Full wwPDB X-ray Structure Validation Report ⓘ

Dec 18, 2023 – 04:22 PM JST

PDB ID	:	8HUI
Title	:	Crystal structure of DFA I-forming Inulin Lyase from Streptomyces peucetius subsp. caesius ATCC 27952 in complex with GF4, DFA I, and fructose
Authors	:	Cheng, M.; Rao, Y.J.; Mu, W.M.
Deposited on	:	2022-12-24
Resolution	:	1.44 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:

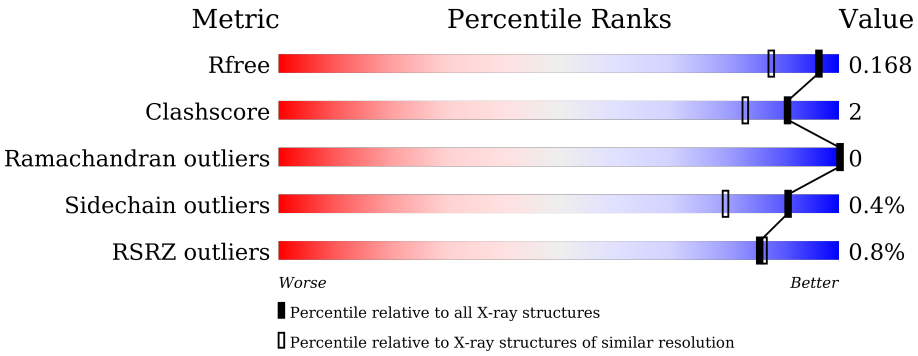
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.44 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}	130704	2021 (1.46-1.42)
Clashscore	141614	2086 (1.46-1.42)
Ramachandran outliers	138981	2047 (1.46-1.42)
Sidechain outliers	138945	2047 (1.46-1.42)
RSRZ outliers	127900	1993 (1.46-1.42)

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 <=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	400	95%
1	B	400	95%
1	C	400	96%
2	D	2	50% 50%
2	E	2	100%
3	F	5	100%

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Mol	Chain	Length	Quality of chain
3	G	5	100%
3	H	5	100%
3	I	5	100%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	FRU	G	5	-	-	-	X
7	PEG	C	501	-	-	X	-

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 19497 atoms, of which 9186 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.

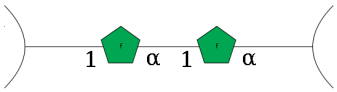
- Molecule 1 is a protein called Fructotransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	393	Total	C	H	N	O	S	0	20	0
			5947	1873	2950	527	591	6			
1	B	393	Total	C	H	N	O	S	0	20	0
			5945	1876	2945	526	592	6			
1	C	399	Total	C	H	N	O	S	0	25	0
			6071	1915	3003	546	601	6			

There are 18 discrepancies between the modelled and reference sequences:

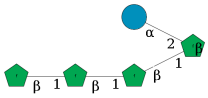
Chain	Residue	Modelled	Actual	Comment	Reference
A	395	HIS	-	expression tag	UNP A0A2D3U3Z1
A	396	HIS	-	expression tag	UNP A0A2D3U3Z1
A	397	HIS	-	expression tag	UNP A0A2D3U3Z1
A	398	HIS	-	expression tag	UNP A0A2D3U3Z1
A	399	HIS	-	expression tag	UNP A0A2D3U3Z1
A	400	HIS	-	expression tag	UNP A0A2D3U3Z1
B	395	HIS	-	expression tag	UNP A0A2D3U3Z1
B	396	HIS	-	expression tag	UNP A0A2D3U3Z1
B	397	HIS	-	expression tag	UNP A0A2D3U3Z1
B	398	HIS	-	expression tag	UNP A0A2D3U3Z1
B	399	HIS	-	expression tag	UNP A0A2D3U3Z1
B	400	HIS	-	expression tag	UNP A0A2D3U3Z1
C	395	HIS	-	expression tag	UNP A0A2D3U3Z1
C	396	HIS	-	expression tag	UNP A0A2D3U3Z1
C	397	HIS	-	expression tag	UNP A0A2D3U3Z1
C	398	HIS	-	expression tag	UNP A0A2D3U3Z1
C	399	HIS	-	expression tag	UNP A0A2D3U3Z1
C	400	HIS	-	expression tag	UNP A0A2D3U3Z1

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



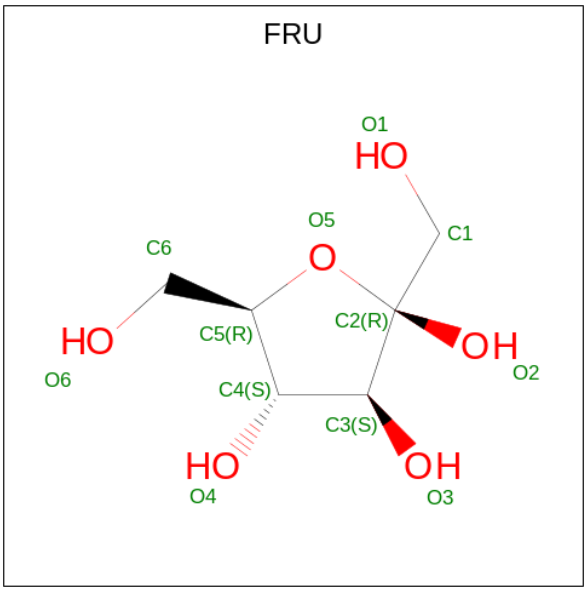
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	2	Total	C	H	O	0	0	0
			42	12	20	10			
2	E	2	Total	C	H	O	0	0	0
			42	12	20	10			

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



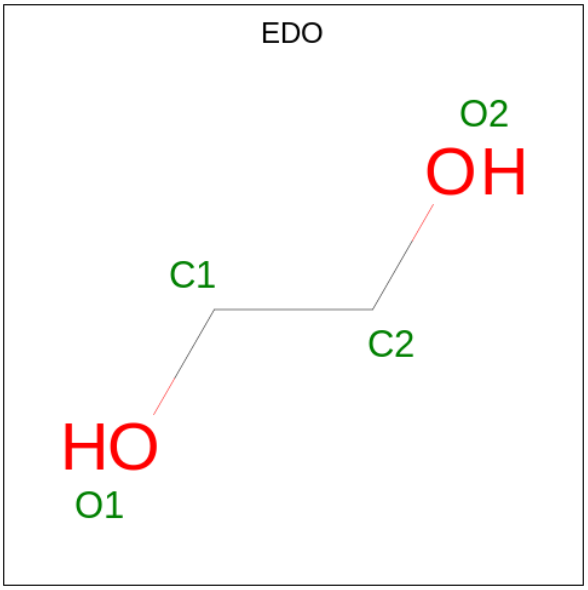
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	5	Total	C	H	O	0	0	0
			108	30	52	26			
3	G	5	Total	C	H	O	0	0	0
			108	30	52	26			
3	H	5	Total	C	H	O	0	0	0
			108	30	52	26			
3	I	5	Total	C	H	O	0	0	0
			108	30	52	26			

- Molecule 4 is beta-D-fructofuranose (three-letter code: FRU) (formula: $C_6H_{12}O_6$) (labeled as "Ligand of Interest" by depositor).



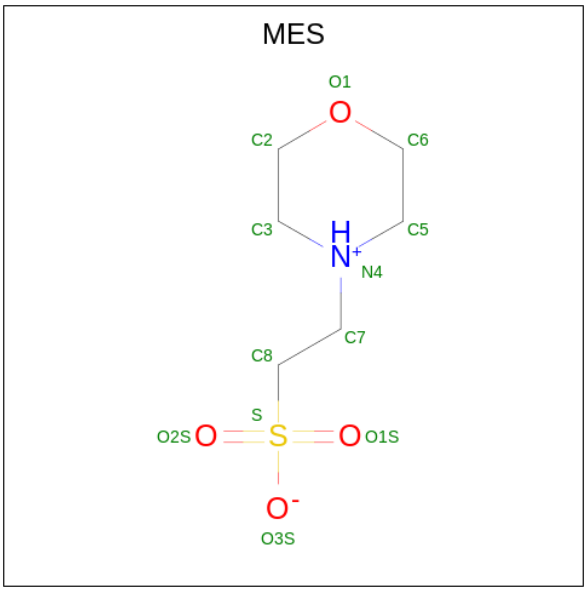
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			24	6	12	6		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



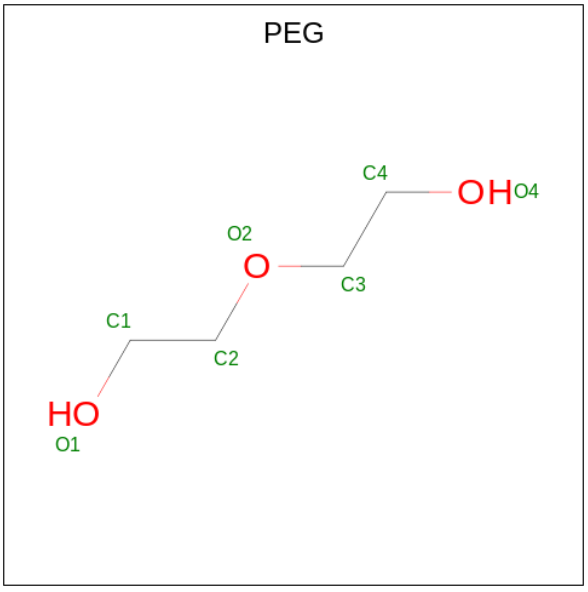
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 6 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	A	1	Total	C	H	N	O	S	0	0
			24	6	12	1	4	1		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 8 is water.

