



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:29 PM UTC

PDB ID : 4IFP / pdb_00004ifp
Title : X-ray Crystal Structure of Human NLRP1 CARD Domain
Authors : Jin, T.; Curry, J.; Smith, P.; Jiang, J.; Xiao, T.
Deposited on : 2012-12-14
Resolution : 1.99 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

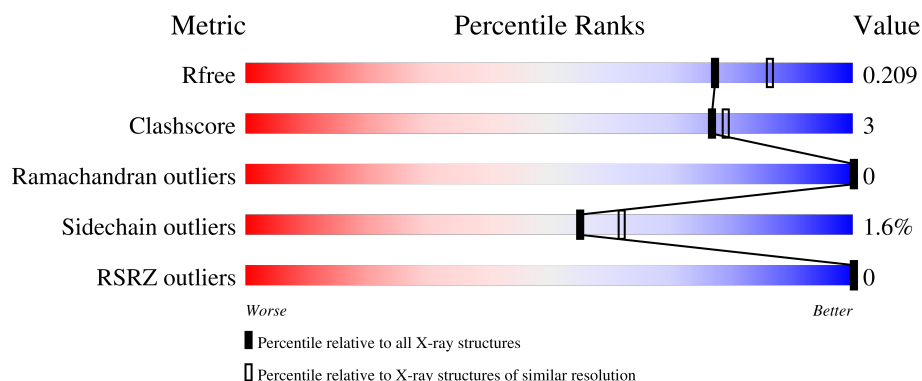
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	466	
1	B	466	
1	C	466	
2	D	2	
2	E	2	

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Mol	Chain	Length	Quality of chain
2	F	2	 100%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12414 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Maltose-binding periplasmic protein,NACHT, LRR and PYD domains-containing protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	458	Total	C	N	O	S	0	4	0
			3596	2309	599	679	9			
1	B	458	Total	C	N	O	S	0	3	0
			3588	2304	595	680	9			
1	C	458	Total	C	N	O	S	0	5	0
			3602	2314	600	679	9			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P0AEX9
A	83	ALA	ASP	conflict	UNP P0AEX9
A	84	ALA	LYS	conflict	UNP P0AEX9
A	173	ALA	GLU	conflict	UNP P0AEX9
A	174	ALA	ASN	conflict	UNP P0AEX9
A	240	ALA	LYS	conflict	UNP P0AEX9
A	360	ALA	GLU	conflict	UNP P0AEX9
A	363	ALA	LYS	conflict	UNP P0AEX9
A	364	ALA	ASP	conflict	UNP P0AEX9
A	368	ASN	-	linker	UNP P0AEX9
A	369	ALA	-	linker	UNP P0AEX9
A	370	VAL	-	linker	UNP P0AEX9
A	371	ASP	-	linker	UNP P0AEX9
A	456	ALA	-	expression tag	UNP Q9C000
A	457	ALA	-	expression tag	UNP Q9C000
A	458	ALA	-	expression tag	UNP Q9C000
A	459	LEU	-	expression tag	UNP Q9C000
A	460	GLU	-	expression tag	UNP Q9C000
A	461	HIS	-	expression tag	UNP Q9C000
A	462	HIS	-	expression tag	UNP Q9C000
A	463	HIS	-	expression tag	UNP Q9C000
A	464	HIS	-	expression tag	UNP Q9C000

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Chain	Residue	Modelled	Actual	Comment	Reference
A	465	HIS	-	expression tag	UNP Q9C000
A	466	HIS	-	expression tag	UNP Q9C000
B	1	MET	-	initiating methionine	UNP P0AEX9
B	83	ALA	ASP	conflict	UNP P0AEX9
B	84	ALA	LYS	conflict	UNP P0AEX9
B	173	ALA	GLU	conflict	UNP P0AEX9
B	174	ALA	ASN	conflict	UNP P0AEX9
B	240	ALA	LYS	conflict	UNP P0AEX9
B	360	ALA	GLU	conflict	UNP P0AEX9
B	363	ALA	LYS	conflict	UNP P0AEX9
B	364	ALA	ASP	conflict	UNP P0AEX9
B	368	ASN	-	linker	UNP P0AEX9
B	369	ALA	-	linker	UNP P0AEX9
B	370	VAL	-	linker	UNP P0AEX9
B	371	ASP	-	linker	UNP P0AEX9
B	456	ALA	-	expression tag	UNP Q9C000
B	457	ALA	-	expression tag	UNP Q9C000
B	458	ALA	-	expression tag	UNP Q9C000
B	459	LEU	-	expression tag	UNP Q9C000
B	460	GLU	-	expression tag	UNP Q9C000
B	461	HIS	-	expression tag	UNP Q9C000
B	462	HIS	-	expression tag	UNP Q9C000
B	463	HIS	-	expression tag	UNP Q9C000
B	464	HIS	-	expression tag	UNP Q9C000
B	465	HIS	-	expression tag	UNP Q9C000
B	466	HIS	-	expression tag	UNP Q9C000
C	1	MET	-	initiating methionine	UNP P0AEX9
C	83	ALA	ASP	conflict	UNP P0AEX9
C	84	ALA	LYS	conflict	UNP P0AEX9
C	173	ALA	GLU	conflict	UNP P0AEX9
C	174	ALA	ASN	conflict	UNP P0AEX9
C	240	ALA	LYS	conflict	UNP P0AEX9
C	360	ALA	GLU	conflict	UNP P0AEX9
C	363	ALA	LYS	conflict	UNP P0AEX9
C	364	ALA	ASP	conflict	UNP P0AEX9
C	368	ASN	-	linker	UNP P0AEX9
C	369	ALA	-	linker	UNP P0AEX9
C	370	VAL	-	linker	UNP P0AEX9
C	371	ASP	-	linker	UNP P0AEX9
C	456	ALA	-	expression tag	UNP Q9C000
C	457	ALA	-	expression tag	UNP Q9C000
C	458	ALA	-	expression tag	UNP Q9C000

