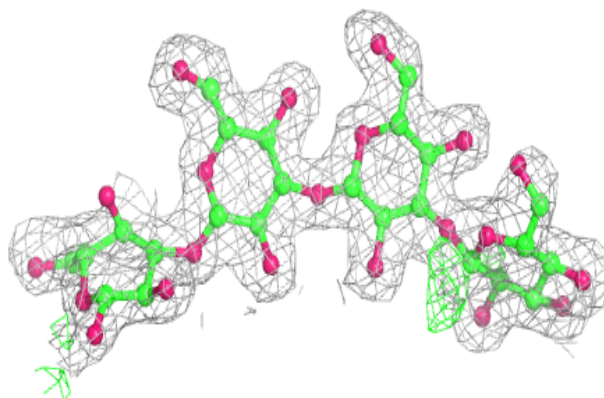


Electron density around Chain B:

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

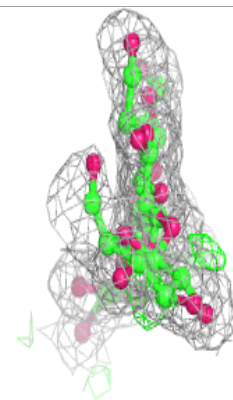
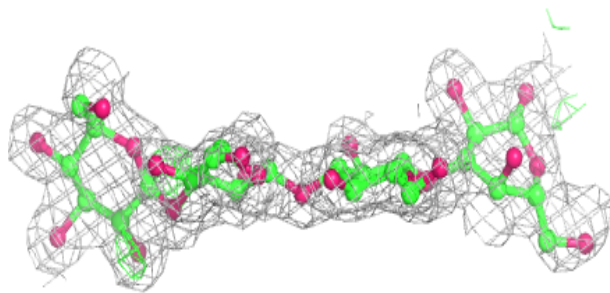
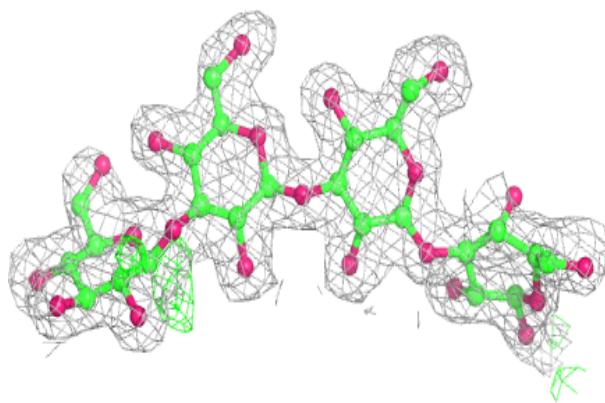
**Electron density around Chain B:**

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

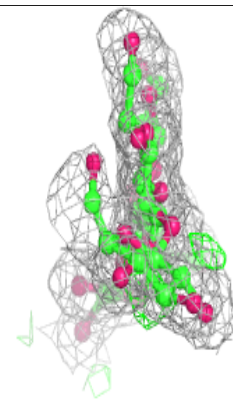
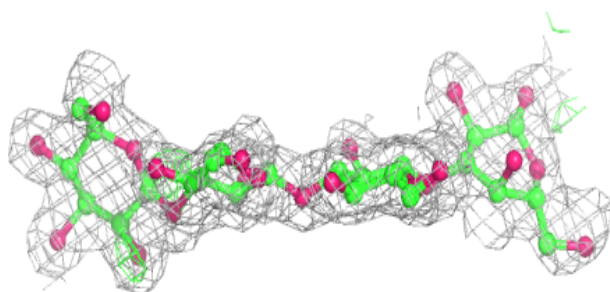
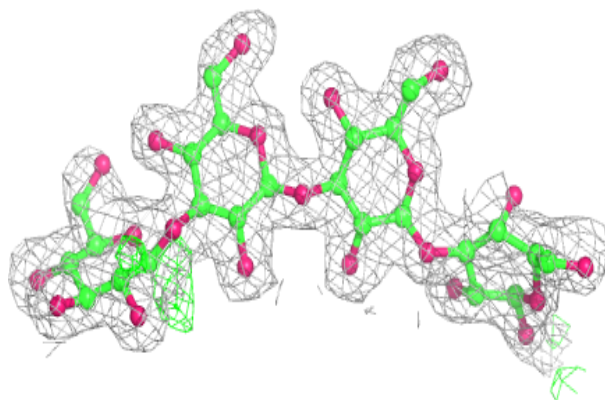


Electron density around Chain B:

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

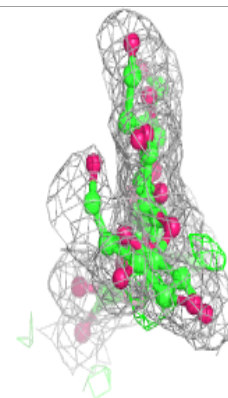
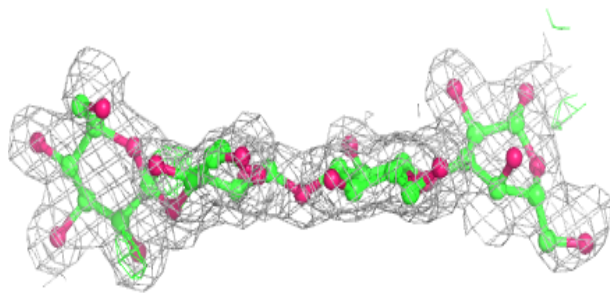
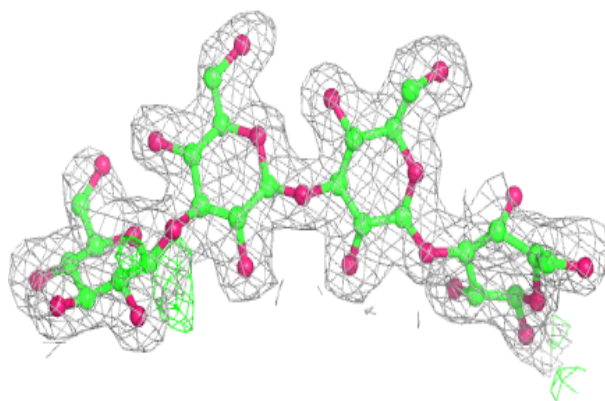
**Electron density around Chain B:**

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

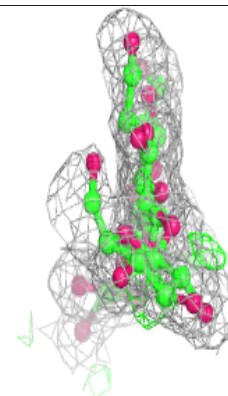
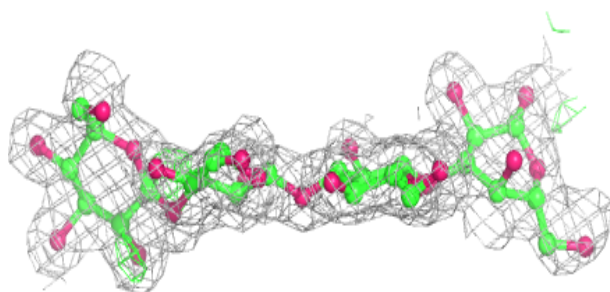
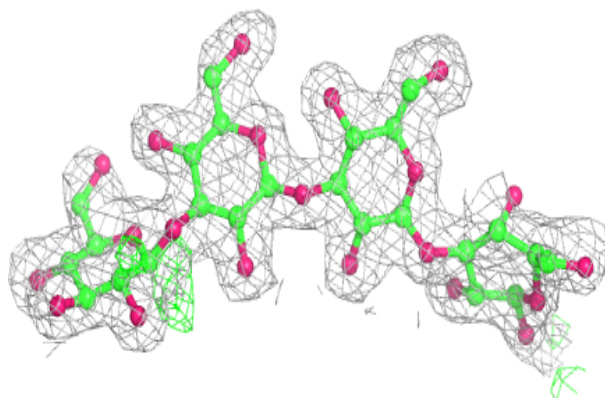


Electron density around Chain B:

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

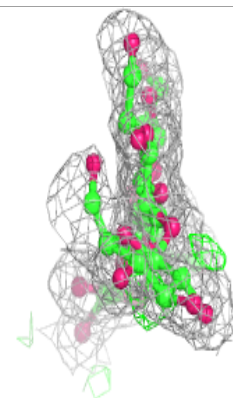
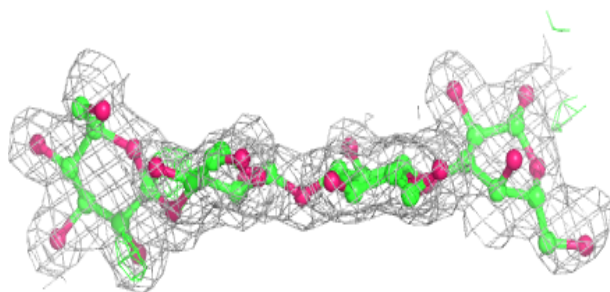
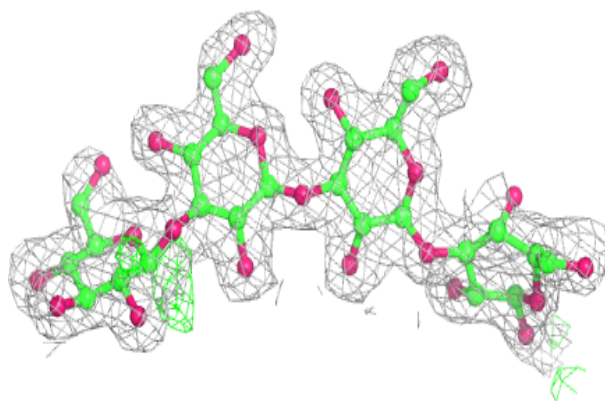
**Electron density around Chain B:**