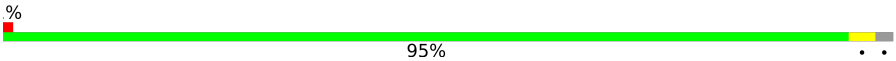
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	306	Total O 306 306	0	0
8	B	320	Total O 320 320	0	0
8	C	317	Total O 317 317	0	0

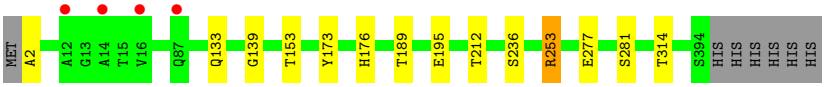
3 Residue-property plots (i)

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: Fructotransferase

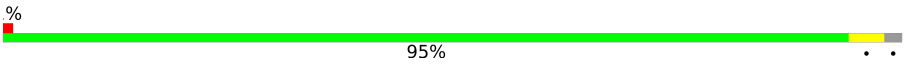
Chain A:

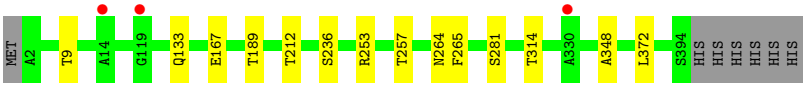




• Molecule 1: Fructotransferase

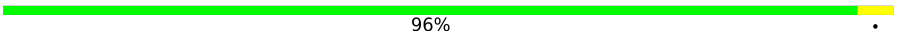
Chain B:





• Molecule 1: Fructotransferase

Chain C:

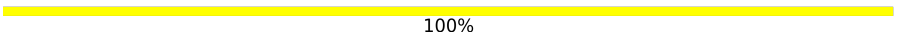


• Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-fructofuranose

Chain D:

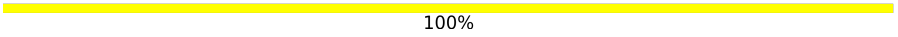
• Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-fructofuranose

Chain E:



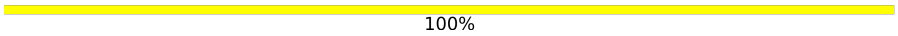
• Molecule 3: beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain F:



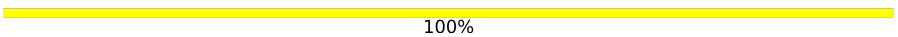
• Molecule 3: beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain G:



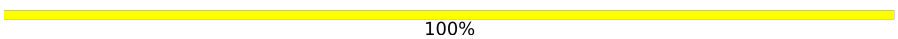
• Molecule 3: beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain H:



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

Chain I:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	97.76Å 105.60Å 131.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.84 – 1.44 45.84 – 1.44	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.84-1.44) 99.7 (45.84-1.44)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.85 (at 1.43Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.154 , 0.168 0.153 , 0.168	Depositor DCC
R_{free} test set	1991 reflections (0.81%)	wwPDB-VP
Wilson B-factor (Å^2)	11.2	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.43 , 45.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	19497	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: Z9N, GLC, MES, FRU, EDO, PEG

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	0.39	0/3119	0.68	0/4253
1	B	0.41	0/3141	0.69	0/4286
1	C	0.39	0/3239	0.69	0/4419
All	All	0.40	0/9499	0.68	0/12958

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

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	2997	2950	2879	10	0
1	B	3000	2945	2865	9	0
1	C	3068	3003	2900	11	0
2	D	22	20	10	1	0
2	E	22	20	10	0	0
3	F	56	52	52	0	0
3	G	56	52	52	0	0
3	H	56	52	52	0	0
3	I	56	52	52	0	0
4	A	12	12	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	4	6	6	0	0
6	A	12	12	12	1	0
7	C	7	10	10	4	0
8	A	306	0	0	0	0
8	B	320	0	0	0	0
8	C	317	0	0	4	0
All	All	10311	9186	8912	32	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 2.

All (32) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:GLU:OE2	2:D:2:FRU:H12	2.02	0.60
7:C:501:PEG:H32	8:C:602:HOH:O	2.05	0.55
1:A:2:ALA:N	1:B:9:THR:HG1	2.04	0.54
1:C:36:LYS:HE3	7:C:501:PEG:H12	1.92	0.51
1:C:400:HIS:HD2	8:C:866:HOH:O	1.92	0.51
7:C:501:PEG:C2	8:C:602:HOH:O	2.60	0.48
6:A:503:MES:H81	1:C:71:LYS:HD3	1.96	0.48
1:A:253:ARG:NH1	1:A:277:GLU:OE1	2.45	0.46
1:B:281:SER:HA	1:B:314:THR:O	2.16	0.46
1:B:167:GLU:HA	1:B:189:THR:O	2.17	0.45
1:B:212:THR:HA	1:B:236:SER:O	2.17	0.45
1:C:139:GLY:HA3	1:C:173:TYR:O	2.16	0.44
1:B:257:THR:HA	1:B:281:SER:O	2.18	0.44
1:A:281:SER:HA	1:A:314:THR:O	2.18	0.43
1:C:71:LYS:HA	1:C:133[B]:GLN:O	2.19	0.43
1:A:139:GLY:HA3	1:A:173:TYR:O	2.18	0.43
1:B:264:ASN:HB2	1:B:265:PHE:CE1	2.55	0.42
1:C:281:SER:HA	1:C:314:THR:O	2.20	0.42
1:B:189:THR:HA	1:B:212:THR:O	2.20	0.42
1:C:189:THR:HA	1:C:212:THR:O	2.20	0.41
1:B:133:GLN:HA	1:B:167:GLU:O	2.20	0.41
1:C:328:THR:HA	1:C:329:PRO:C	2.41	0.41
1:A:153:THR:HA	1:A:176:HIS:O	2.20	0.41
1:A:189:THR:HA	1:A:212:THR:O	2.21	0.41
1:C:314:THR:HA	1:C:348:ALA:O	2.22	0.40
1:A:212:THR:HA	1:A:236:SER:O	2.22	0.40
1:B:348:ALA:HA	1:B:372:LEU:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:C:501:PEG:C3	8:C:602:HOH:O	2.66	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	407/400 (102%)	392 (96%)	15 (4%)	0	100	100
1	B	411/400 (103%)	391 (95%)	20 (5%)	0	100	100
1	C	420/400 (105%)	401 (96%)	19 (4%)	0	100	100
All	All	1238/1200 (103%)	1184 (96%)	54 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	335/328 (102%)	334 (100%)	1 (0%)	92	82
1	B	335/328 (102%)	334 (100%)	1 (0%)	92	82
1	C	348/328 (106%)	346 (99%)	2 (1%)	86	68
All	All	1018/984 (104%)	1014 (100%)	4 (0%)	91	80

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	253	ARG
1	B	253	ARG
1	C	233	LEU
1	C	253	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	207	GLN
1	C	400	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

24 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	Z9N	D	1	2	11,11,12	0.99	0	15,15,18	0.89	0
2	FRU	D	2	2	11,11,12	1.60	2 (18%)	15,15,18	1.19	1 (6%)
2	Z9N	E	1	2	11,11,12	1.51	1 (9%)	15,15,18	2.00	5 (33%)
2	FRU	E	2	2	11,11,12	1.09	1 (9%)	15,15,18	1.21	1 (6%)
3	GLC	F	1	3	12,12,12	2.06	4 (33%)	17,17,17	1.95	4 (23%)
3	FRU	F	2	3	11,11,12	1.40	2 (18%)	15,15,18	1.01	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	FRU	F	3	3	11,11,12	1.19	1 (9%)	15,15,18	0.94	0
3	FRU	F	4	3	11,11,12	1.09	1 (9%)	15,15,18	0.71	0
3	FRU	F	5	3	11,11,12	1.18	2 (18%)	15,15,18	0.91	0
3	GLC	G	1	3	12,12,12	2.05	3 (25%)	17,17,17	1.71	4 (23%)
3	FRU	G	2	3	11,11,12	1.36	2 (18%)	15,15,18	1.26	3 (20%)
3	FRU	G	3	3	11,11,12	1.14	1 (9%)	15,15,18	1.11	2 (13%)
3	FRU	G	4	3	11,11,12	1.10	1 (9%)	15,15,18	0.84	0
3	FRU	G	5	3	11,11,12	1.15	1 (9%)	15,15,18	1.28	2 (13%)
3	GLC	H	1	3	12,12,12	2.11	4 (33%)	17,17,17	1.38	1 (5%)
3	FRU	H	2	3	11,11,12	1.38	2 (18%)	15,15,18	1.55	2 (13%)
3	FRU	H	3	3	11,11,12	1.15	1 (9%)	15,15,18	0.86	0
3	FRU	H	4	3	11,11,12	0.99	1 (9%)	15,15,18	0.96	0
3	FRU	H	5	3	11,11,12	1.14	1 (9%)	15,15,18	1.10	0
3	GLC	I	1	3	12,12,12	2.06	4 (33%)	17,17,17	1.63	4 (23%)
3	FRU	I	2	3	11,11,12	1.44	2 (18%)	15,15,18	1.19	1 (6%)
3	FRU	I	3	3	11,11,12	1.18	1 (9%)	15,15,18	0.95	0
3	FRU	I	4	3	11,11,12	1.09	1 (9%)	15,15,18	0.93	0
3	FRU	I	5	3	11,11,12	1.17	1 (9%)	15,15,18	1.09	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
2	Z9N	D	1	2	-	2/4/20/24	0/1/1/1
2	FRU	D	2	2	-	4/4/20/24	0/1/1/1
2	Z9N	E	1	2	-	2/4/20/24	0/1/1/1
2	FRU	E	2	2	-	2/4/20/24	0/1/1/1
3	GLC	F	1	3	-	0/2/22/22	0/1/1/1
3	FRU	F	2	3	-	0/4/20/24	0/1/1/1
3	FRU	F	3	3	-	0/4/20/24	0/1/1/1
3	FRU	F	4	3	-	0/4/20/24	0/1/1/1
3	FRU	F	5	3	-	2/4/20/24	0/1/1/1
3	GLC	G	1	3	-	0/2/22/22	0/1/1/1
3	FRU	G	2	3	-	0/4/20/24	0/1/1/1
3	FRU	G	3	3	-	0/4/20/24	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	FRU	G	4	3	-	0/4/20/24	0/1/1/1
3	FRU	G	5	3	-	0/4/20/24	0/1/1/1
3	GLC	H	1	3	-	0/2/22/22	0/1/1/1
3	FRU	H	2	3	-	0/4/20/24	0/1/1/1
3	FRU	H	3	3	-	0/4/20/24	0/1/1/1
3	FRU	H	4	3	-	0/4/20/24	0/1/1/1
3	FRU	H	5	3	-	0/4/20/24	0/1/1/1
3	GLC	I	1	3	-	0/2/22/22	0/1/1/1
3	FRU	I	2	3	-	0/4/20/24	0/1/1/1
3	FRU	I	3	3	-	1/4/20/24	0/1/1/1
3	FRU	I	4	3	-	1/4/20/24	0/1/1/1
3	FRU	I	5	3	-	2/4/20/24	0/1/1/1