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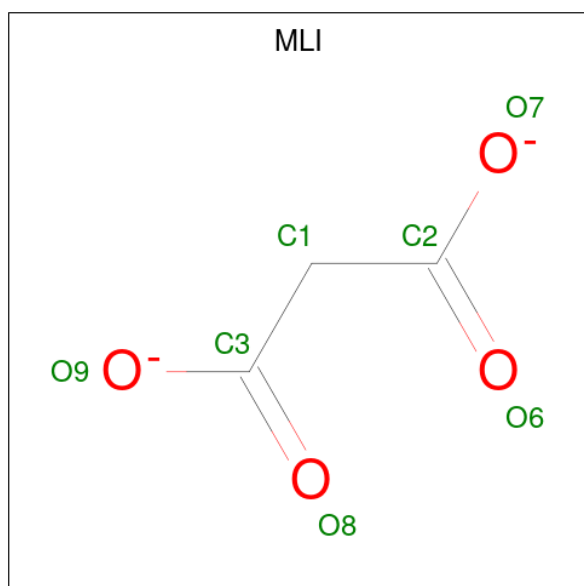
Chain	Residue	Modelled	Actual	Comment	Reference
C	459	LEU	-	expression tag	UNP Q9C000
C	460	GLU	-	expression tag	UNP Q9C000
C	461	HIS	-	expression tag	UNP Q9C000
C	462	HIS	-	expression tag	UNP Q9C000
C	463	HIS	-	expression tag	UNP Q9C000
C	464	HIS	-	expression tag	UNP Q9C000
C	465	HIS	-	expression tag	UNP Q9C000
C	466	HIS	-	expression tag	UNP Q9C000

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	D	2	Total	C	O	0	0	0
			23	12	11			
2	E	2	Total	C	O	0	0	0
			23	12	11			
2	F	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is MALONATE ION (CCD ID: MLI) (formula: C₃H₂O₄).

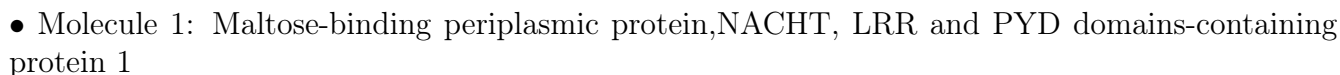
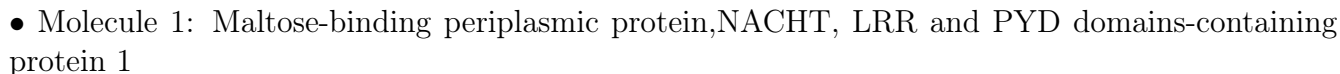


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 7 3 4	0	0
3	B	1	Total C O 7 3 4	0	0
3	C	1	Total C O 7 3 4	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	512	Total O 512 512	0	0
4	B	522	Total O 522 522	0	0
4	C	504	Total O 504 504	0	0

- Molecule 1: Maltose-binding periplasmic protein, NACHT, LRR and PYD domains-containing protein 1



GLC1
GLC2

- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain E:  100%

GLC1
GLC2

- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain F:  100%

GLC1
GLC2

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	187.01Å 108.77Å 121.27Å 90.00° 120.98° 90.00°	Depositor
Resolution (Å)	37.49 – 1.99 37.49 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.7 (37.49-1.99) 94.6 (37.49-2.00)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1217)	Depositor
R, R_{free}	0.170 , 0.208 0.172 , 0.209	Depositor DCC
R_{free} test set	7015 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.417 for 1/2*h+3/2*k,1/2*h-1/2*k,-1/2*h-1/2*k-l 0.417 for 1/2*h-3/2*k,-1/2*h-1/2*k,-1/2*h+1/2*k-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	12414	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MLI, GLC

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	0.47	0/3689	0.73	1/5008 (0.0%)
1	B	0.45	0/3678	0.74	0/4994
1	C	0.46	0/3698	0.74	0/5019
All	All	0.46	0/11065	0.74	1/15021 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	413	ASN	N-CA-C	5.27	118.18	111.69

