



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 06:28 PM UTC

PDB ID : 5V6Y / pdb_00005v6y
Title : Crystal structure of the human CLR:RAMP1 extracellular domain heterodimer with bound high-affinity and altered selectivity adrenomedullin variant
Authors : Pioszak, A.; Booe, J.
Deposited on : 2017-03-17
Resolution : 2.80 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 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

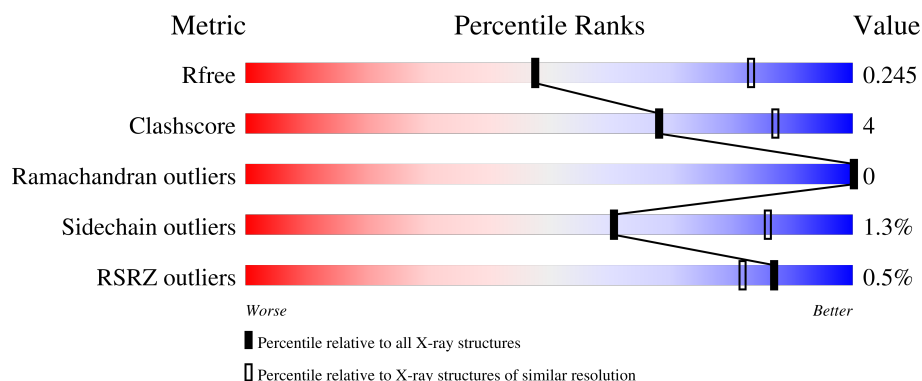
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	593	85% 7% 7%
1	B	593	83% 10% 7%
1	C	593	82% 10% 7%
1	D	593	83% 9% 8%
2	E	17	76% 18% 6%

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Mol	Chain	Length	Quality of chain
2	F	17	6% 88% 6% 6%
2	G	17	76% 6% 18%
2	H	17	59% 24% 18%
3	I	2	50% 50%
3	J	2	100%
3	K	2	50% 50%
3	L	2	50% 50%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17855 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, Receptor activity-modifying protein 1, Calcitonin gene-related peptide type 1 receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	550	Total	C	N	O	S	0	0	0
			4341	2773	720	825	23			
1	B	551	Total	C	N	O	S	0	0	0
			4349	2779	721	826	23			
1	C	549	Total	C	N	O	S	0	0	0
			4335	2770	718	824	23			
1	D	547	Total	C	N	O	S	0	0	0
			4278	2727	710	818	23			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P0AEY0
A	369	ASN	-	linker	UNP P0AEY0
A	370	ALA	-	linker	UNP P0AEY0
A	371	ALA	-	linker	UNP P0AEY0
A	372	ALA	-	linker	UNP P0AEY0
A	373	GLU	-	linker	UNP P0AEY0
A	374	PHE	-	linker	UNP P0AEY0
A	2020	GLY	-	linker	UNP O60894
A	2021	SER	-	linker	UNP O60894
A	2022	ALA	-	linker	UNP O60894
A	2023	GLY	-	linker	UNP O60894
A	2024	SER	-	linker	UNP O60894
A	2025	ALA	-	linker	UNP O60894
A	2026	GLY	-	linker	UNP O60894
A	2027	SER	-	linker	UNP O60894
A	2028	ALA	-	linker	UNP O60894
A	2145	HIS	-	expression tag	UNP Q16602
A	2146	HIS	-	expression tag	UNP Q16602
A	2147	HIS	-	expression tag	UNP Q16602
A	2148	HIS	-	expression tag	UNP Q16602

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Chain	Residue	Modelled	Actual	Comment	Reference
A	2149	HIS	-	expression tag	UNP Q16602
A	2150	HIS	-	expression tag	UNP Q16602
B	1	MET	-	initiating methionine	UNP P0AEY0
B	369	ASN	-	linker	UNP P0AEY0
B	370	ALA	-	linker	UNP P0AEY0
B	371	ALA	-	linker	UNP P0AEY0
B	372	ALA	-	linker	UNP P0AEY0
B	373	GLU	-	linker	UNP P0AEY0
B	374	PHE	-	linker	UNP P0AEY0
B	2020	GLY	-	linker	UNP O60894
B	2021	SER	-	linker	UNP O60894
B	2022	ALA	-	linker	UNP O60894
B	2023	GLY	-	linker	UNP O60894
B	2024	SER	-	linker	UNP O60894
B	2025	ALA	-	linker	UNP O60894
B	2026	GLY	-	linker	UNP O60894
B	2027	SER	-	linker	UNP O60894
B	2028	ALA	-	linker	UNP O60894
B	2145	HIS	-	expression tag	UNP Q16602
B	2146	HIS	-	expression tag	UNP Q16602
B	2147	HIS	-	expression tag	UNP Q16602
B	2148	HIS	-	expression tag	UNP Q16602
B	2149	HIS	-	expression tag	UNP Q16602
B	2150	HIS	-	expression tag	UNP Q16602
C	1	MET	-	initiating methionine	UNP P0AEY0
C	369	ASN	-	linker	UNP P0AEY0
C	370	ALA	-	linker	UNP P0AEY0
C	371	ALA	-	linker	UNP P0AEY0
C	372	ALA	-	linker	UNP P0AEY0
C	373	GLU	-	linker	UNP P0AEY0
C	374	PHE	-	linker	UNP P0AEY0
C	2020	GLY	-	linker	UNP O60894
C	2021	SER	-	linker	UNP O60894
C	2022	ALA	-	linker	UNP O60894
C	2023	GLY	-	linker	UNP O60894
C	2024	SER	-	linker	UNP O60894
C	2025	ALA	-	linker	UNP O60894
C	2026	GLY	-	linker	UNP O60894
C	2027	SER	-	linker	UNP O60894
C	2028	ALA	-	linker	UNP O60894
C	2145	HIS	-	expression tag	UNP Q16602
C	2146	HIS	-	expression tag	UNP Q16602