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

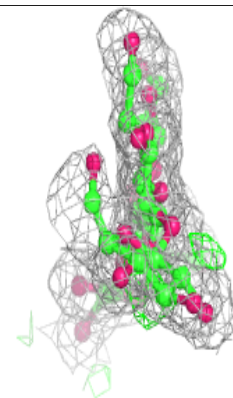
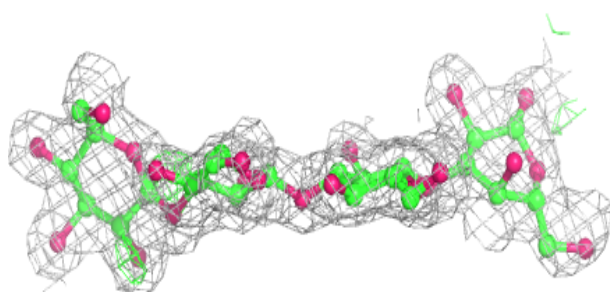
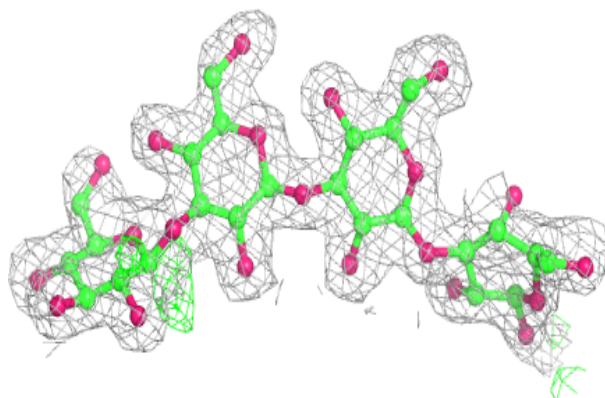


Electron density around Chain B:

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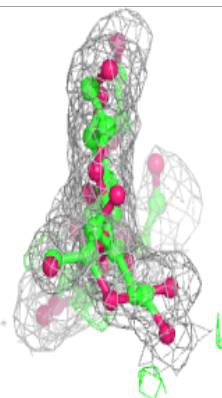
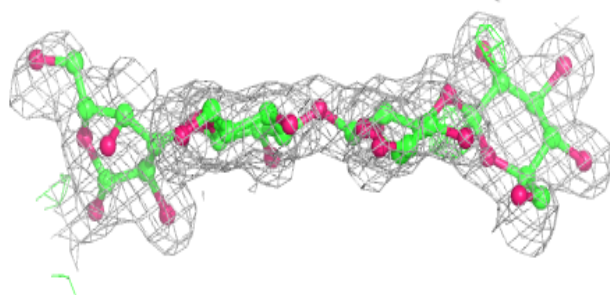
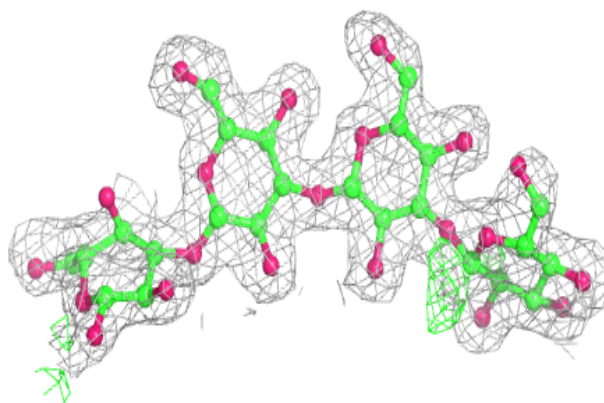
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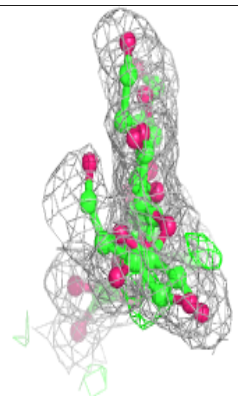
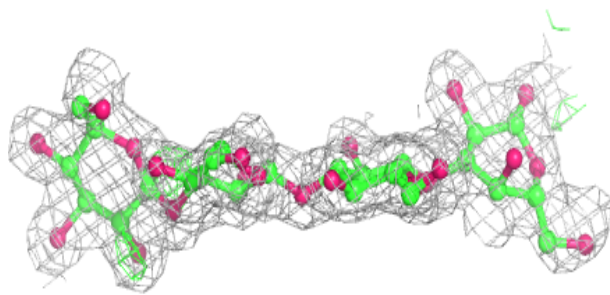
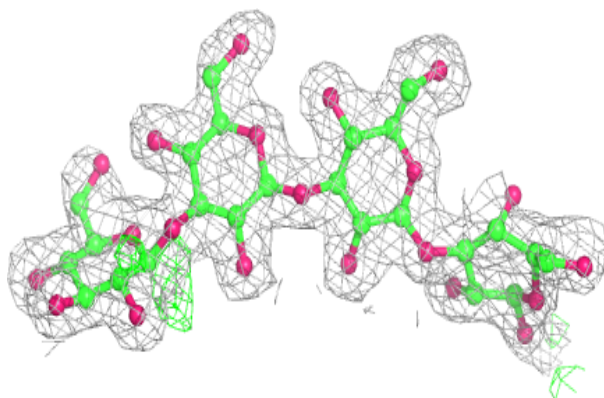


Electron density around Chain B:

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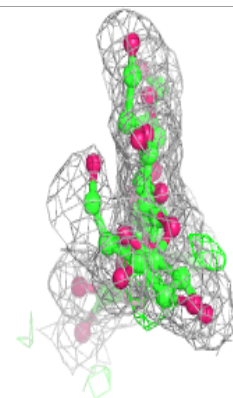
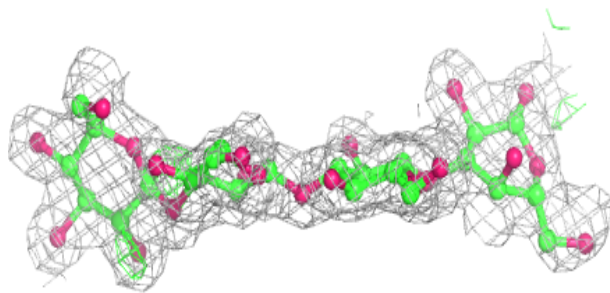
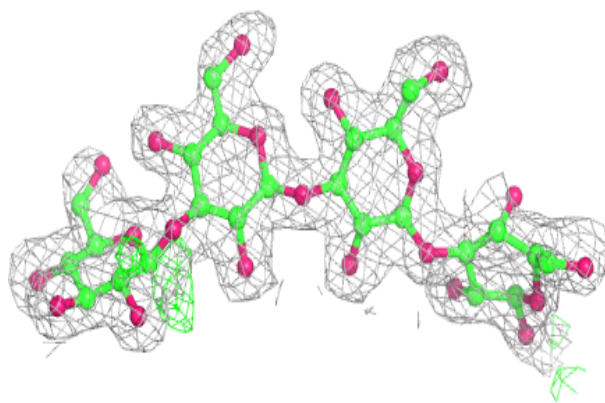
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 mF_o-DF_c (at 3 rmsd) in purple (negative)
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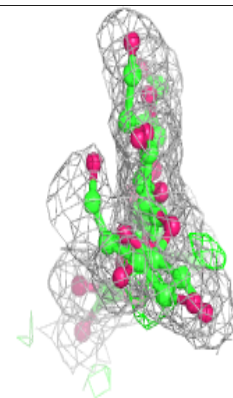
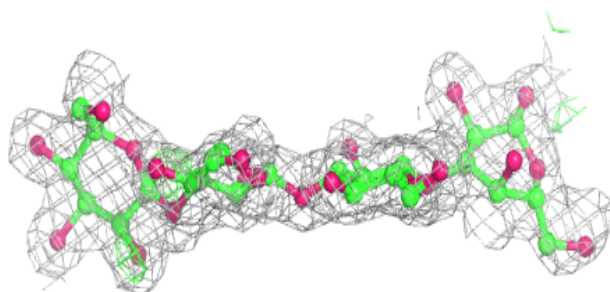
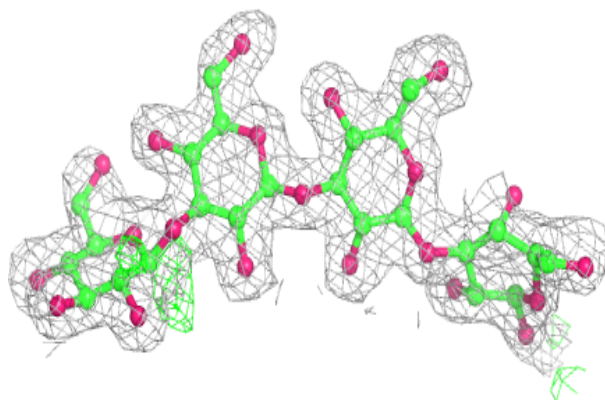


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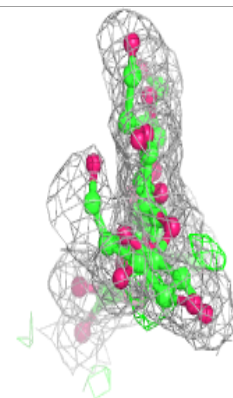
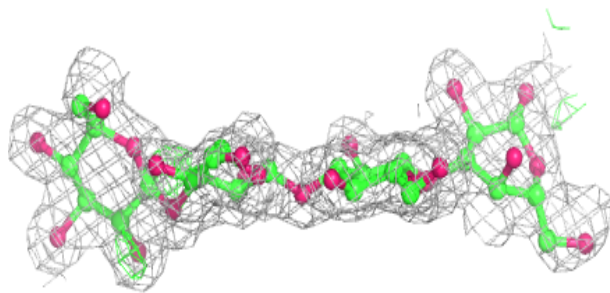
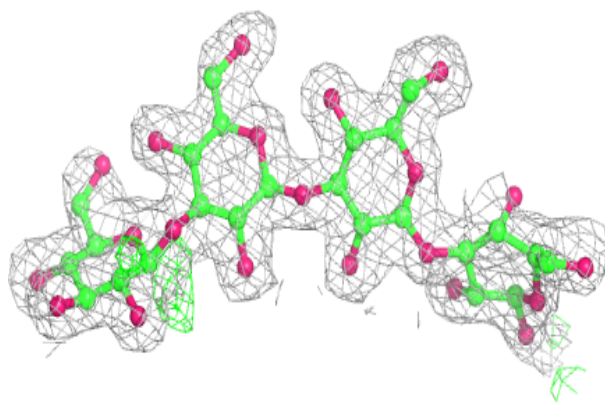
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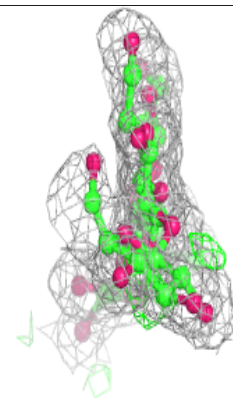
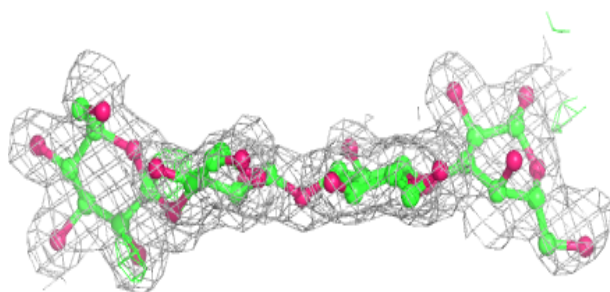
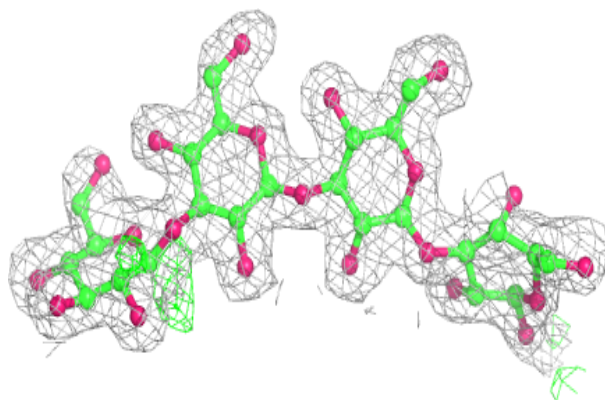