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3596	0	3577	18	0
1	B	3588	0	3562	24	0
1	C	3602	0	3590	20	0
2	D	23	0	21	0	0
2	E	23	0	21	0	0
2	F	23	0	21	0	0
3	A	7	0	2	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	7	0	2	0	0
3	C	7	0	2	0	0
4	A	512	0	0	4	0
4	B	522	0	0	7	0
4	C	504	0	0	5	0
All	All	12414	0	10798	62	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (62) 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:345:ARG:NH1	4:A:785:HOH:O	2.27	0.67
1:B:306:LYS:NZ	4:B:1080:HOH:O	2.26	0.64
1:A:45:GLU:HB2	1:A:67:ARG:HD2	1.82	0.60
1:C:80:ILE:HG22	1:C:82:PRO:HD3	1.83	0.60
1:C:74:SER:HA	1:C:459:LEU:HD13	1.82	0.60
1:A:239[A]:SER:OG	1:A:241:VAL:HG23	2.00	0.60
1:A:143:LYS:NZ	4:A:1040:HOH:O	2.34	0.58
1:B:143:LYS:NZ	4:B:723:HOH:O	2.39	0.56
1:B:96:ASP:OD1	1:B:99:ARG:NH1	2.34	0.56
1:C:89:LYS:NZ	4:C:902:HOH:O	2.39	0.55
1:A:74:SER:HA	1:A:459:LEU:HD13	1.88	0.55
1:A:7:LYS:O	1:A:273:ASN:ND2	2.41	0.54
1:B:230:PRO:HA	1:B:233:TRP:CE2	2.43	0.54
1:C:239[A]:SER:OG	1:C:241:VAL:HG23	2.09	0.53
1:B:429:TRP:HB3	1:B:433:CYS:HB2	1.91	0.53
1:A:69:GLY:HA3	1:A:333:ASN:O	2.09	0.52
1:B:74:SER:HA	1:B:459:LEU:HD13	1.91	0.52
1:C:45:GLU:HB2	1:C:67:ARG:HD2	1.92	0.52
1:B:292:GLU:HG2	1:B:296:LYS:HD3	1.92	0.51
1:C:143:LYS:NZ	4:C:1048:HOH:O	2.33	0.51
1:B:69:GLY:HA3	1:B:333:ASN:O	2.11	0.50
1:C:298:LYS:NZ	4:C:820:HOH:O	2.37	0.50
1:A:371:ASP:HB3	1:A:431:ARG:NH2	2.27	0.49
1:C:69:GLY:HA3	1:C:333:ASN:O	2.12	0.48
1:C:96:ASP:OD1	1:C:99:ARG:NH1	2.43	0.48
1:B:327:LYS:NZ	4:B:817:HOH:O	2.46	0.48
1:C:230:PRO:HA	1:C:233:TRP:CE2	2.49	0.48
1:C:355:ARG:NH2	4:C:1044:HOH:O	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:GLU:HB2	1:B:67:ARG:HD2	1.97	0.47
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.50	0.47
1:C:63:TRP:CD1	1:C:67:ARG:HG3	2.49	0.47
1:B:63:TRP:CD1	1:B:67:ARG:HG3	2.50	0.46
1:A:322:MET:HA	1:A:322:MET:HE2	1.96	0.46
1:A:56:ASP:CG	1:A:57:GLY:H	2.24	0.46
1:B:41:PRO:HG2	1:B:44:LEU:HD13	1.98	0.46
1:B:259:PHE:CG	1:B:331:MET:HG2	2.50	0.45
1:B:372:LEU:HD23	1:B:377:GLN:HB2	1.99	0.45
1:A:80:ILE:HG22	1:A:82:PRO:HD3	1.99	0.44
1:B:35:LYS:NZ	4:B:1100:HOH:O	2.51	0.44
1:A:47:LYS:HE3	4:A:1089:HOH:O	2.18	0.44
1:A:393:LEU:HD21	1:A:409:VAL:HG11	2.00	0.44
1:C:252:LYS:NZ	4:C:1056:HOH:O	2.50	0.43
1:B:371:ASP:HB3	1:B:431:ARG:NH2	2.34	0.43
1:C:93:PHE:CZ	1:C:111:VAL:HG21	2.53	0.43
1:C:426:SER:HA	1:C:429:TRP:CD2	2.54	0.42
1:C:429:TRP:HB3	1:C:433:CYS:HB2	2.01	0.42
1:A:240:ALA:O	4:A:1056:HOH:O	2.21	0.42
1:A:426:SER:HB2	1:A:429:TRP:CE3	2.54	0.42
1:A:45:GLU:OE1	1:A:45:GLU:N	2.35	0.42
1:C:355:ARG:NH2	1:C:428:SER:OG	2.53	0.42
1:C:29:GLU:OE2	1:C:35:LYS:HD3	2.20	0.41
1:B:350:ASN:OD1	1:B:355:ARG:NH1	2.52	0.41
1:B:90:LEU:HD13	1:B:108:PRO:HG2	2.02	0.41
1:C:259:PHE:CG	1:C:331:MET:HG2	2.56	0.41
1:B:443:GLU:OE2	4:B:992:HOH:O	2.22	0.41
1:B:368:ASN:ND2	1:B:427:GLN:O	2.53	0.41
1:B:123:LEU:HD21	1:B:127:PRO:HD3	2.01	0.41
1:B:56:ASP:OD1	1:B:57:GLY:N	2.48	0.40
1:B:252:LYS:NZ	4:B:1121:HOH:O	2.53	0.40
1:C:169:ALA:O	1:C:182:VAL:HA	2.21	0.40
1:A:271:SER:O	1:A:274:LYS:NZ	2.50	0.40
1:B:240:ALA:O	4:B:1050:HOH:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	460/466 (99%)	449 (98%)	11 (2%)	0	100	100
1	B	459/466 (98%)	449 (98%)	10 (2%)	0	100	100
1	C	461/466 (99%)	453 (98%)	8 (2%)	0	100	100
All	All	1380/1398 (99%)	1351 (98%)	29 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	372/376 (99%)	366 (98%)	6 (2%)	55	62
1	B	371/376 (99%)	363 (98%)	8 (2%)	45	50
1	C	373/376 (99%)	365 (98%)	8 (2%)	47	52
All	All	1116/1128 (99%)	1094 (98%)	22 (2%)	55	54