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Chain	Residue	Modelled	Actual	Comment	Reference
C	2147	HIS	-	expression tag	UNP Q16602
C	2148	HIS	-	expression tag	UNP Q16602
C	2149	HIS	-	expression tag	UNP Q16602
C	2150	HIS	-	expression tag	UNP Q16602
D	1	MET	-	initiating methionine	UNP P0AEY0
D	369	ASN	-	linker	UNP P0AEY0
D	370	ALA	-	linker	UNP P0AEY0
D	371	ALA	-	linker	UNP P0AEY0
D	372	ALA	-	linker	UNP P0AEY0
D	373	GLU	-	linker	UNP P0AEY0
D	374	PHE	-	linker	UNP P0AEY0
D	2020	GLY	-	linker	UNP O60894
D	2021	SER	-	linker	UNP O60894
D	2022	ALA	-	linker	UNP O60894
D	2023	GLY	-	linker	UNP O60894
D	2024	SER	-	linker	UNP O60894
D	2025	ALA	-	linker	UNP O60894
D	2026	GLY	-	linker	UNP O60894
D	2027	SER	-	linker	UNP O60894
D	2028	ALA	-	linker	UNP O60894
D	2145	HIS	-	expression tag	UNP Q16602
D	2146	HIS	-	expression tag	UNP Q16602
D	2147	HIS	-	expression tag	UNP Q16602
D	2148	HIS	-	expression tag	UNP Q16602
D	2149	HIS	-	expression tag	UNP Q16602
D	2150	HIS	-	expression tag	UNP Q16602

- Molecule 2 is a protein called ADM.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	16	Total	C	N	O	0	0	1
			118	80	19	19			
2	F	16	Total	C	N	O	0	1	1
			138	94	24	20			
2	G	14	Total	C	N	O	0	0	1
			111	76	20	15			
2	H	14	Total	C	N	O	0	0	1
			90	61	15	14			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	45	TRP	SER	engineered mutation	UNP P35318
E	46	LEU	LYS	engineered mutation	UNP P35318
E	50	TRP	GLN	engineered mutation	UNP P35318
E	52	PHE	TYR	engineered mutation	UNP P35318
E	53	NH2	-	amidation	UNP P35318
F	45	TRP	SER	engineered mutation	UNP P35318
F	46	LEU	LYS	engineered mutation	UNP P35318
F	50	TRP	GLN	engineered mutation	UNP P35318
F	52	PHE	TYR	engineered mutation	UNP P35318
F	53	NH2	-	amidation	UNP P35318
G	45	TRP	SER	engineered mutation	UNP P35318
G	46	LEU	LYS	engineered mutation	UNP P35318
G	50	TRP	GLN	engineered mutation	UNP P35318
G	52	PHE	TYR	engineered mutation	UNP P35318
G	53	NH2	-	amidation	UNP P35318
H	45	TRP	SER	engineered mutation	UNP P35318
H	46	LEU	LYS	engineered mutation	UNP P35318
H	50	TRP	GLN	engineered mutation	UNP P35318
H	52	PHE	TYR	engineered mutation	UNP P35318
H	53	NH2	-	amidation	UNP P35318

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	I	2	Total	C	O	0	0	0
			23	12	11			
3	J	2	Total	C	O	0	0	0
			23	12	11			
3	K	2	Total	C	O	0	0	0
			23	12	11			
3	L	2	Total	C	O	0	0	0
			23	12	11			

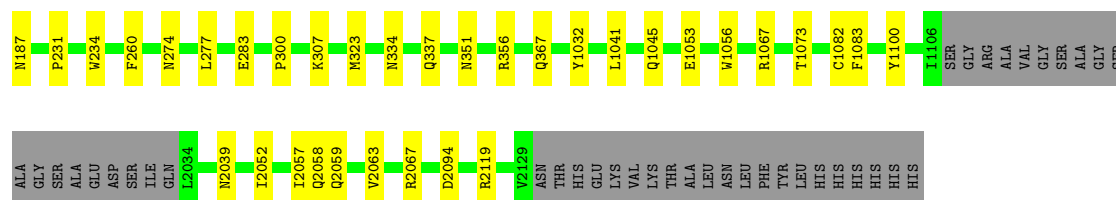
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		

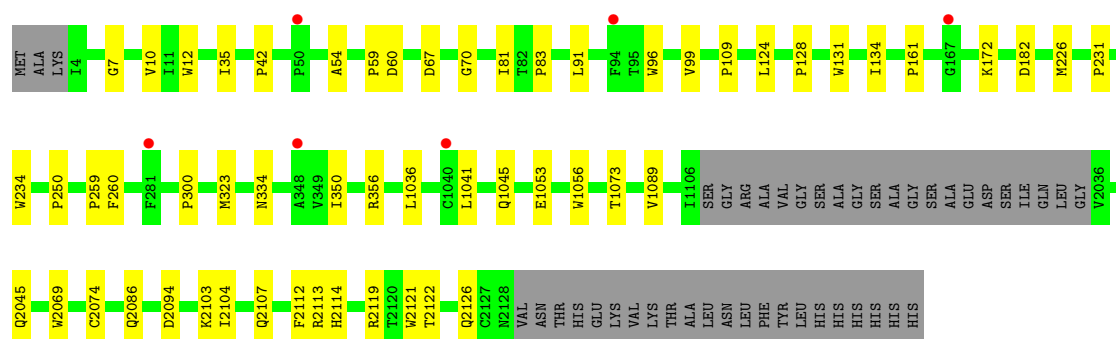
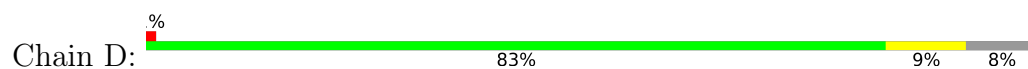
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	O	0	0
			1	1		



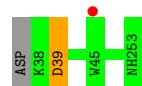
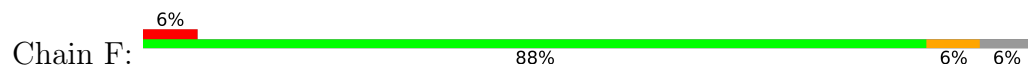
- Molecule 1: Maltose-binding periplasmic protein, Receptor activity-modifying protein 1, Calcitonin gene-related peptide type 1 receptor