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	1	GLC	C1-C2	4.35	1.62	1.52
3	F	1	GLC	C1-C2	4.32	1.62	1.52
3	I	1	GLC	C1-C2	4.31	1.62	1.52
3	G	1	GLC	C1-C2	4.26	1.62	1.52
3	F	1	GLC	C3-C2	4.12	1.62	1.52
3	H	1	GLC	C3-C2	4.11	1.62	1.52
3	G	1	GLC	C3-C2	4.02	1.62	1.52
3	I	1	GLC	C3-C2	4.00	1.62	1.52
3	I	2	FRU	C1-C2	3.82	1.64	1.51
3	F	2	FRU	C1-C2	3.66	1.64	1.51
3	H	2	FRU	C1-C2	3.60	1.63	1.51
3	G	2	FRU	C1-C2	3.59	1.63	1.51
2	E	1	Z9N	O6-C6	-3.36	1.28	1.42
3	I	3	FRU	C1-C2	3.20	1.62	1.51
3	F	3	FRU	C1-C2	3.10	1.62	1.51
3	G	3	FRU	C1-C2	3.10	1.62	1.51
3	H	3	FRU	C1-C2	3.04	1.62	1.51
3	G	4	FRU	C1-C2	2.89	1.61	1.51
3	I	4	FRU	C1-C2	2.77	1.61	1.51
3	F	4	FRU	C1-C2	2.64	1.60	1.51
3	G	1	GLC	C6-C5	2.64	1.60	1.51
3	H	1	GLC	C6-C5	2.62	1.60	1.51
3	G	5	FRU	C1-C2	2.59	1.60	1.51
3	I	1	GLC	C6-C5	2.56	1.60	1.51
3	F	5	FRU	C1-C2	2.53	1.60	1.51
3	H	4	FRU	C1-C2	2.50	1.60	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	1	GLC	C6-C5	2.49	1.60	1.51
3	I	5	FRU	C1-C2	2.48	1.60	1.51
2	D	2	FRU	O4-C4	-2.45	1.37	1.43
3	H	5	FRU	C1-C2	2.24	1.59	1.51
3	H	1	GLC	O1-C1	2.24	1.46	1.39
3	F	2	FRU	C6-C5	2.23	1.59	1.51
3	G	2	FRU	C6-C5	2.17	1.59	1.51
3	H	2	FRU	C6-C5	2.14	1.59	1.51
3	I	2	FRU	C6-C5	2.13	1.59	1.51
3	I	1	GLC	O1-C1	2.10	1.46	1.39
2	D	2	FRU	C3-C2	2.09	1.58	1.53
3	F	1	GLC	O1-C1	2.08	1.46	1.39
2	E	2	FRU	O3-C3	-2.05	1.38	1.43
3	F	5	FRU	C6-C5	2.03	1.58	1.51

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	F	1	GLC	C1-O5-C5	5.08	123.24	113.66
3	G	1	GLC	C1-O5-C5	4.04	121.29	113.66
3	I	1	GLC	C1-O5-C5	3.97	121.15	113.66
2	E	1	Z9N	C3-C4-C5	-3.68	95.50	102.64
3	F	1	GLC	C1-C2-C3	-3.60	102.85	110.31
3	H	2	FRU	C4-C3-C2	-3.56	95.73	102.64
2	E	1	Z9N	O4-C4-C5	3.52	121.24	111.05
3	G	1	GLC	C1-C2-C3	-3.30	103.46	110.31
2	E	1	Z9N	C4-C3-C2	-3.21	96.41	102.64
2	E	1	Z9N	O5-C5-C6	-2.91	102.92	109.21
2	E	1	Z9N	O6-C6-C5	-2.87	101.45	111.29
3	H	2	FRU	O1-C1-C2	-2.80	101.68	111.29
3	G	2	FRU	O1-C1-C2	-2.79	101.73	111.29
3	F	1	GLC	O3-C3-C2	-2.74	104.00	110.35
3	I	2	FRU	O1-C1-C2	-2.72	101.95	111.29
3	F	2	FRU	O1-C1-C2	-2.72	101.96	111.29
3	H	1	GLC	C1-O5-C5	2.69	118.74	113.66
3	G	5	FRU	C1-C2-C3	-2.68	108.62	115.09
2	E	2	FRU	O5-C2-C1	-2.65	103.48	109.21
3	G	5	FRU	O5-C2-C1	2.60	114.84	109.21
3	I	1	GLC	C1-C2-C3	-2.50	105.14	110.31
3	G	2	FRU	C4-C3-C2	-2.37	98.05	102.64
3	G	3	FRU	O1-C1-C2	-2.34	103.27	111.29
3	G	1	GLC	C6-C5-C4	-2.27	107.68	113.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	2	FRU	O4-C4-C3	-2.26	104.50	111.82
3	I	1	GLC	C6-C5-C4	-2.21	107.82	113.00
3	I	1	GLC	O2-C2-C1	-2.13	104.22	109.16
3	F	1	GLC	O5-C1-C2	-2.11	106.53	110.28
3	G	1	GLC	O1-C1-O5	-2.08	104.14	110.38
3	G	2	FRU	O5-C2-C1	-2.04	104.81	109.21
3	G	3	FRU	C4-C3-C2	-2.00	98.75	102.64

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	Z9N	C4-C5-C6-O6
2	E	1	Z9N	O5-C5-C6-O6
3	I	5	FRU	C4-C5-C6-O6
2	E	2	FRU	O5-C5-C6-O6
2	E	2	FRU	C4-C5-C6-O6
2	D	2	FRU	O5-C5-C6-O6
2	D	2	FRU	C4-C5-C6-O6
3	I	5	FRU	O5-C5-C6-O6
2	D	1	Z9N	O5-C5-C6-O6
3	F	5	FRU	C4-C5-C6-O6
2	D	2	FRU	O1-C1-C2-C3
2	D	1	Z9N	C4-C5-C6-O6
3	F	5	FRU	O5-C5-C6-O6
3	I	4	FRU	O1-C1-C2-O5
2	D	2	FRU	O1-C1-C2-O5
3	I	3	FRU	O1-C1-C2-C3

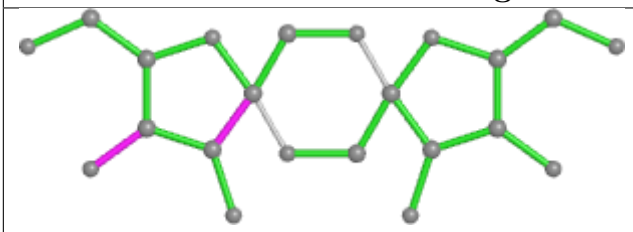
There are no ring outliers.