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

Mol	Chain	Res	Type
1	A	239[A]	SER
1	A	239[B]	SER
1	A	259	PHE
1	A	372	LEU
1	A	400	VAL

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Mol	Chain	Res	Type
1	A	431	ARG
1	B	30	LYS
1	B	88[A]	ASP
1	B	88[B]	ASP
1	B	239[A]	SER
1	B	239[B]	SER
1	B	259	PHE
1	B	372	LEU
1	B	399	GLN
1	C	239[A]	SER
1	C	239[B]	SER
1	C	259	PHE
1	C	372	LEU
1	C	377	GLN
1	C	399	GLN
1	C	426	SER
1	C	432	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	HIS
1	B	40	HIS
1	B	283	ASN
1	B	427	GLN
1	B	447	HIS
1	C	405	GLN
1	C	447	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

6 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	GLC	D	1	2	12,12,12	2.46	6 (50%)	17,17,17	1.08	2 (11%)
2	GLC	D	2	2	11,11,12	1.50	2 (18%)	15,15,17	0.76	0
2	GLC	E	1	2	12,12,12	2.41	6 (50%)	17,17,17	1.18	2 (11%)
2	GLC	E	2	2	11,11,12	1.57	2 (18%)	15,15,17	0.71	0
2	GLC	F	1	2	12,12,12	2.43	6 (50%)	17,17,17	1.29	3 (17%)
2	GLC	F	2	2	11,11,12	1.53	2 (18%)	15,15,17	0.68	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	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	0/2/19/22	0/1/1/1
2	GLC	E	1	2	-	0/2/22/22	0/1/1/1
2	GLC	E	2	2	-	0/2/19/22	0/1/1/1
2	GLC	F	1	2	-	0/2/22/22	0/1/1/1
2	GLC	F	2	2	-	0/2/19/22	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	GLC	C6-C5	-4.52	1.36	1.51
2	E	1	GLC	C6-C5	-4.43	1.36	1.51
2	F	1	GLC	C6-C5	-4.41	1.37	1.51
2	E	1	GLC	C3-C2	-3.99	1.42	1.52
2	D	1	GLC	C3-C2	-3.91	1.42	1.52
2	F	1	GLC	C3-C2	-3.86	1.42	1.52
2	E	2	GLC	O5-C1	3.13	1.49	1.43
2	D	1	GLC	O5-C5	3.09	1.51	1.44
2	F	2	GLC	O5-C1	3.03	1.48	1.43
2	D	2	GLC	O5-C1	2.98	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1	GLC	O5-C5	2.97	1.51	1.44
2	D	1	GLC	O2-C2	2.88	1.50	1.43
2	E	1	GLC	O5-C5	2.86	1.51	1.44
2	F	1	GLC	O2-C2	2.86	1.50	1.43
2	E	1	GLC	O2-C2	2.82	1.49	1.43
2	F	1	GLC	C1-C2	-2.71	1.46	1.52
2	F	2	GLC	C2-C3	-2.62	1.48	1.52
2	D	1	GLC	C1-C2	-2.60	1.46	1.52
2	E	1	GLC	C1-C2	-2.58	1.46	1.52
2	E	2	GLC	C2-C3	-2.55	1.48	1.52
2	F	1	GLC	O3-C3	2.33	1.48	1.43
2	D	2	GLC	C2-C3	-2.33	1.48	1.52
2	D	1	GLC	O3-C3	2.19	1.48	1.43
2	E	1	GLC	O3-C3	2.09	1.48	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	GLC	C1-C2-C3	2.62	115.70	110.36
2	E	1	GLC	C1-C2-C3	2.50	115.45	110.36
2	F	1	GLC	C4-C3-C2	2.43	115.09	110.83
2	D	1	GLC	C4-C3-C2	2.32	114.90	110.83
2	F	1	GLC	O5-C1-C2	2.27	114.29	110.30
2	D	1	GLC	C1-C2-C3	2.19	114.83	110.36
2	E	1	GLC	C4-C3-C2	2.19	114.67	110.83