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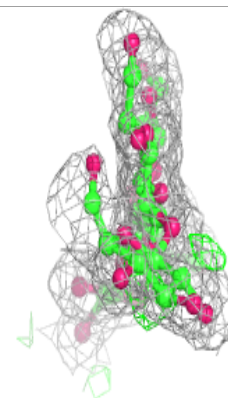
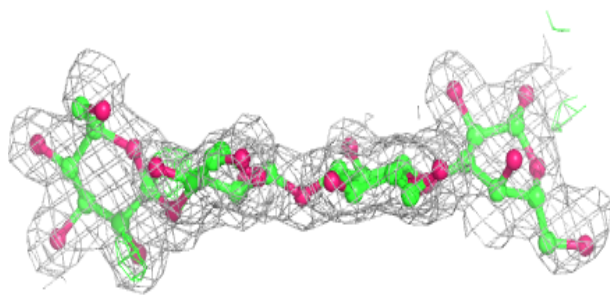
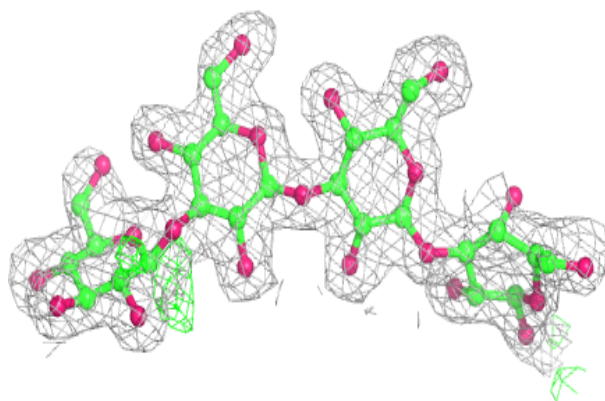
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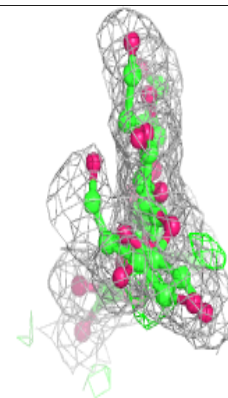
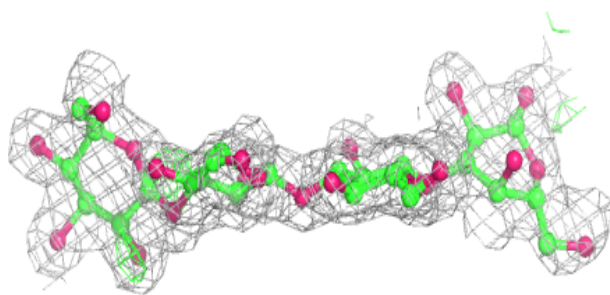
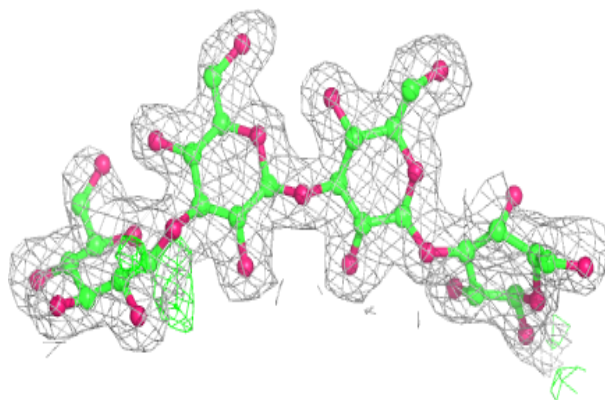