1 monomer is involved in 1 short contact:

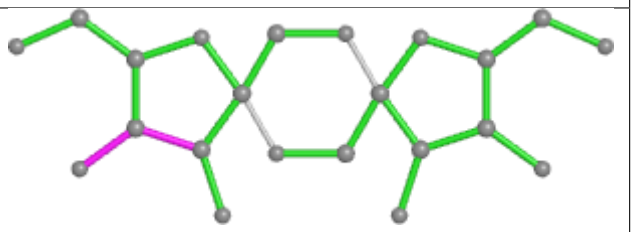
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	FRU	1	0

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

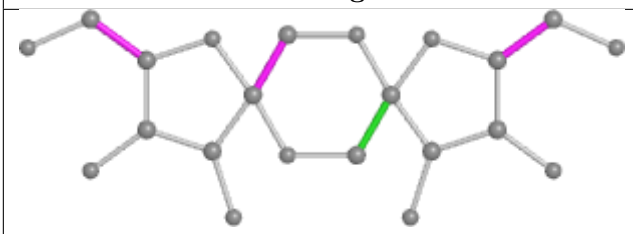
Oligosaccharide Chain D



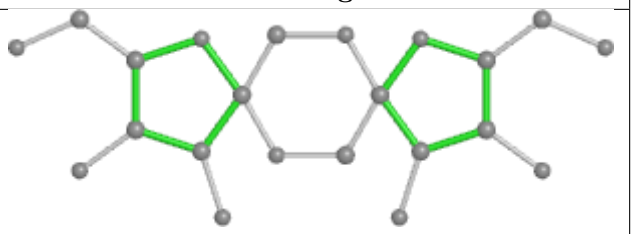
Bond lengths



Bond angles

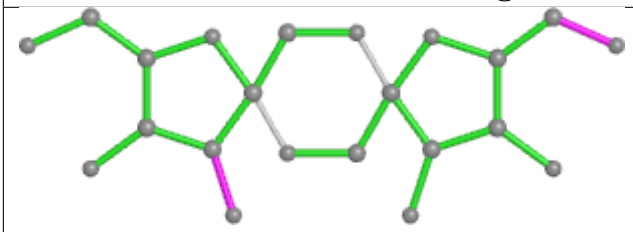


Torsions

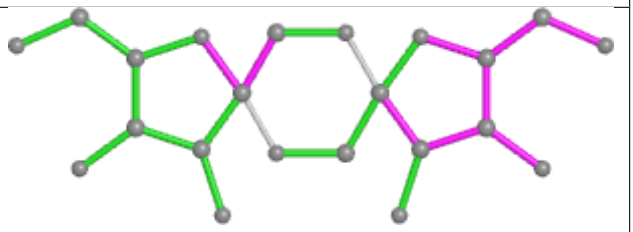


Rings

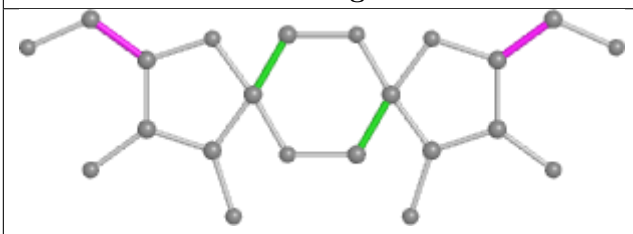
Oligosaccharide Chain E



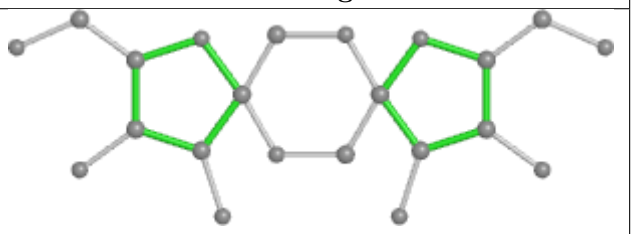
Bond lengths



Bond angles

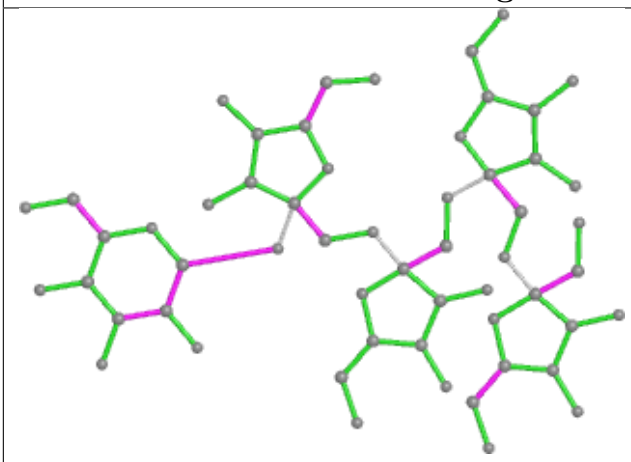


Torsions

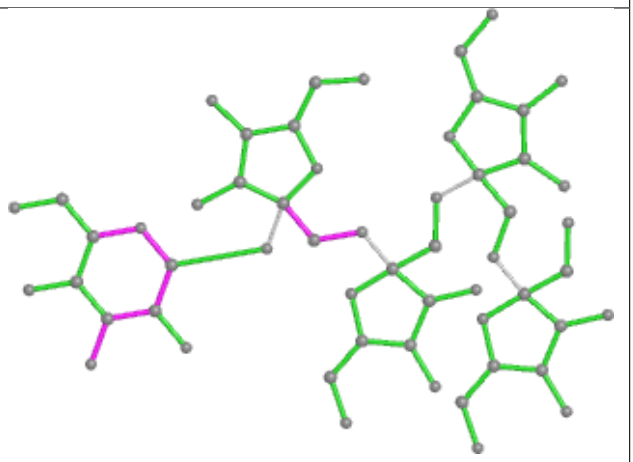


Rings

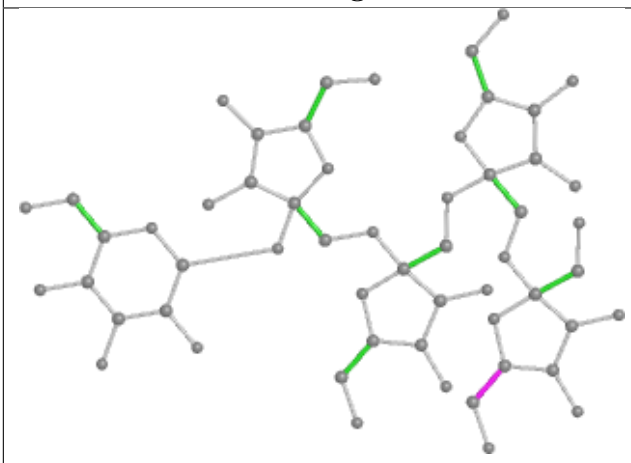
Oligosaccharide Chain F



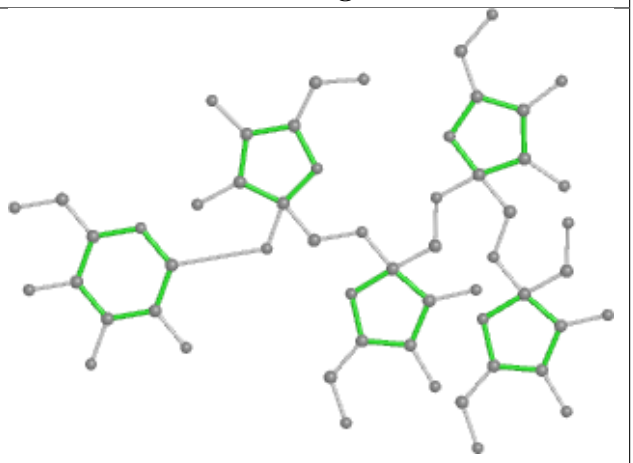
Bond lengths



Bond angles

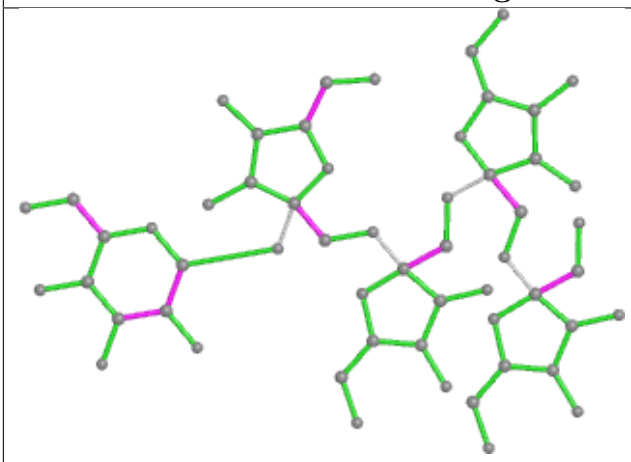


Torsions

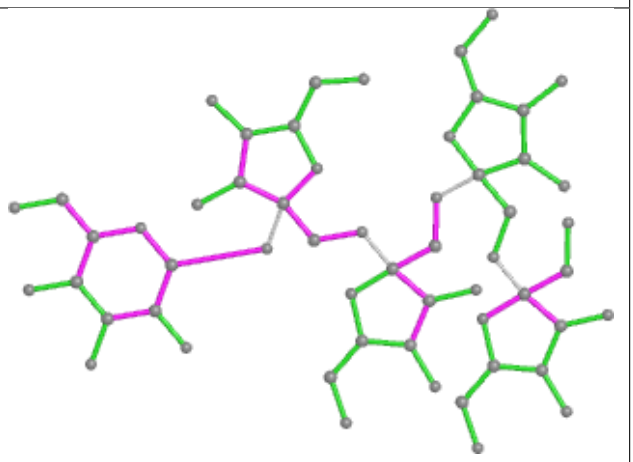


Rings

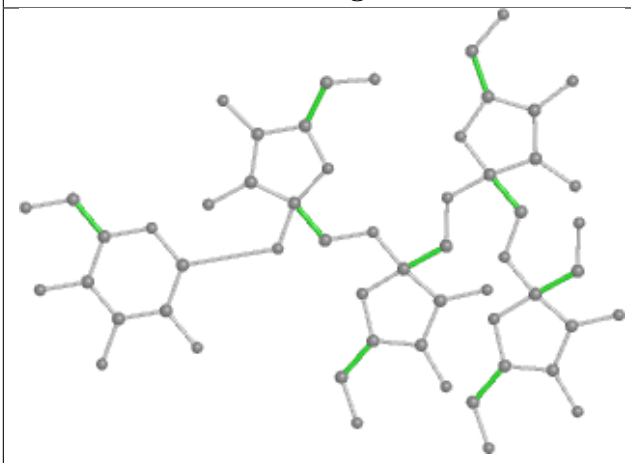
Oligosaccharide Chain G



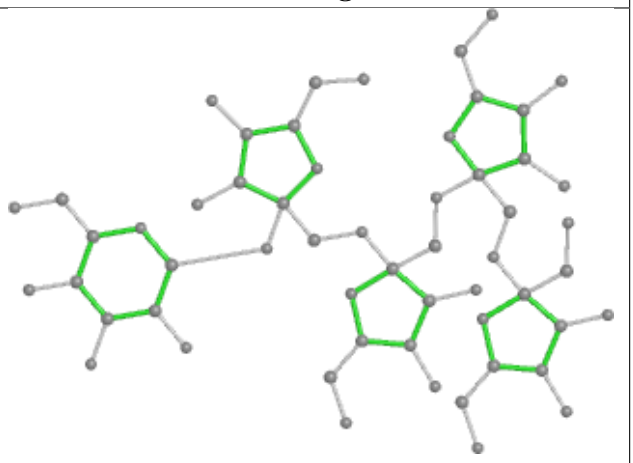
Bond lengths



Bond angles

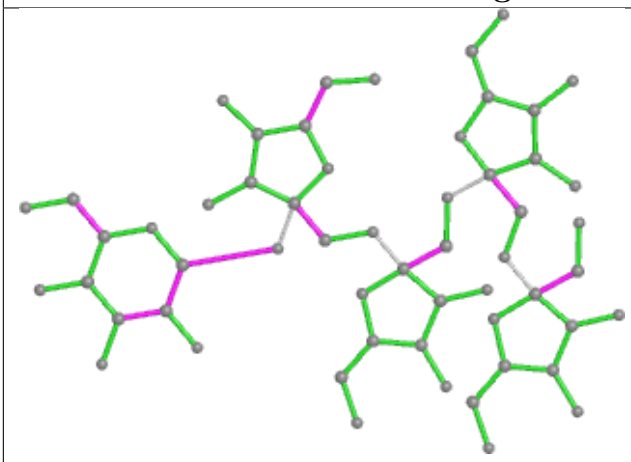


Torsions

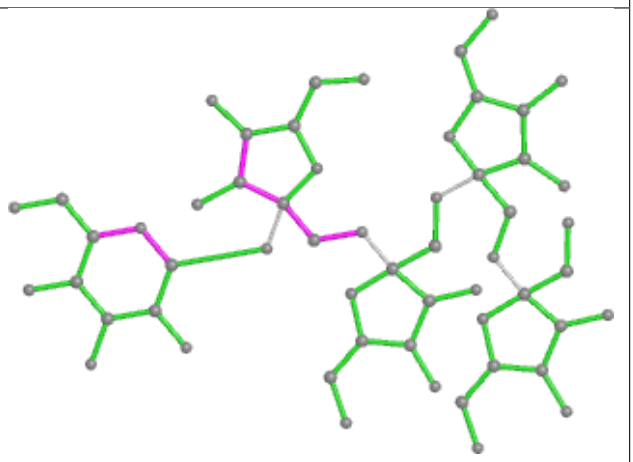