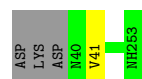
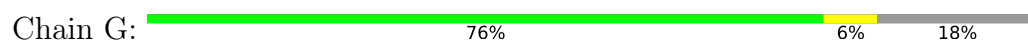
- Molecule 2: ADM



- Molecule 2: ADM



- Molecule 2: ADM



- Molecule 2: ADM

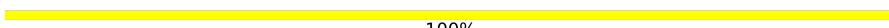


- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain I:  50% 50%



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain J:  100%



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain K:  50% 50%



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain L:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	195.50Å 69.44Å 200.00Å 90.00° 99.43° 90.00°	Depositor
Resolution (Å)	45.11 – 2.80 45.11 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.1 (45.11-2.80) 87.7 (45.11-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.212 , 0.244 0.212 , 0.245	Depositor DCC
R_{free} test set	2000 reflections (3.23%)	wwPDB-VP
Wilson B-factor (Å ²)	80.0	Xtriage
Anisotropy	0.475	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 65.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17855	wwPDB-VP
Average B, all atoms (Å ²)	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.86% 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: NH2, 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	0.09	0/4453	0.26	0/6050
1	B	0.09	0/4461	0.26	0/6061
1	C	0.08	0/4447	0.25	0/6043
1	D	0.09	0/4385	0.25	0/5961
2	E	0.08	0/123	0.27	0/171
2	F	0.10	0/145	0.26	0/201
2	G	0.13	0/116	0.31	0/160
2	H	0.10	0/93	0.32	0/129
All	All	0.09	0/18223	0.26	0/24776

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4341	0	4178	24	1
1	B	4349	0	4186	33	1
1	C	4335	0	4171	33	1
1	D	4278	0	4095	32	0
2	E	118	0	100	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	138	0	120	1	0
2	G	111	0	105	1	0
2	H	90	0	73	2	0
3	I	23	0	21	0	0
3	J	23	0	21	0	0
3	K	23	0	21	0	0
3	L	23	0	21	1	0
4	A	2	0	0	0	0
4	B	1	0	0	0	0
All	All	17855	0	17112	126	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

The worst 5 of 126 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:97:ASP:OD1	1:C:100:ARG:NH1	2.00	0.95
1:B:80:GLU:OE2	1:B:104:LYS:NZ	2.09	0.85
1:C:1067:ARG:NH1	1:C:2039:ASN:OD1	2.09	0.85
1:B:97:ASP:OD1	1:B:100:ARG:NH1	2.24	0.69
1:C:274:ASN:HB3	1:C:277:LEU:HB2	1.78	0.64

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:LYS:NZ	1:B:179:ASP:OD2[4_655]	1.80	0.40
1:A:1050:ALA:O	1:C:36:LYS:NZ[3_545]	1.97	0.23

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	546/593 (92%)	535 (98%)	11 (2%)	0	100	100
1	B	547/593 (92%)	534 (98%)	13 (2%)	0	100	100
1	C	545/593 (92%)	532 (98%)	13 (2%)	0	100	100
1	D	543/593 (92%)	529 (97%)	14 (3%)	0	100	100
2	E	14/17 (82%)	14 (100%)	0	0	100	100
2	F	15/17 (88%)	14 (93%)	1 (7%)	0	100	100
2	G	12/17 (71%)	12 (100%)	0	0	100	100
2	H	12/17 (71%)	12 (100%)	0	0	100	100
All	All	2234/2440 (92%)	2182 (98%)	52 (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	450/484 (93%)	444 (99%)	6 (1%)	61	86
1	B	451/484 (93%)	447 (99%)	4 (1%)	70	89
1	C	450/484 (93%)	441 (98%)	9 (2%)	48	80
1	D	440/484 (91%)	437 (99%)	3 (1%)	76	91
2	E	11/14 (79%)	11 (100%)	0	100	100
2	F	13/14 (93%)	12 (92%)	1 (8%)	12	36
2	G	11/14 (79%)	11 (100%)	0	100	100
2	H	7/14 (50%)	7 (100%)	0	100	100
All	All	1833/1992 (92%)	1810 (99%)	23 (1%)	61	86

5 of 23 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	260	PHE
1	C	2059	GLN

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Mol	Chain	Res	Type
1	C	2058	GLN
1	C	2063	VAL
1	B	5	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	2045	GLN
1	D	284	ASN
2	F	40	ASN
1	D	367	GLN
1	D	203	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 ⓘ

8 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GLC	I	1	3	12,12,12	0.77	0	17,17,17	0.82	1 (5%)
3	GLC	I	2	3	11,11,12	0.95	0	15,15,17	0.78	0
3	GLC	J	1	3	12,12,12	0.78	0	17,17,17	0.83	1 (5%)
3	GLC	J	2	3	11,11,12	0.91	0	15,15,17	0.73	1 (6%)
3	GLC	K	1	3	12,12,12	0.77	0	17,17,17	0.84	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GLC	K	2	3	11,11,12	0.92	0	15,15,17	0.81	0
3	GLC	L	1	3	12,12,12	0.77	0	17,17,17	0.82	1 (5%)
3	GLC	L	2	3	11,11,12	0.97	0	15,15,17	0.81	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
3	GLC	I	1	3	-	0/2/22/22	0/1/1/1
3	GLC	I	2	3	-	0/2/19/22	0/1/1/1
3	GLC	J	1	3	-	0/2/22/22	0/1/1/1
3	GLC	J	2	3	-	0/2/19/22	0/1/1/1
3	GLC	K	1	3	-	0/2/22/22	0/1/1/1
3	GLC	K	2	3	-	0/2/19/22	0/1/1/1
3	GLC	L	1	3	-	0/2/22/22	0/1/1/1
3	GLC	L	2	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	1	GLC	C1-O5-C5	-2.54	108.74	113.65
3	I	1	GLC	C1-O5-C5	-2.52	108.77	113.65
3	J	1	GLC	C1-O5-C5	-2.50	108.81	113.65
3	K	1	GLC	C1-O5-C5	-2.49	108.83	113.65
3	L	2	GLC	C1-O5-C5	-2.08	109.39	112.19

There are no chirality outliers.