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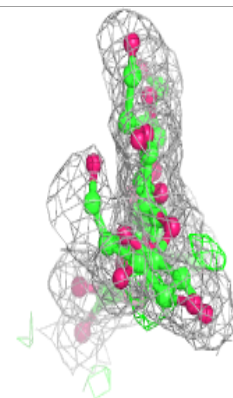
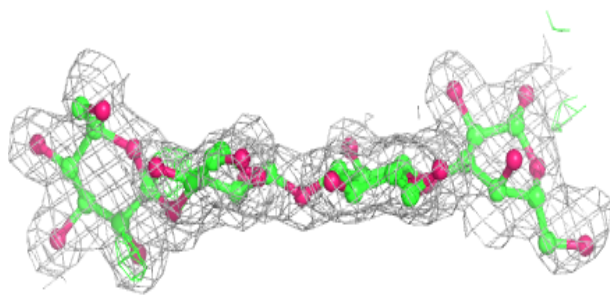
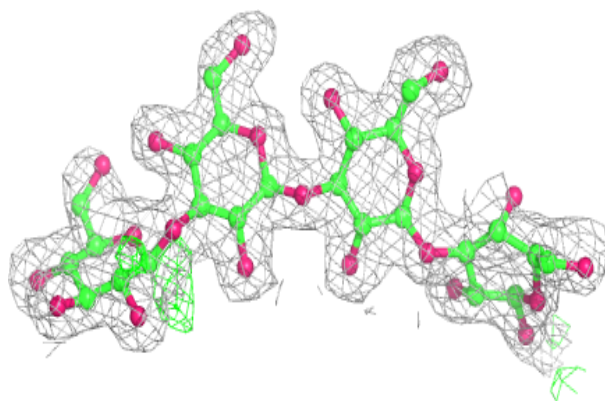
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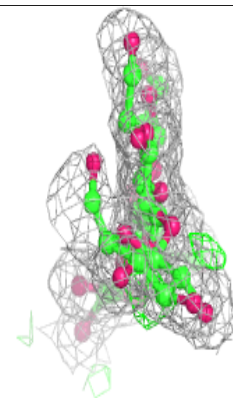
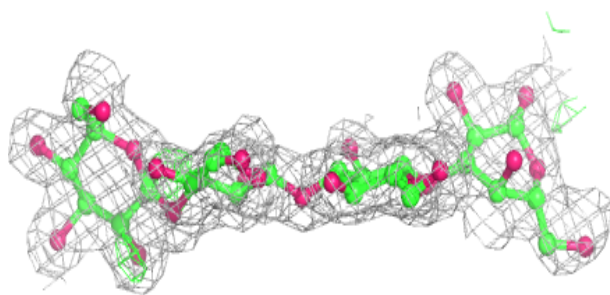
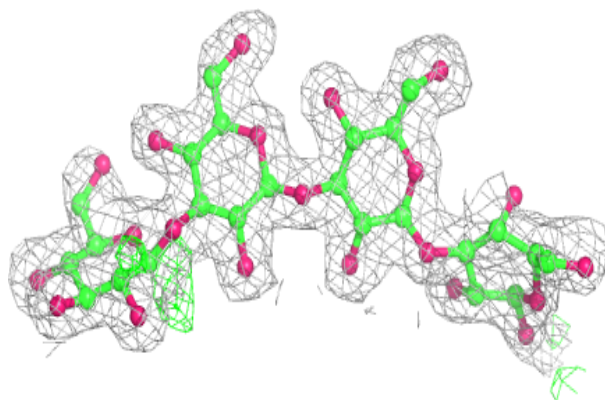


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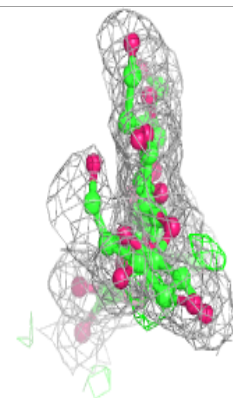
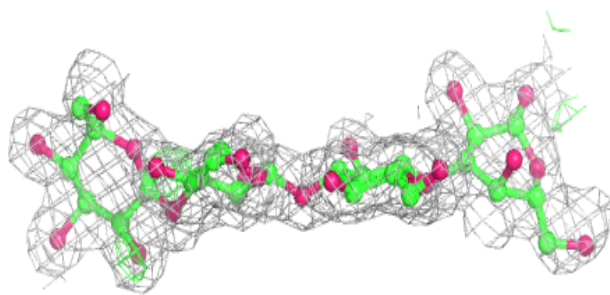
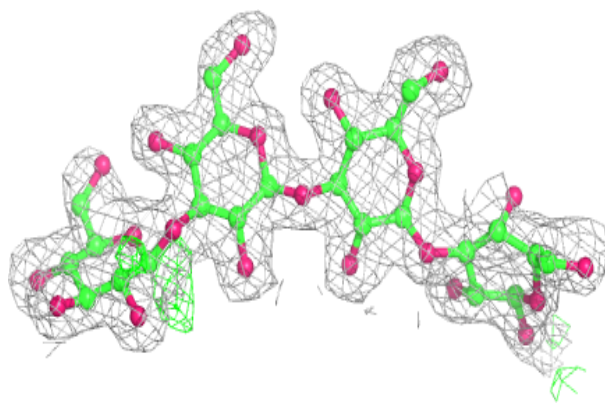
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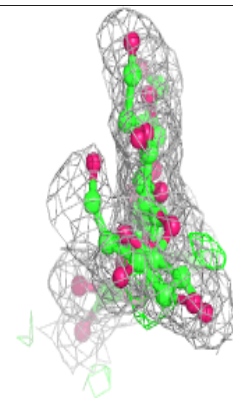
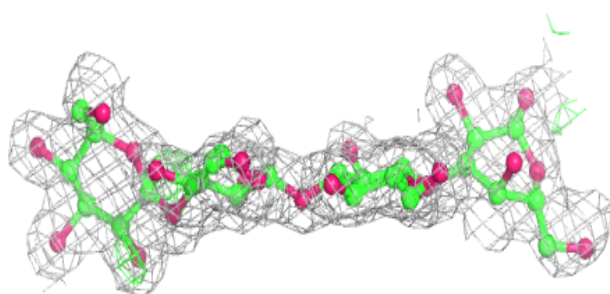
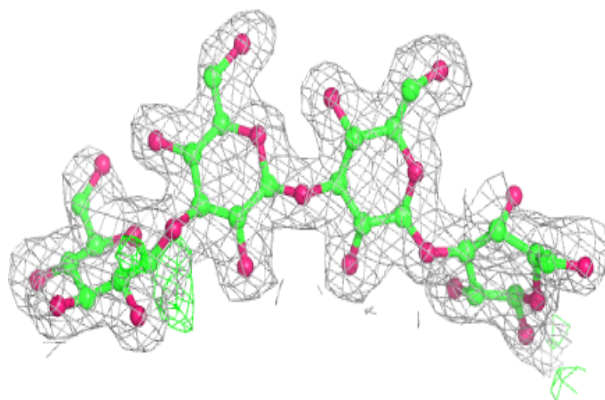