Rings

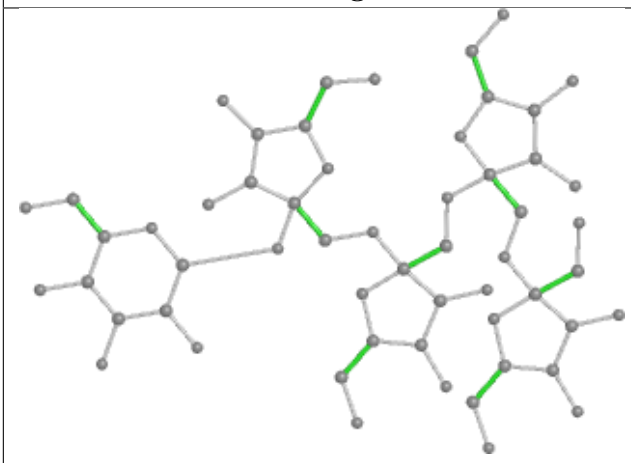
Oligosaccharide Chain H



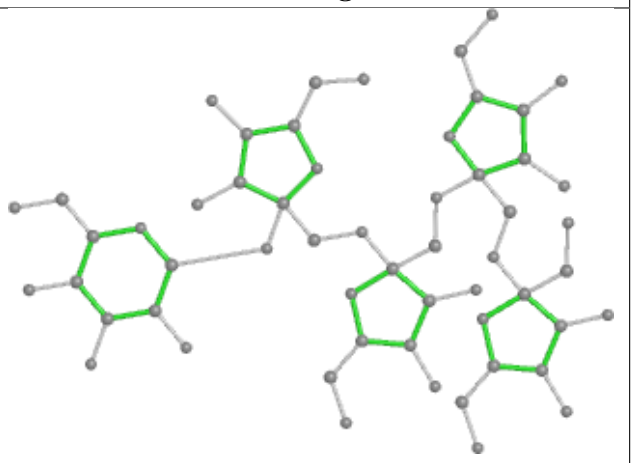
Bond lengths



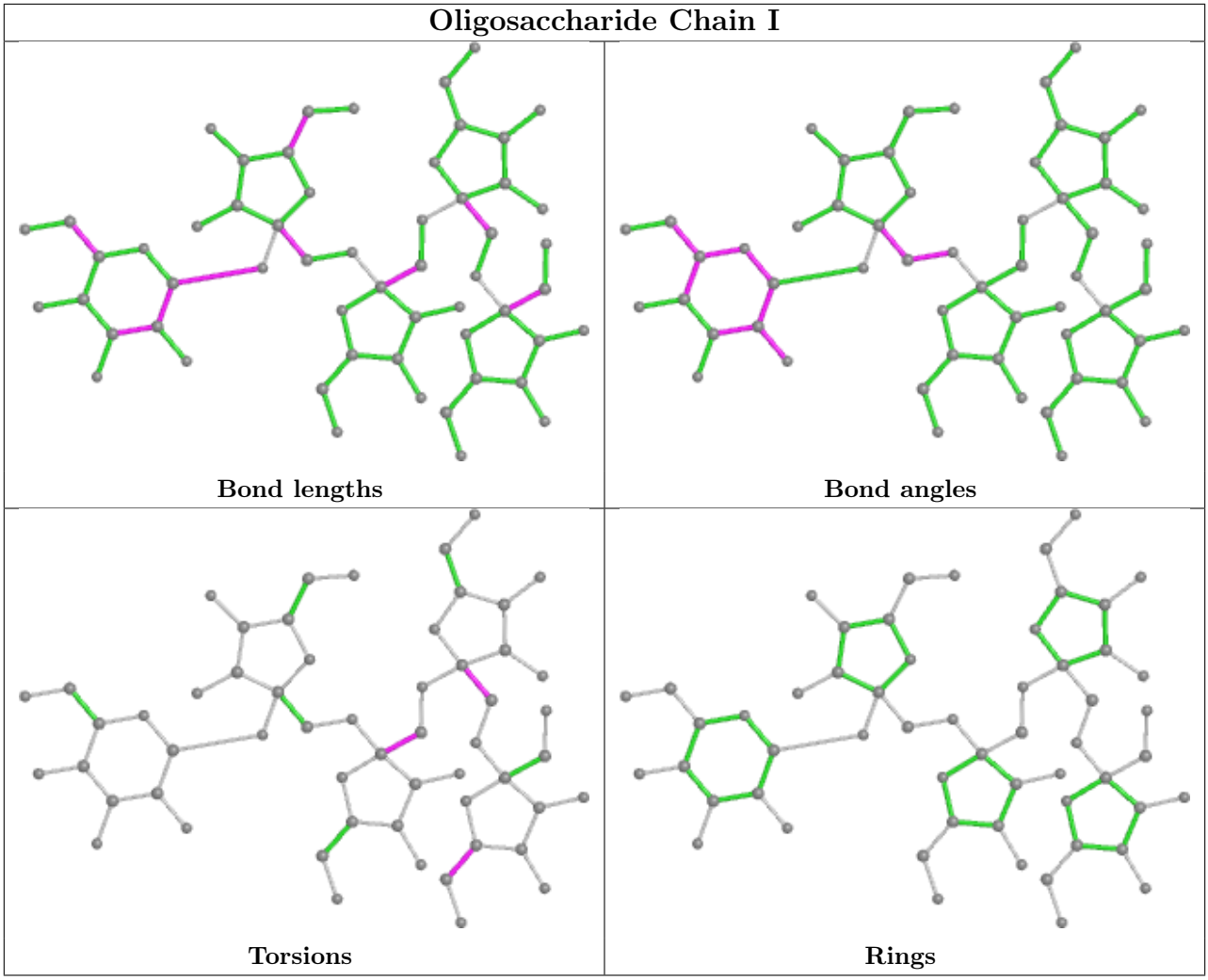
Bond angles



Torsions



Rings



5.6 Ligand geometry (i)

4 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	FRU	A	501	-	11,12,12	0.60	0	10,18,18	1.00	1 (10%)
7	PEG	C	501	-	6,6,6	0.25	0	5,5,5	0.25	0
6	MES	A	503	-	12,12,12	2.22	1 (8%)	14,16,16	2.28	4 (28%)
5	EDO	A	502	-	3,3,3	0.49	0	2,2,2	0.34	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FRU	A	501	-	-	0/5/24/24	0/1/1/1
7	PEG	C	501	-	-	2/4/4/4	-
6	MES	A	503	-	-	2/6/14/14	0/1/1/1
5	EDO	A	502	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	503	MES	C8-S	-7.36	1.67	1.77

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	A	503	MES	C7-N4-C3	5.04	124.12	111.23
6	A	503	MES	C5-N4-C3	3.91	117.63	108.83
6	A	503	MES	O3S-S-C8	3.86	112.01	105.77
6	A	503	MES	O2S-S-C8	2.65	110.11	106.92
4	A	501	FRU	O6-C6-C5	-2.11	104.06	111.29

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	503	MES	C8-C7-N4-C3
7	C	501	PEG	O1-C1-C2-O2
7	C	501	PEG	O2-C3-C4-O4
6	A	503	MES	N4-C7-C8-S

There are no ring outliers.