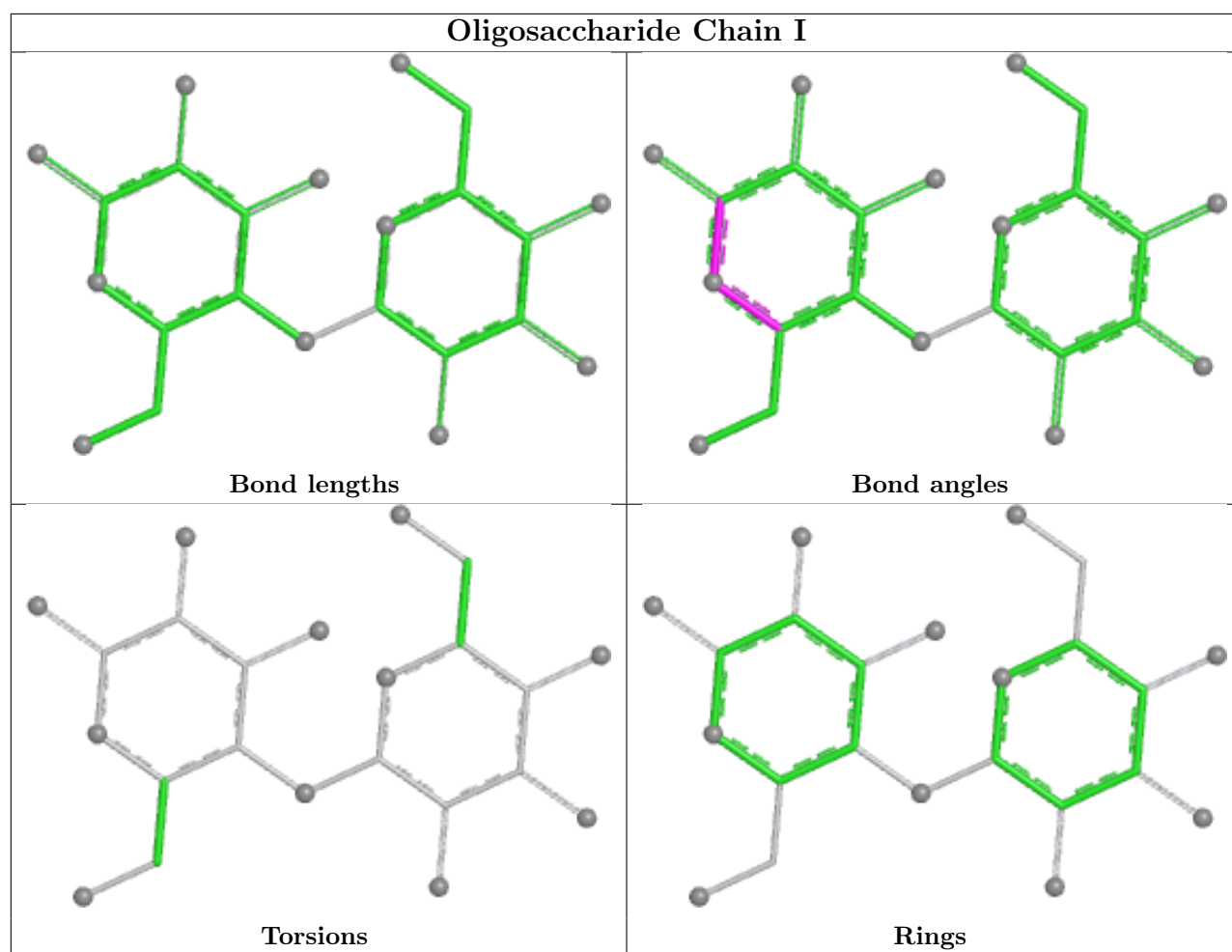
There are no torsion outliers.

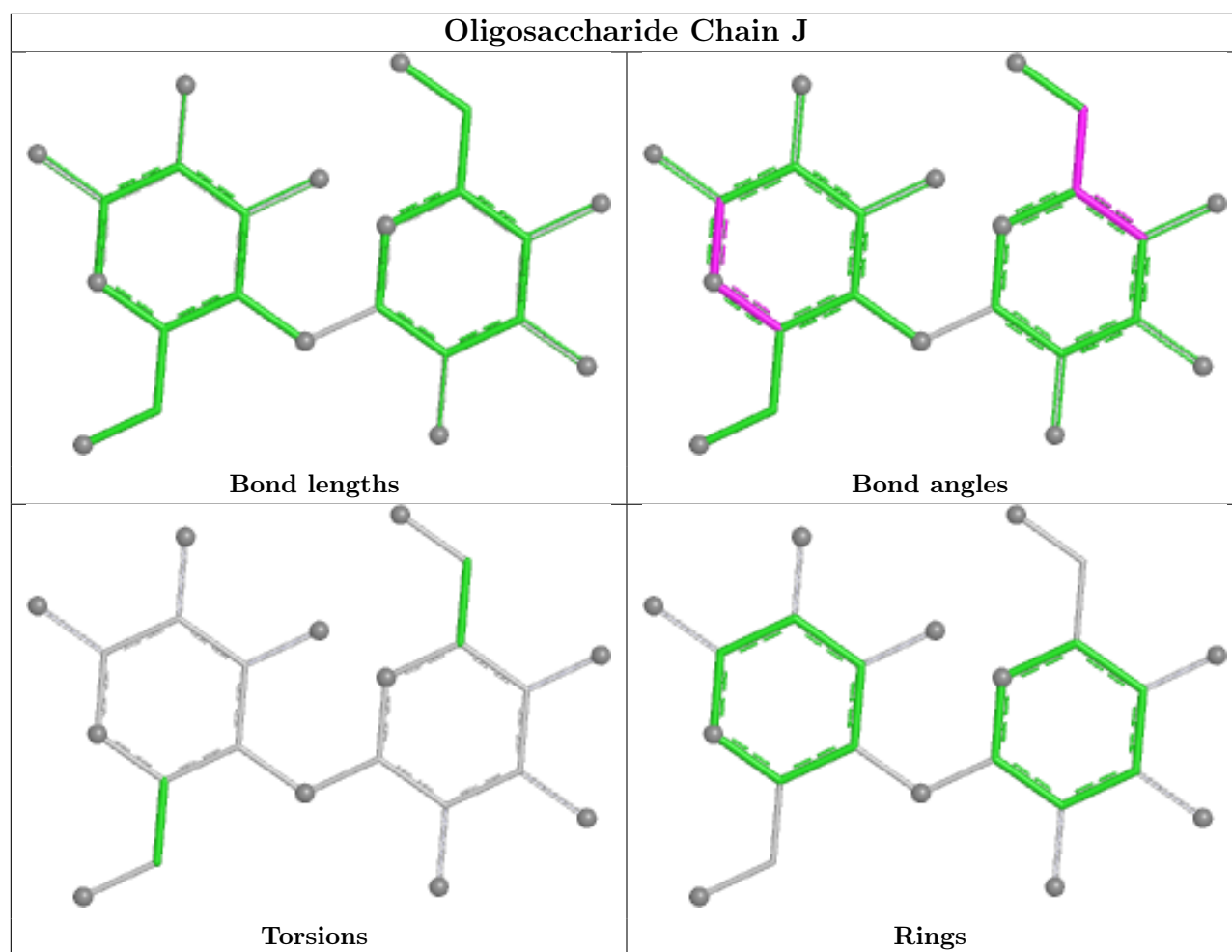
There are no ring outliers.