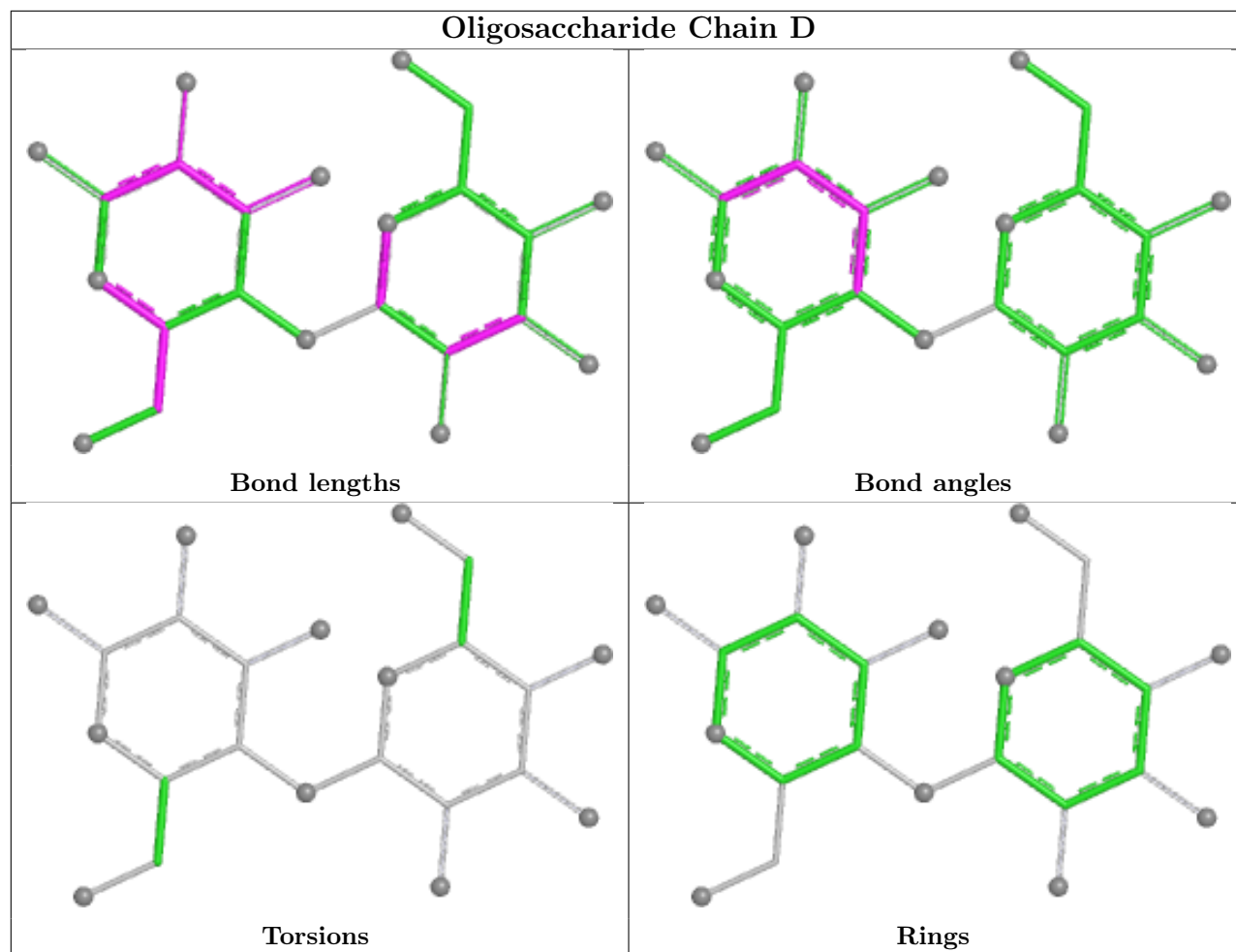
There are no chirality outliers.

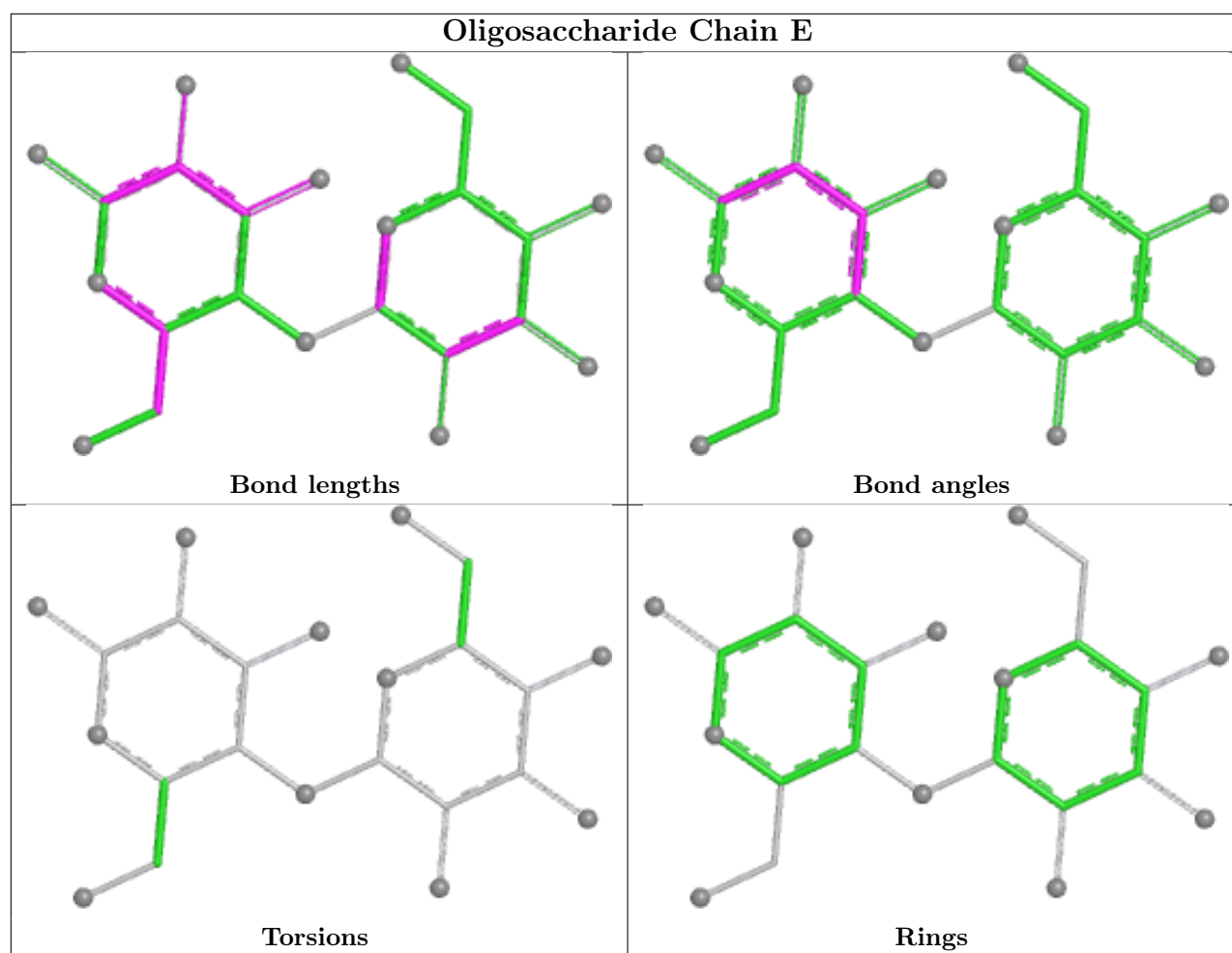
There are no torsion outliers.