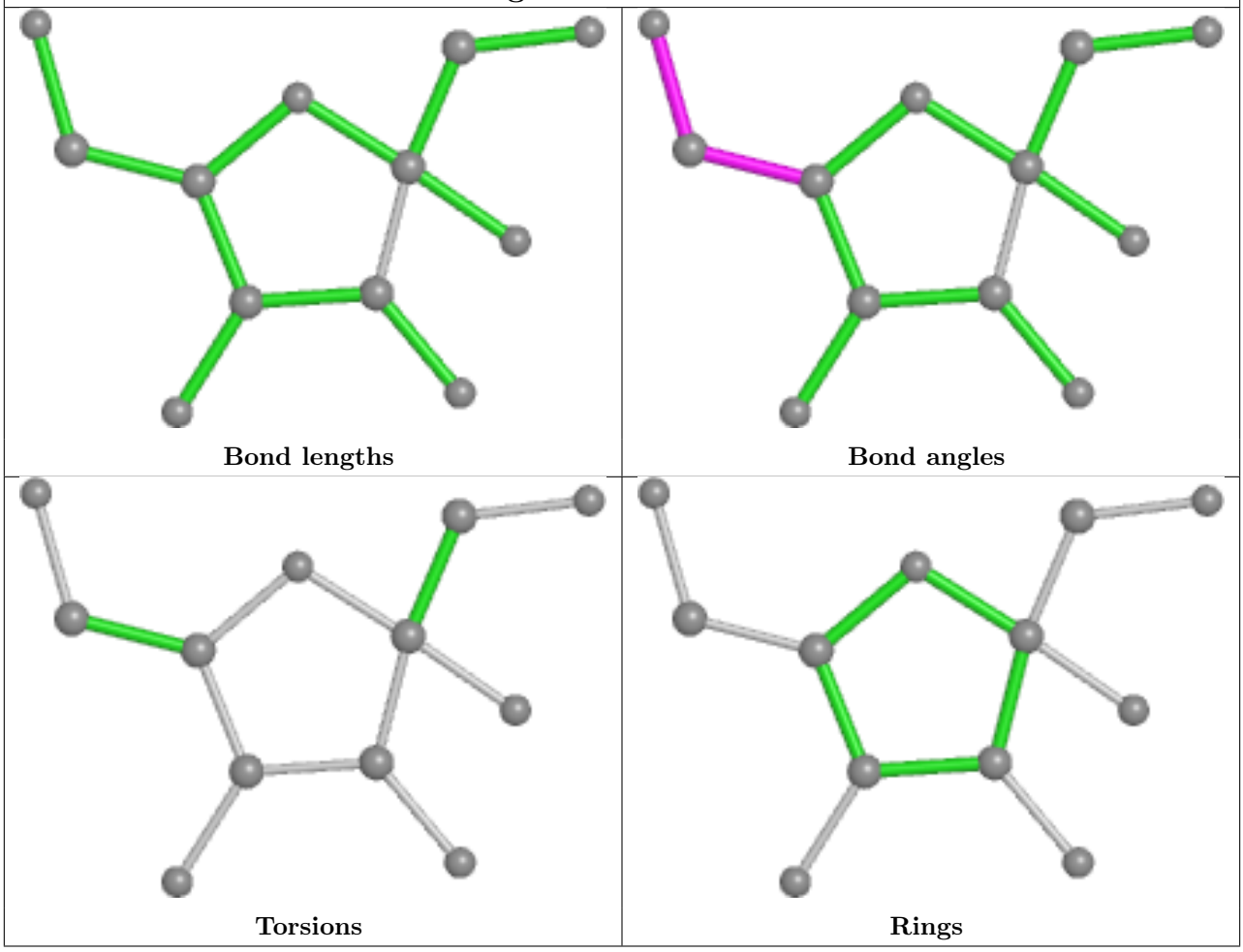
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	501	PEG	4	0
6	A	503	MES	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 then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand FRU A 501



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, 95^{th} 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	393/400 (98%)	-0.07	4 (1%)	82	82	7, 10, 19, 27	0
1	B	393/400 (98%)	-0.16	3 (0%)	86	86	7, 10, 17, 28	0
1	C	399/400 (99%)	-0.08	2 (0%)	91	92	7, 11, 18, 39	0
All	All	1185/1200 (98%)	-0.10	9 (0%)	86	86	7, 10, 18, 39	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	ALA	3.0
1	A	16	VAL	2.7
1	A	14	ALA	2.7
1	C	14	ALA	2.3
1	B	14	ALA	2.2
1	B	119[A]	GLY	2.2
1	B	330	ALA	2.1
1	C	397	HIS	2.1
1	A	87	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

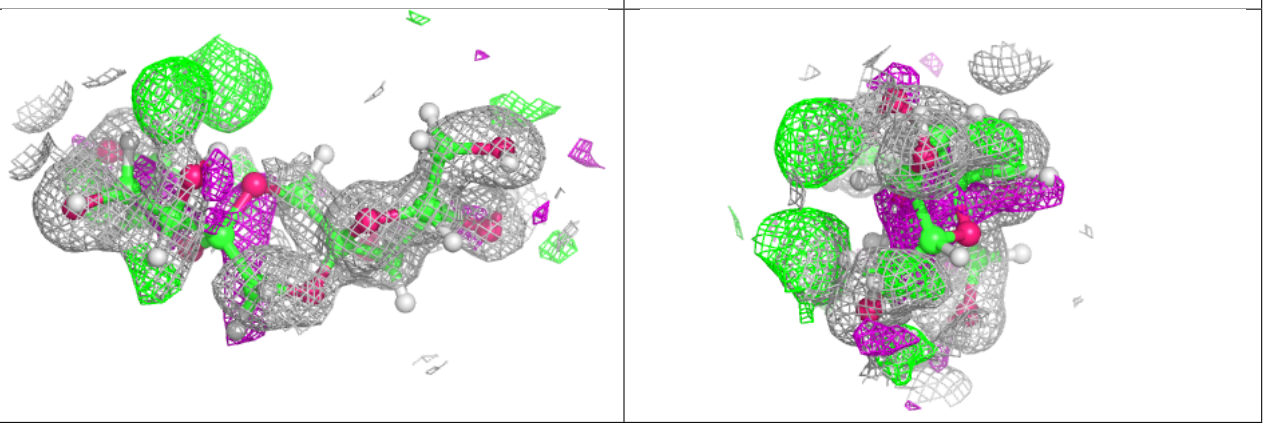
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	FRU	G	5	11/12	0.59	0.43	36,49,59,66	0
3	GLC	H	1	12/12	0.67	0.26	25,37,44,53	0
3	FRU	I	5	11/12	0.69	0.34	32,42,53,60	0
3	FRU	F	5	11/12	0.77	0.28	20,31,43,55	0
2	FRU	D	2	11/12	0.79	0.34	15,41,57,63	0
2	Z9N	E	1	11/12	0.86	0.36	18,30,59,59	0
3	GLC	F	1	12/12	0.86	0.14	14,26,33,35	0
2	FRU	E	2	11/12	0.88	0.22	20,29,44,45	0
3	GLC	G	1	12/12	0.89	0.22	19,36,49,58	0
2	Z9N	D	1	11/12	0.89	0.18	14,23,31,52	0
3	FRU	I	4	11/12	0.90	0.12	12,16,24,31	0
3	GLC	I	1	12/12	0.92	0.31	18,39,48,57	0
3	FRU	H	5	11/12	0.93	0.11	13,17,21,26	0
3	FRU	F	4	11/12	0.93	0.10	14,17,21,23	0
3	FRU	I	2	11/12	0.93	0.11	14,20,26,28	0
3	FRU	G	2	11/12	0.93	0.10	11,15,20,24	0
3	FRU	H	2	11/12	0.93	0.10	11,19,33,48	0
3	FRU	G	4	11/12	0.95	0.14	10,14,22,34	0
3	FRU	G	3	11/12	0.96	0.08	10,13,16,19	0
3	FRU	I	3	11/12	0.96	0.10	12,15,19,21	0
3	FRU	F	2	11/12	0.97	0.07	10,12,15,17	0
3	FRU	H	3	11/12	0.97	0.07	9,12,15,17	0
3	FRU	H	4	11/12	0.97	0.08	7,11,14,16	0
3	FRU	F	3	11/12	0.98	0.06	8,11,13,14	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.