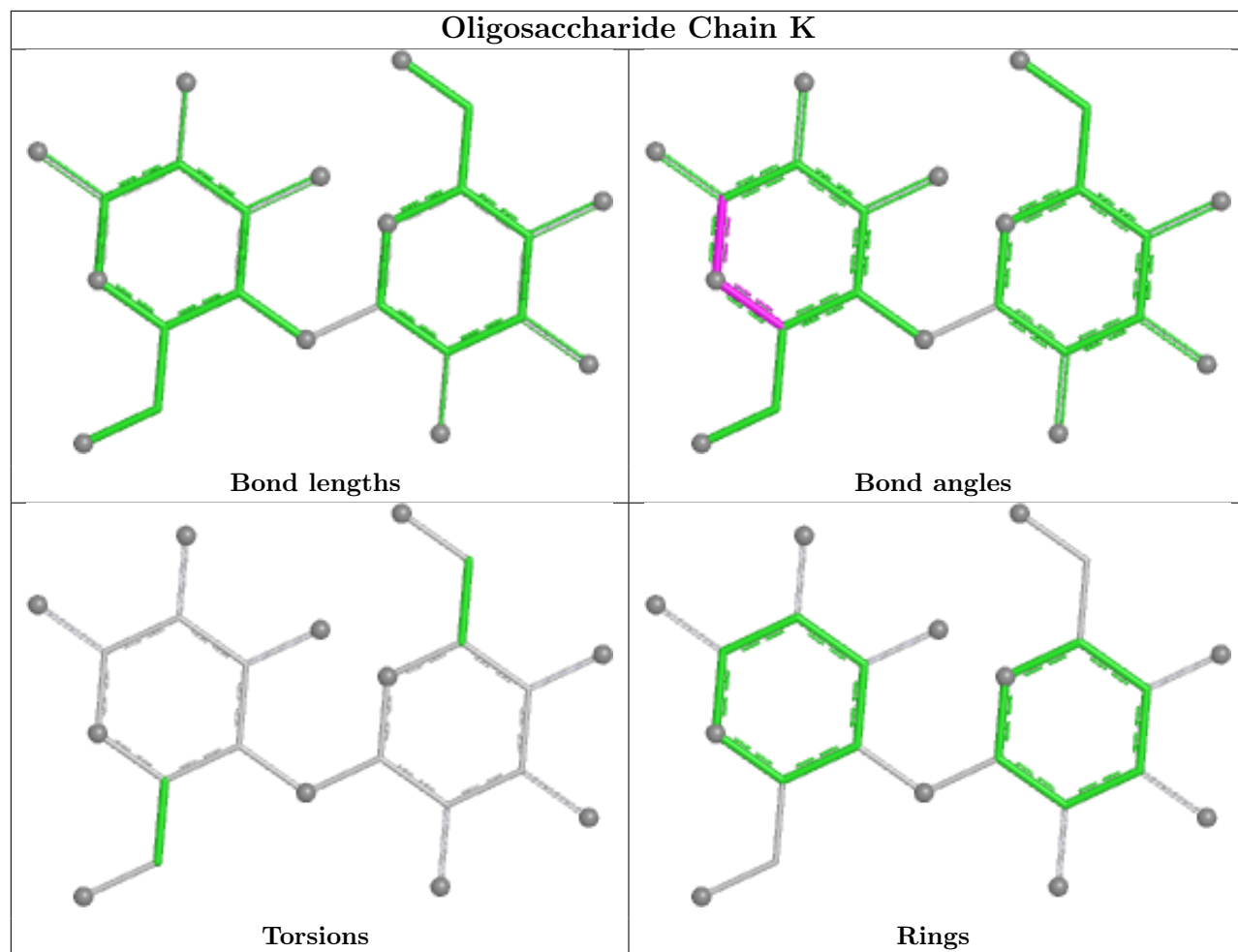
1 monomer is involved in 1 short contact:

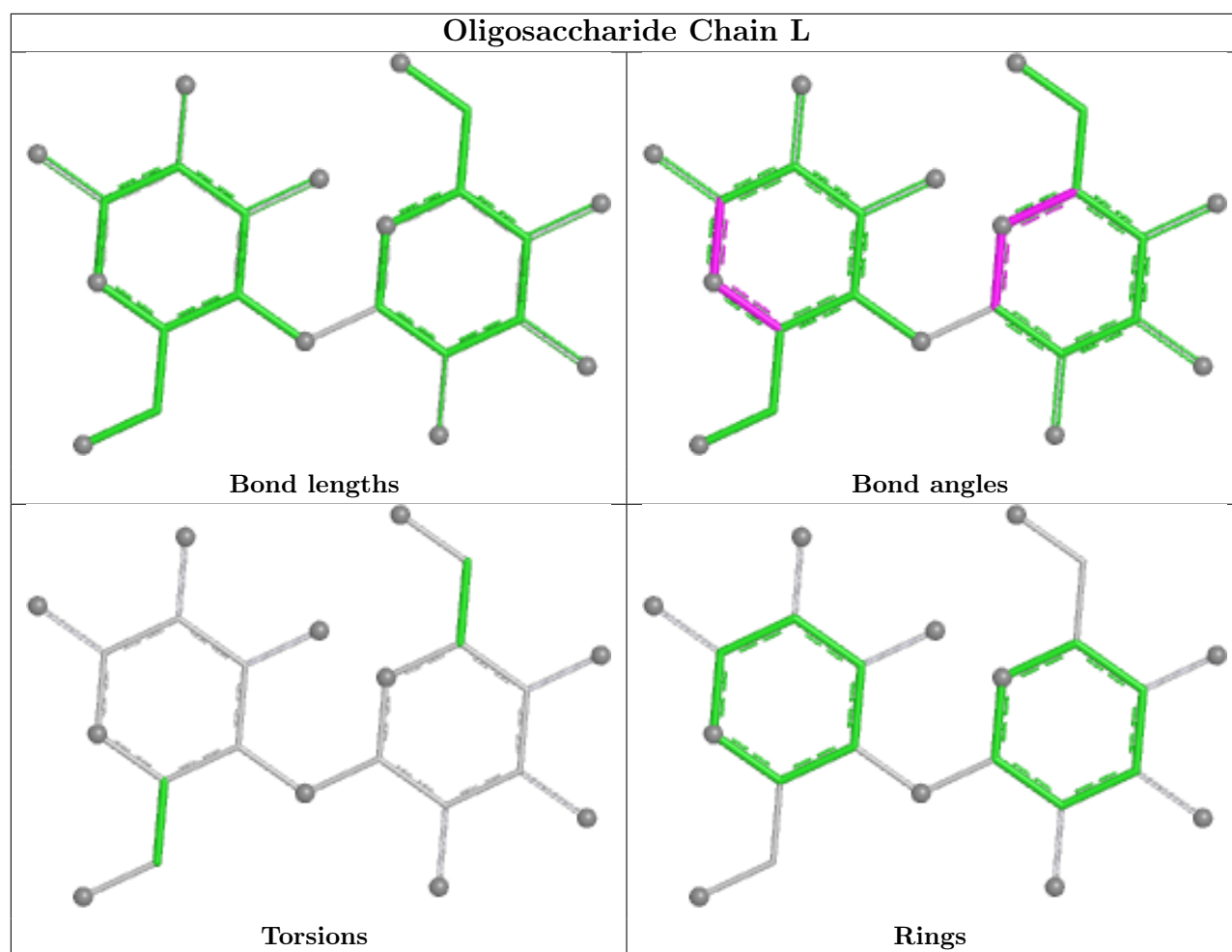
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	2	GLC	1	0

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









5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	550/593 (92%)	-0.32	2 (0%) 88 84	71, 108, 156, 199	0
1	B	551/593 (92%)	-0.29	2 (0%) 88 84	58, 111, 180, 214	0
1	C	549/593 (92%)	-0.31	1 (0%) 91 88	58, 103, 161, 211	0
1	D	547/593 (92%)	-0.14	6 (1%) 78 70	82, 138, 186, 246	0
2	E	15/17 (88%)	-0.08	0 100 100	97, 113, 175, 192	0
2	F	15/17 (88%)	0.03	1 (6%) 24 17	56, 80, 144, 148	1 (6%)
2	G	13/17 (76%)	-0.08	0 100 100	55, 67, 128, 148	0
2	H	13/17 (76%)	0.28	0 100 100	145, 177, 218, 254	0
All	All	2253/2440 (92%)	-0.26	12 (0%) 87 82	55, 115, 175, 254	1 (0%)

The worst 5 of 12 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	348	ALA	4.1
1	B	369	ASN	3.6
1	D	50	PRO	3.0
1	B	255	GLN	2.9
1	C	165	ALA	2.7