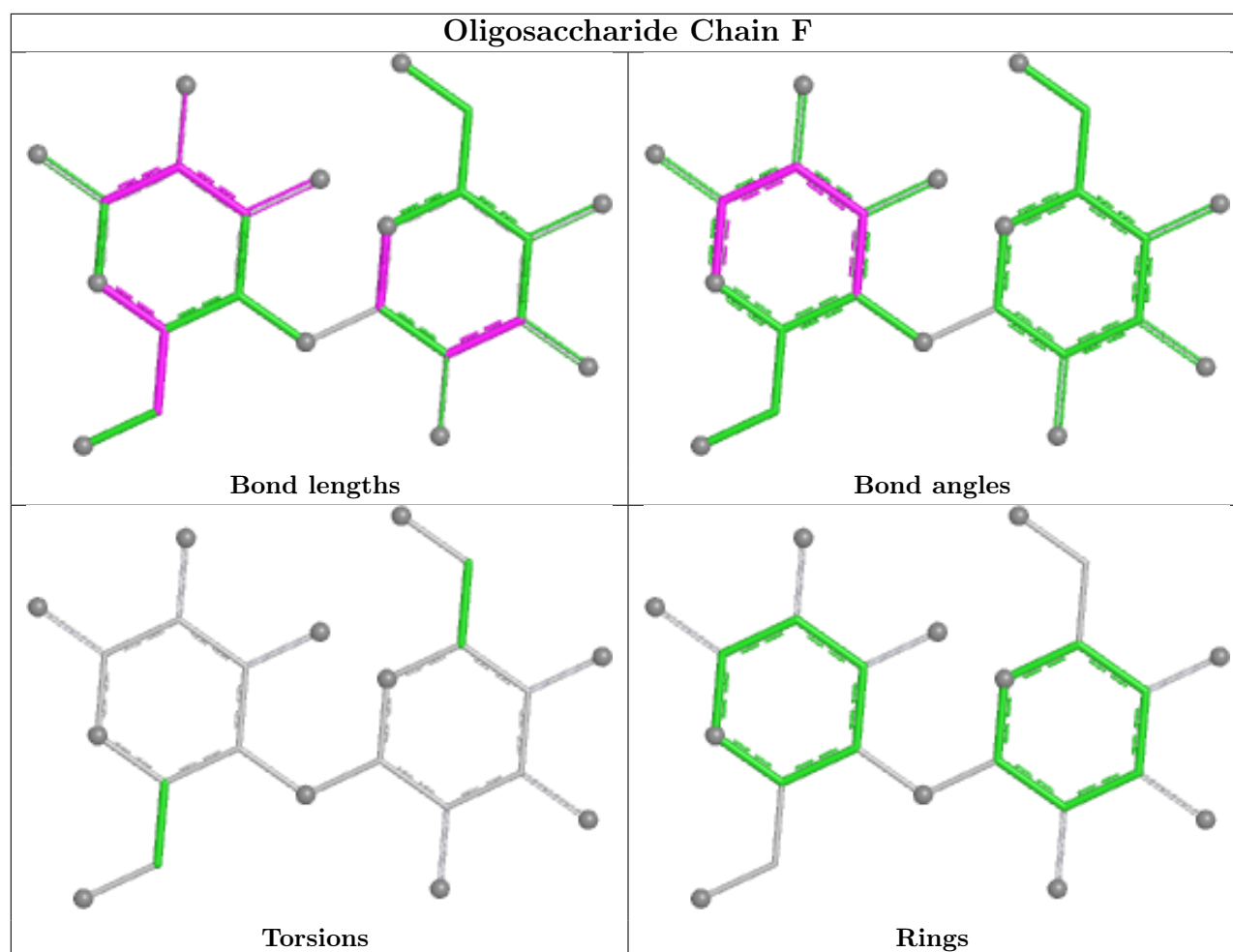
There are no ring outliers.