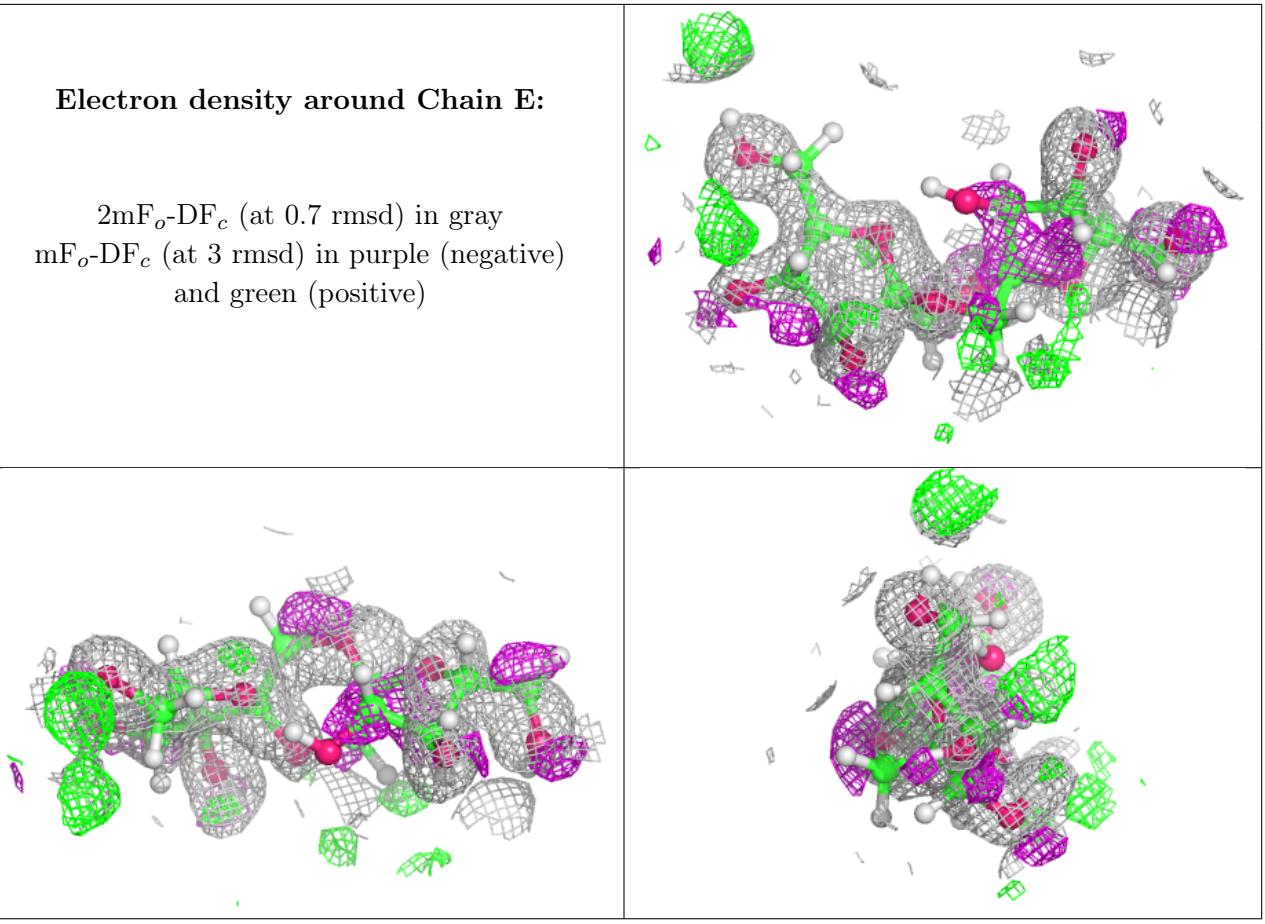
Electron density around Chain D:

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



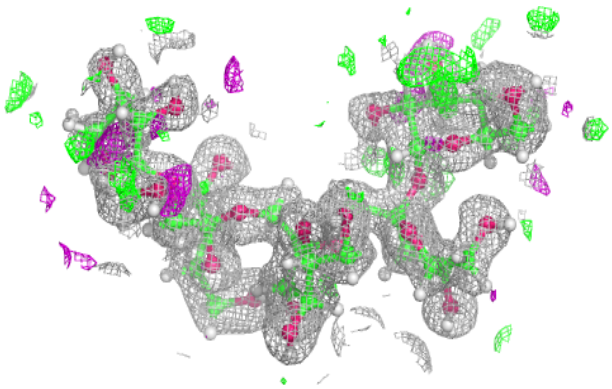
Electron density around Chain E:

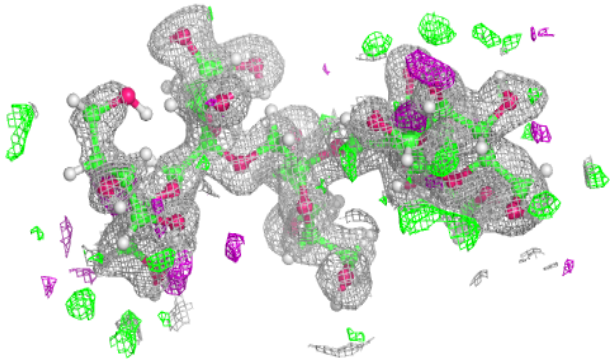
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around Chain F:

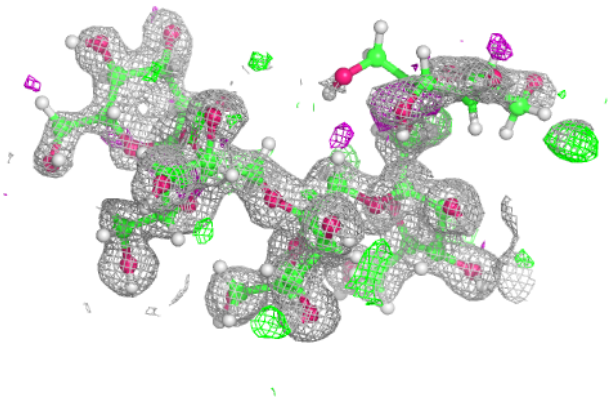
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

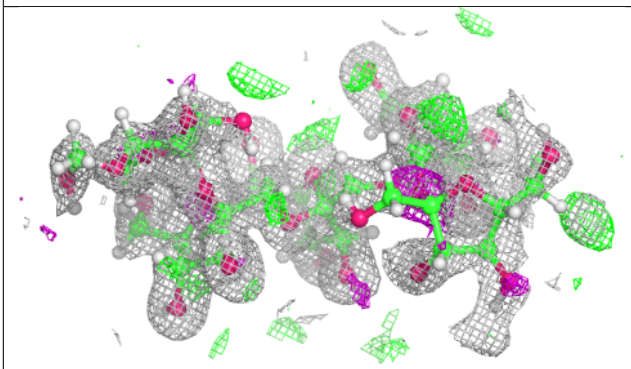


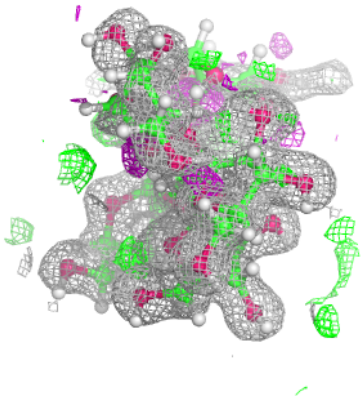


Electron density around Chain G:

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

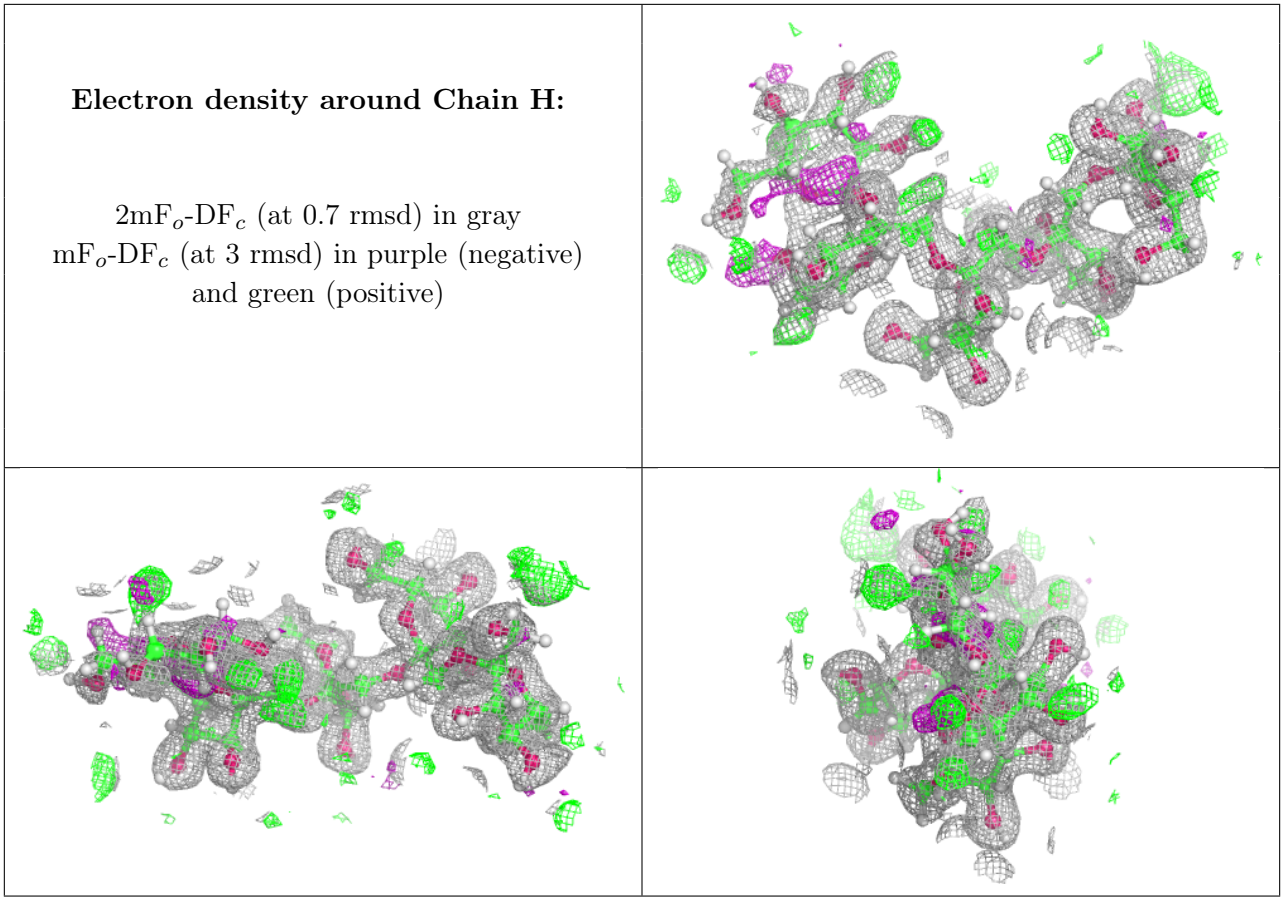






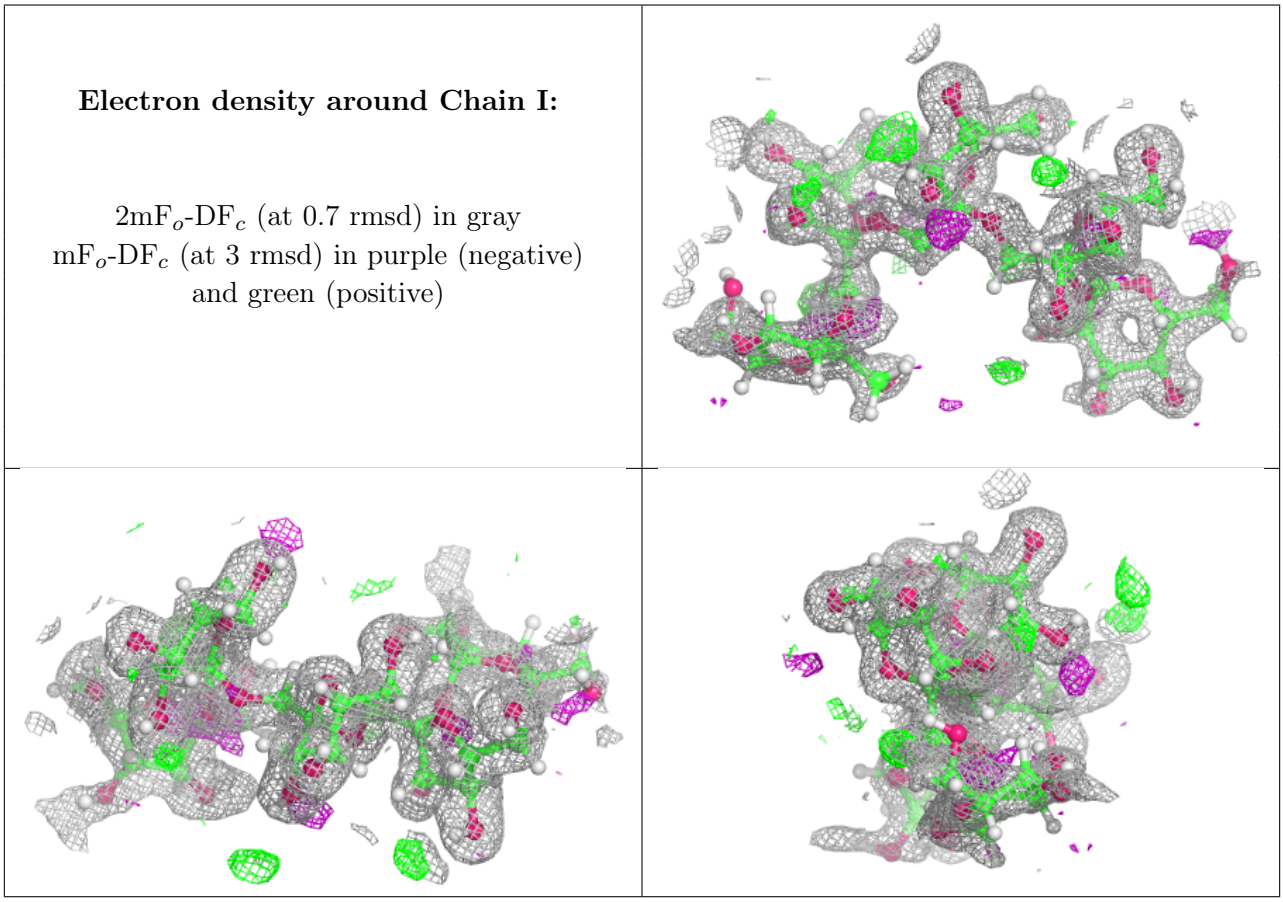
Electron density around Chain H:

$2mF_o$-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

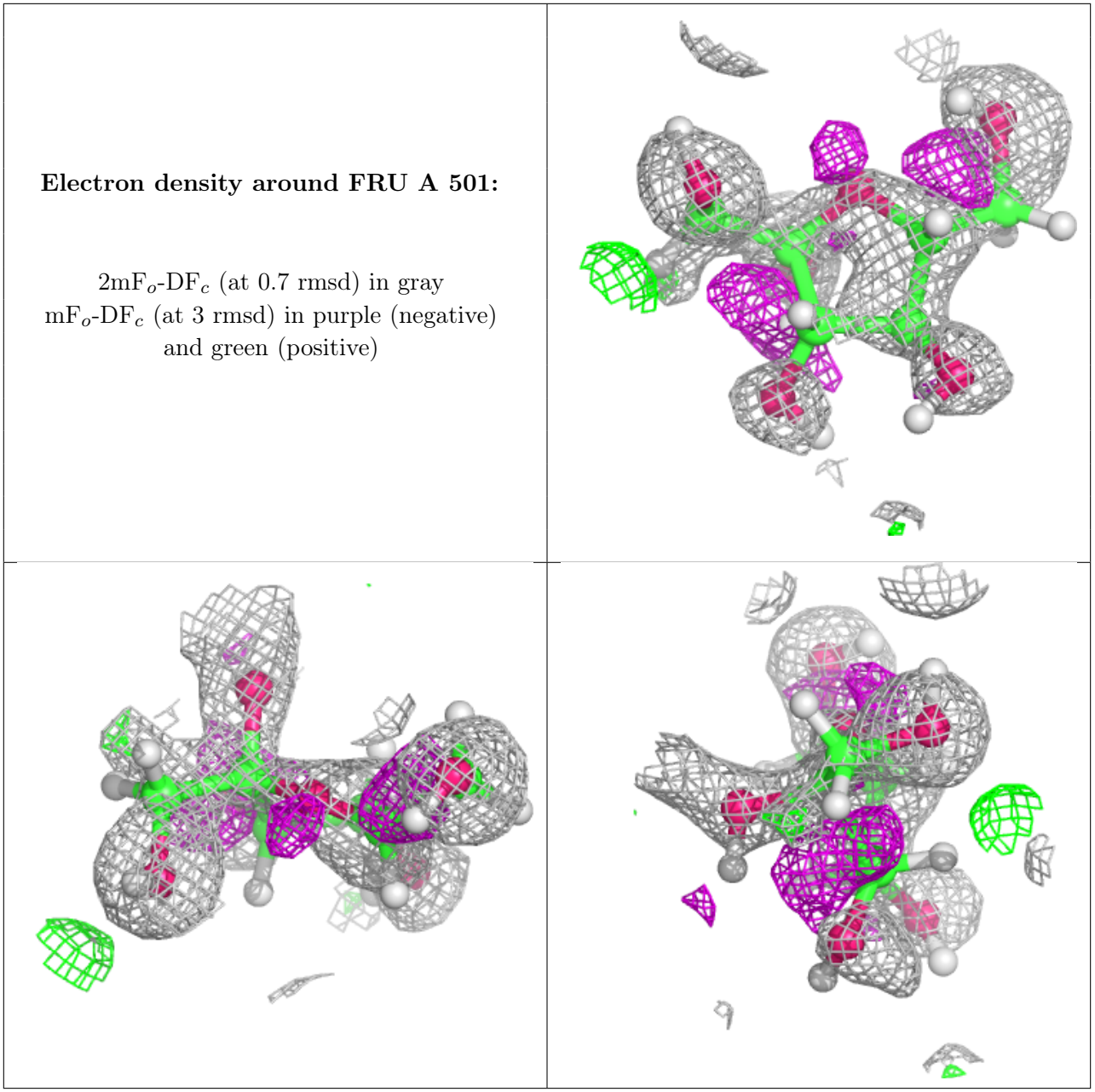
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
6	MES	A	503	12/12	0.54	0.35	39,53,79,101	0
7	PEG	C	501	7/7	0.77	0.40	26,37,45,46	0
4	FRU	A	501	12/12	0.78	0.35	19,39,47,56	0
5	EDO	A	502	4/4	0.85	0.17	29,39,47,48	0

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

Electron density around FRU A 501:

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



6.5 Other polymers (i)

There are no such residues in this entry.