Electron density around Chain B:

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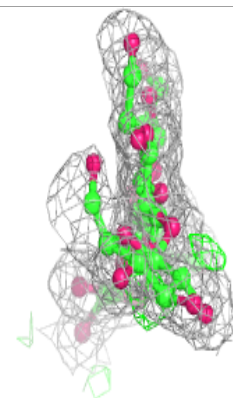
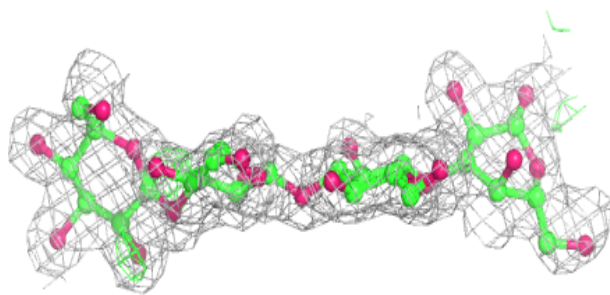
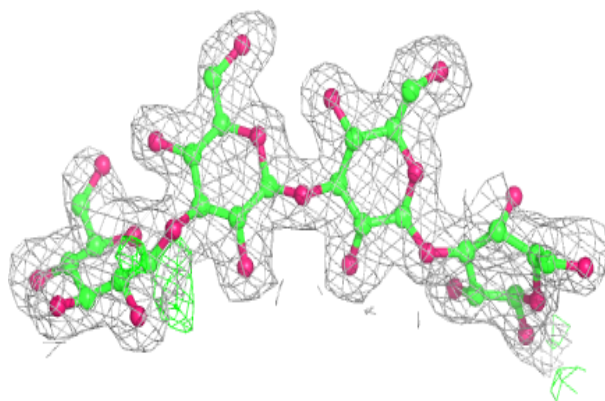
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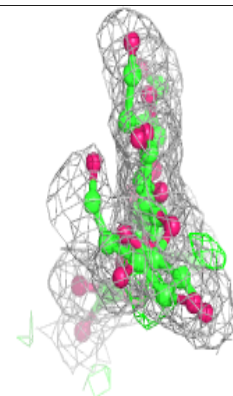
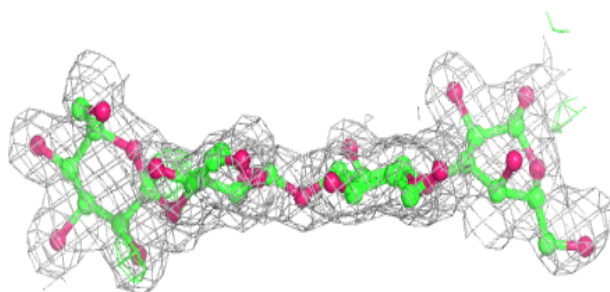
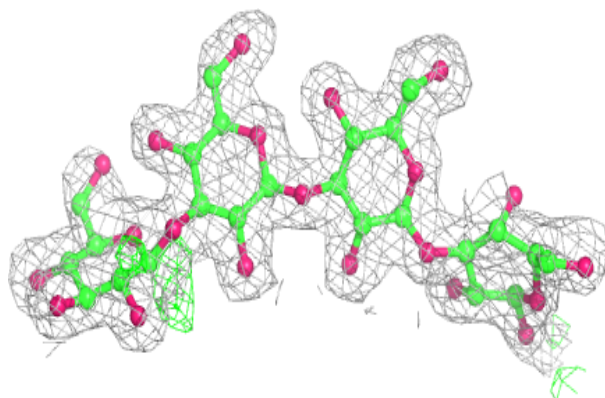