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

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

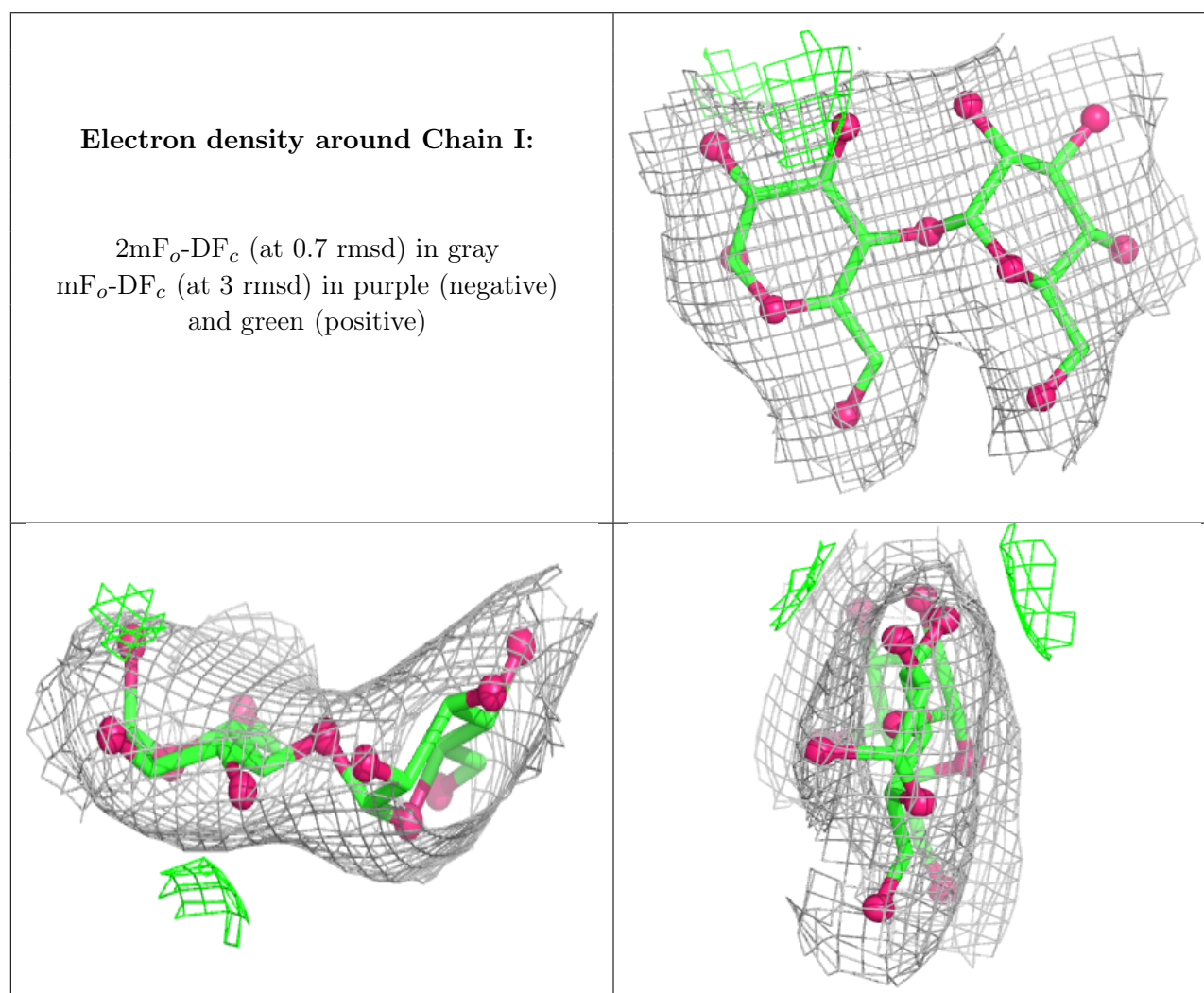
6.3 Carbohydrates [i](#)

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.

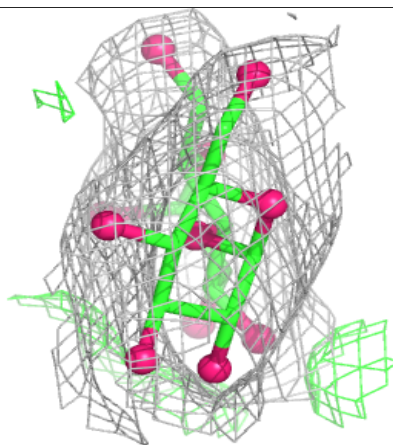
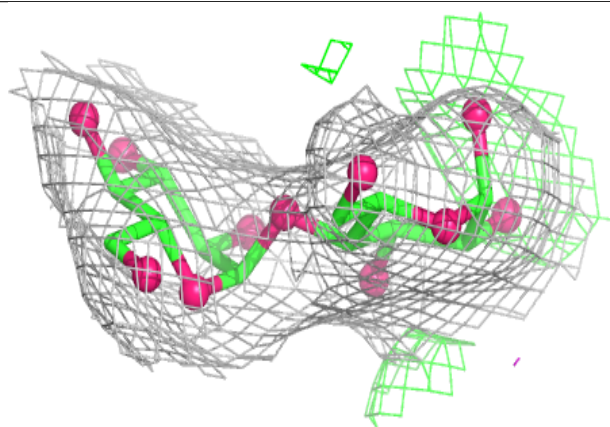
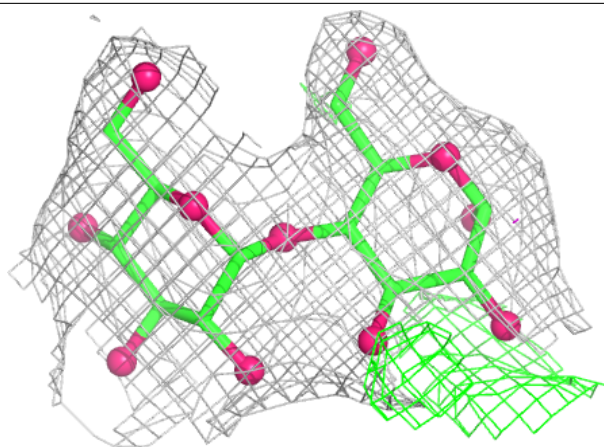
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GLC	J	1	12/12	0.91	0.08	67,79,86,92	0
3	GLC	K	1	12/12	0.94	0.10	72,78,86,89	0
3	GLC	L	1	12/12	0.94	0.07	95,102,110,113	0
3	GLC	I	1	12/12	0.95	0.06	72,76,85,86	0
3	GLC	K	2	11/12	0.97	0.06	61,65,69,70	0
3	GLC	L	2	11/12	0.97	0.07	90,93,103,105	0
3	GLC	I	2	11/12	0.98	0.04	70,73,78,81	0
3	GLC	J	2	11/12	0.98	0.05	60,63,67,73	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.