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](#)

3 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	$\# Z > 2$
3	MLI	A	502	-	6,6,6	1.29	0	7,7,7	1.19	0
3	MLI	C	502	-	6,6,6	1.19	0	7,7,7	1.41	0
3	MLI	B	502	-	6,6,6	1.30	0	7,7,7	1.27	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MLI	A	502	-	-	2/4/4/4	-
3	MLI	C	502	-	-	2/4/4/4	-
3	MLI	B	502	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	MLI	C3-C1-C2-O6
3	C	502	MLI	C2-C1-C3-O9
3	C	502	MLI	C2-C1-C3-O8
3	A	502	MLI	C3-C1-C2-O7

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	458/466 (98%)	-1.30	0 100 100	16, 28, 52, 80	4 (0%)
1	B	458/466 (98%)	-1.29	0 100 100	16, 29, 53, 79	3 (0%)
1	C	458/466 (98%)	-1.30	0 100 100	15, 29, 53, 75	5 (1%)
All	All	1374/1398 (98%)	-1.30	0 100 100	15, 29, 52, 80	12 (0%)

There are no RSRZ outliers to report.

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](#)

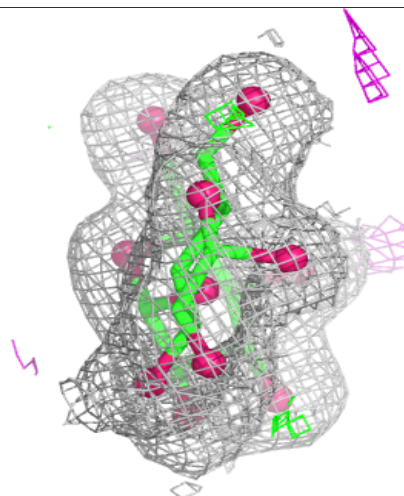
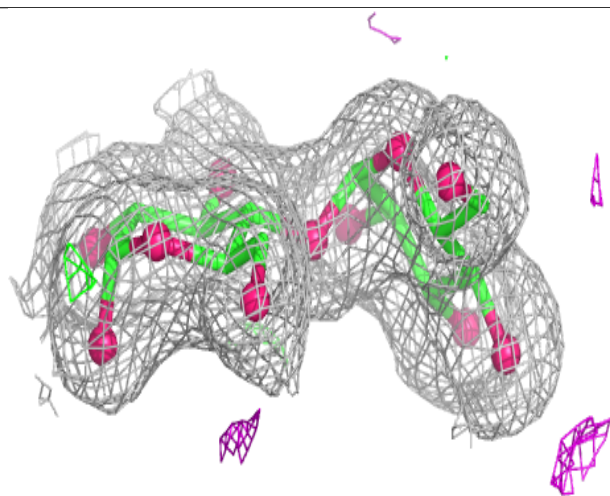
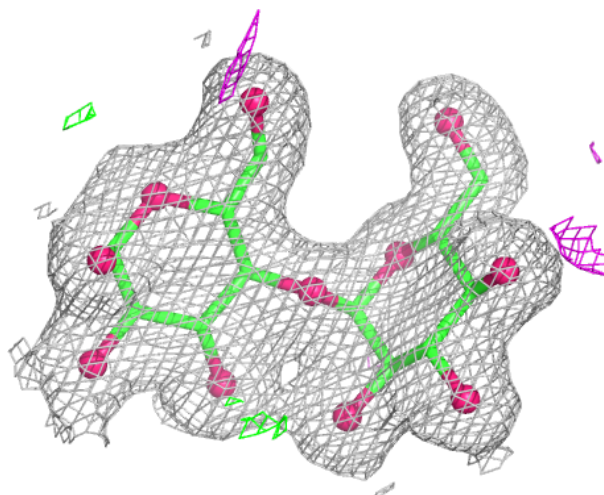
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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	GLC	F	1	12/12	0.99	0.02	18,24,29,31	0
2	GLC	F	2	11/12	0.99	0.02	18,22,24,25	0
2	GLC	E	1	12/12	1.00	0.02	19,23,30,32	0
2	GLC	E	2	11/12	1.00	0.02	17,20,23,24	0
2	GLC	D	1	12/12	1.00	0.02	17,24,28,30	0
2	GLC	D	2	11/12	1.00	0.02	18,20,23,24	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.

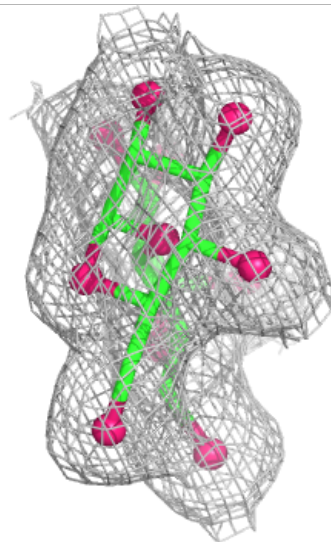
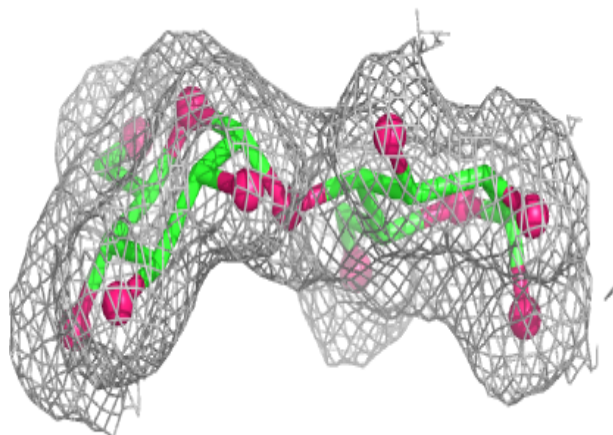
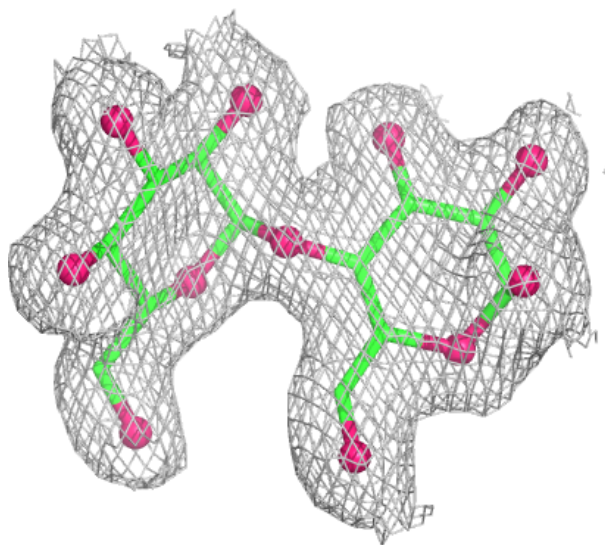
Electron density around Chain D:

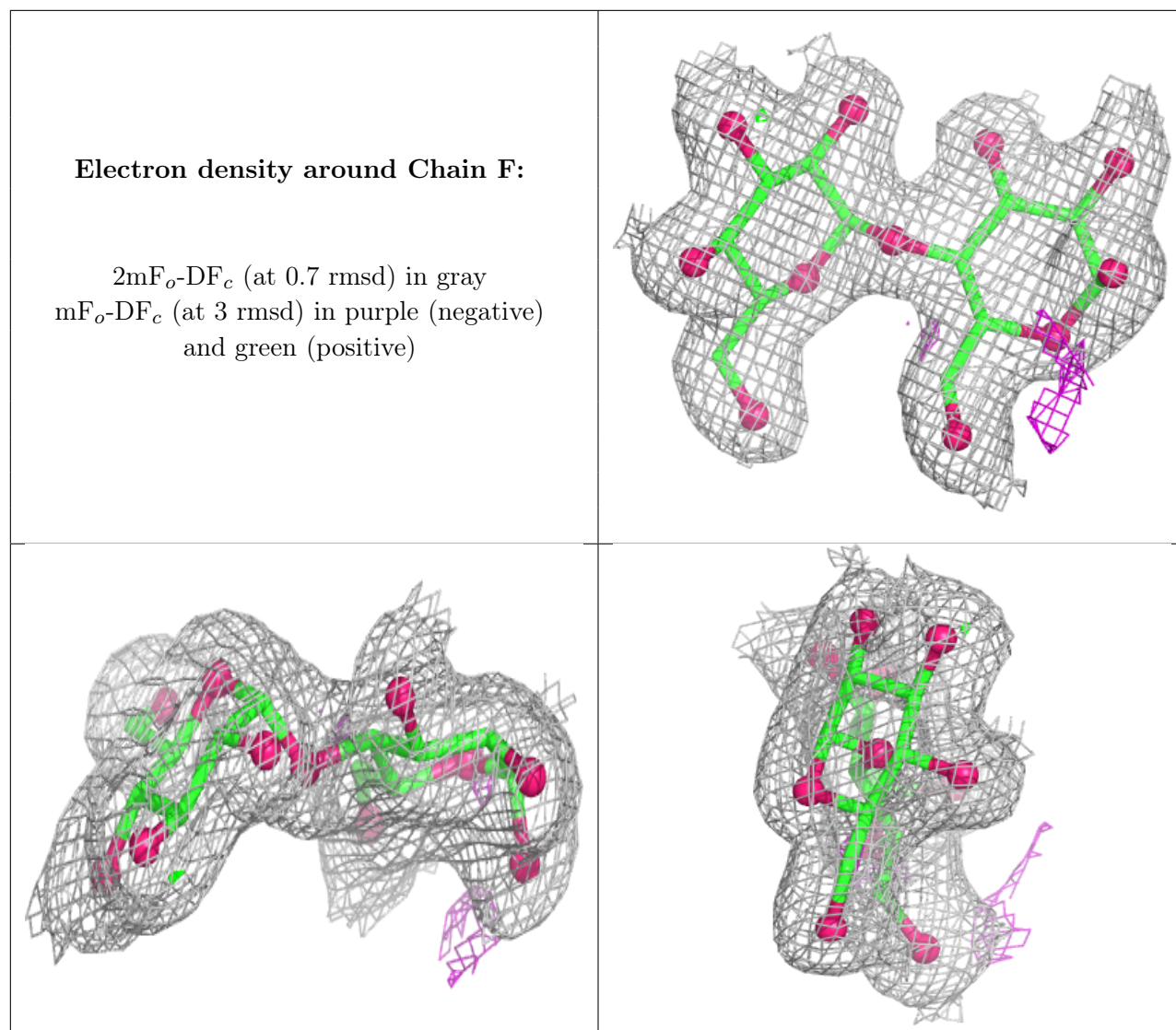
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MLI	A	502	7/7	0.99	0.03	28,31,47,57	0
3	MLI	B	502	7/7	0.99	0.03	28,30,50,61	0
3	MLI	C	502	7/7	0.99	0.03	29,32,54,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.