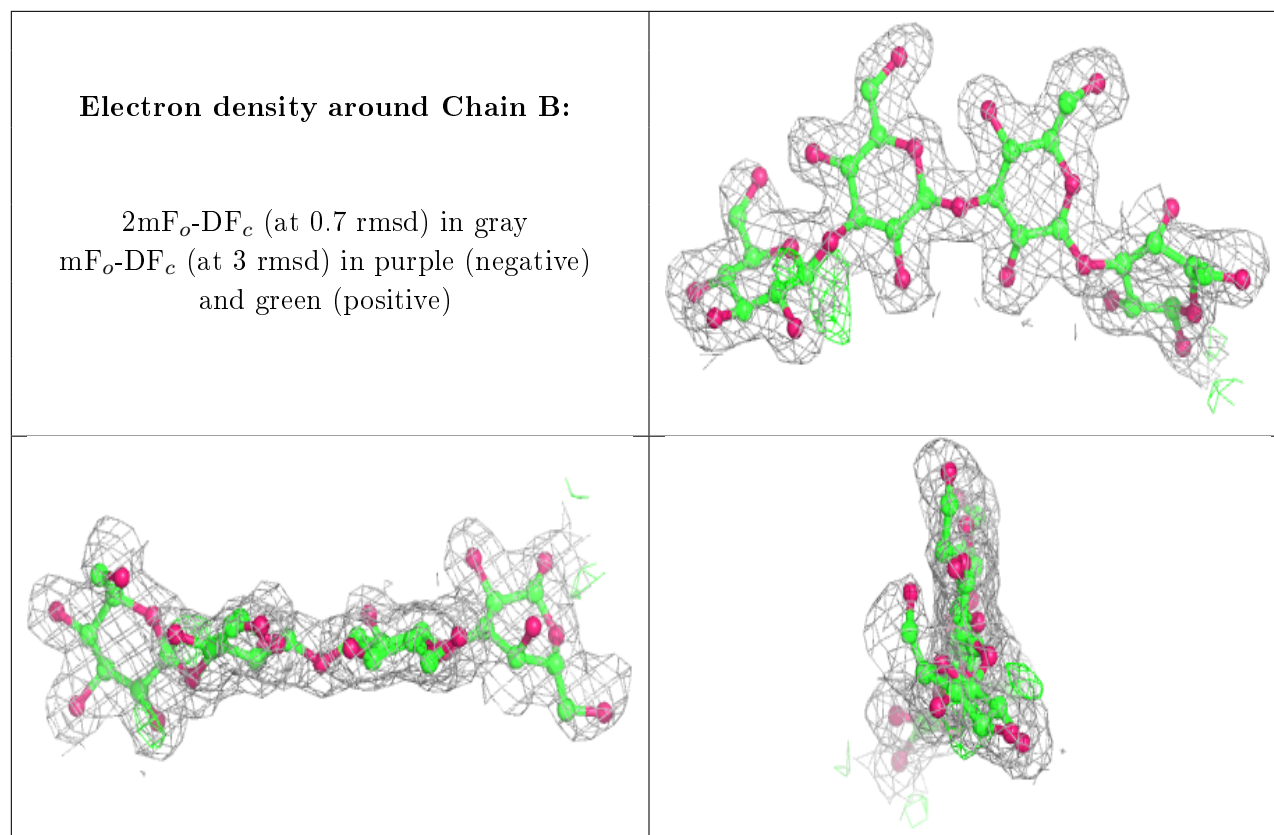
Electron density around Chain B:

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

**Electron density around Chain B:**

$2mF_o-DF_c$ (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
3	BGC	2-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	9-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	6-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	23-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	21-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	19-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	5-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	8-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	13-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	18-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	15-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	3-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	22-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	1-A	701	11/12	0.89	0.29	21,26,31,34	11

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	BGC	4-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	10-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	17-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	12-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	24-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	16-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	11-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	20-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	7-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	14-A	701	11/12	0.89	0.29	21,26,31,34	11
3	BGC	25-A	701	11/12	0.89	0.29	21,26,31,34	11
4	EDO	4-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	7-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	16-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	23-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	20-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	22-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	2-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	3-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	19-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	11-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	24-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	14-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	12-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	15-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	6-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	18-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	21-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	8-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	9-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	17-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	13-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	10-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	5-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	25-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	1-A	704	4/4	0.92	0.30	30,32,35,35	10
4	EDO	2-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	25-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	22-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	16-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	20-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	13-A	703	4/4	0.97	0.07	17,17,22,22	10

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	EDO	10-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	23-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	7-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	24-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	5-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	4-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	17-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	14-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	9-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	19-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	8-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	11-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	3-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	1-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	18-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	12-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	21-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	6-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	15-A	703	4/4	0.97	0.07	17,17,22,22	10
4	EDO	23-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	18-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	17-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	14-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	10-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	11-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	22-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	1-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	6-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	13-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	19-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	16-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	20-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	9-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	4-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	15-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	21-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	2-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	25-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	7-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	12-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	24-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	3-A	702	4/4	0.98	0.11	15,17,21,21	10

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	EDO	5-A	702	4/4	0.98	0.11	15,17,21,21	10
4	EDO	8-A	702	4/4	0.98	0.11	15,17,21,21	10

6.5 Other polymers [i](#)

There are no such residues in this entry.