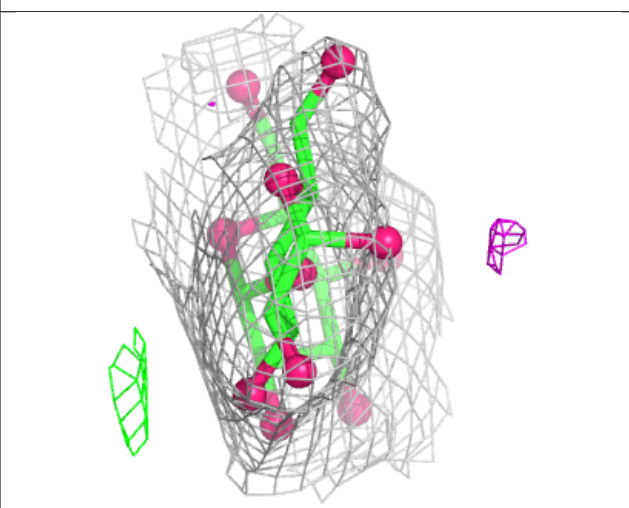
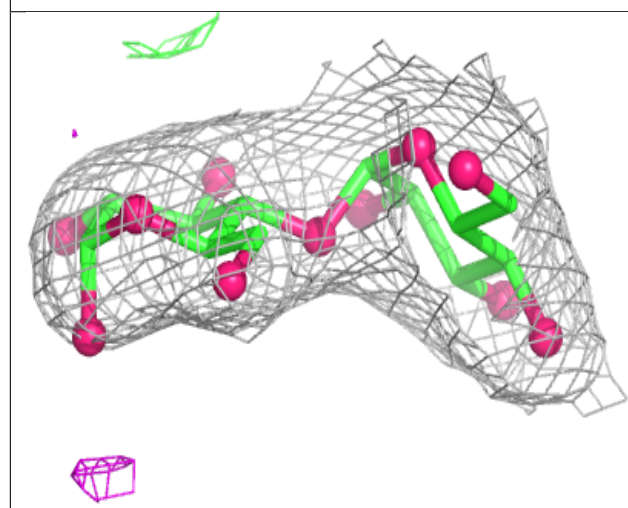
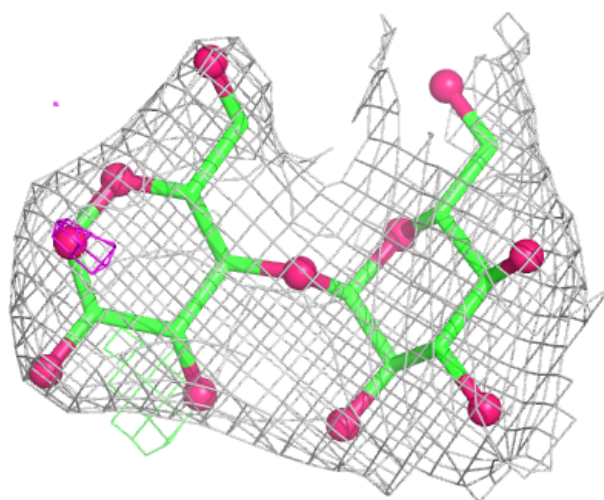
Electron density around Chain J:

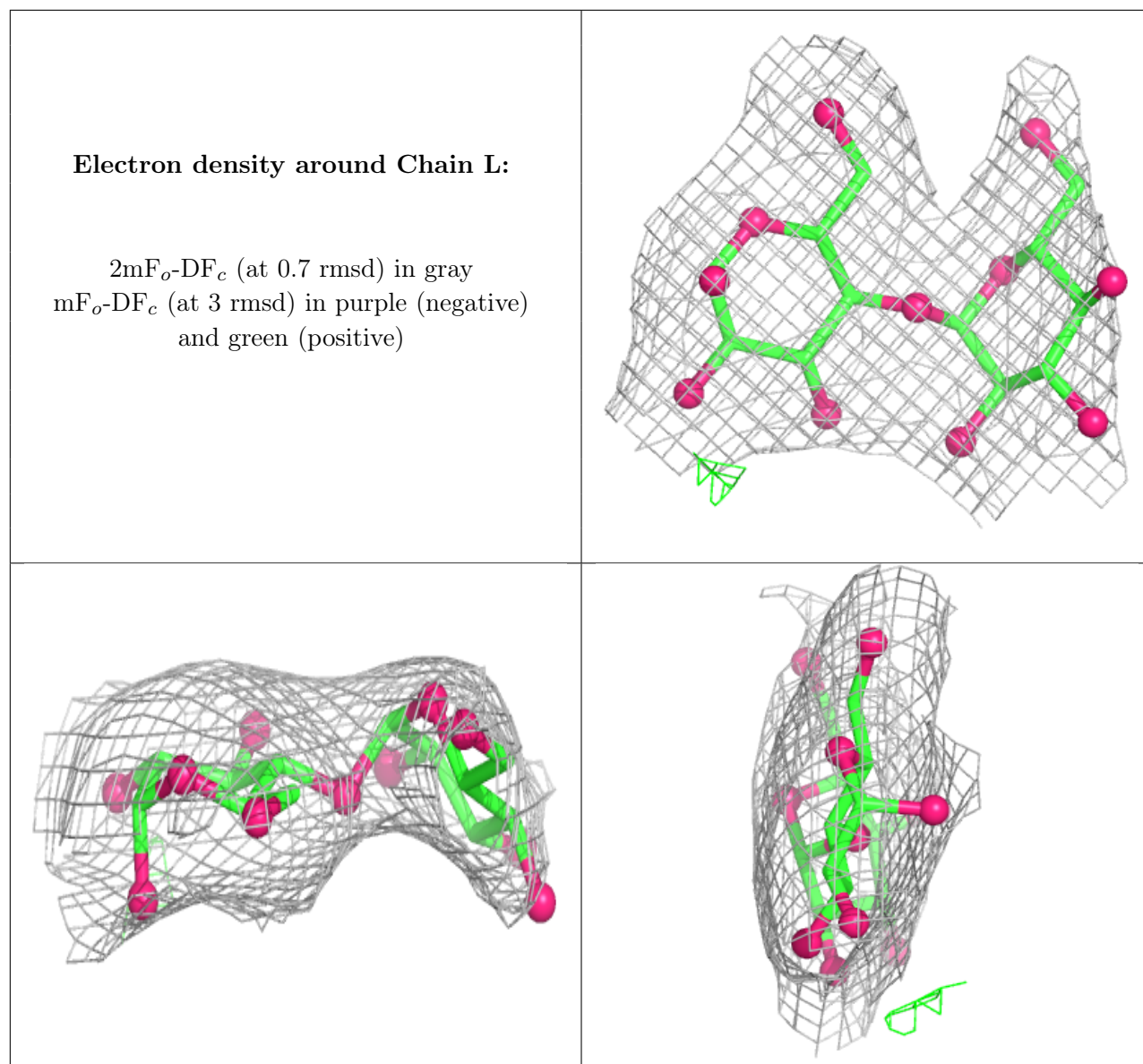
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain K:

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





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.