



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 7, 2026 – 02:34 AM UTC

PDB ID : 3SGJ / pdb_00003sgj
Title : Unique carbohydrate-carbohydrate interactions are required for high affinity binding between FcγIII and antibodies lacking core fucose
Authors : Ferrara, C.; Grau, S.; Jaeger, C.; Sondermann, P.; Bruenker, P.; Waldhauer, I.; Hennig, M.; Ruf, A.; Rufer, A.C.; Stihle, M.; Umana, P.; Benz, J.
Deposited on : 2011-06-15
Resolution : 2.20 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

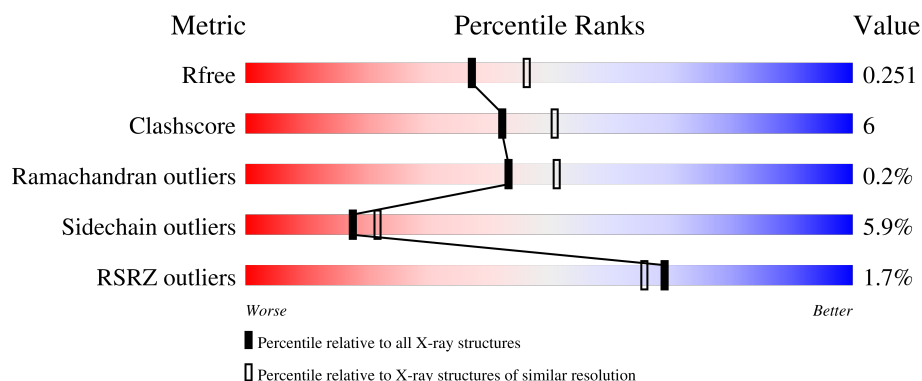
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

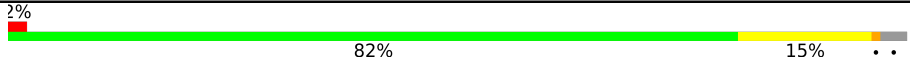
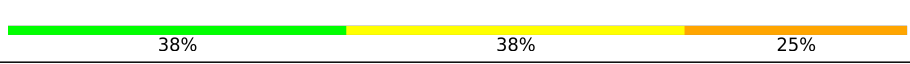
The reported resolution of this entry is 2.20 Å.

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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	225	
1	B	225	
2	C	204	
3	D	8	
3	E	8	

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Mol	Chain	Length	Quality of chain
4	F	5	
5	G	2	

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	FUC	D	8	X	-	-	-
3	MAN	E	4	X	-	-	-
4	MAN	F	4	X	-	-	-
4	MAN	F	5	X	-	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5443 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 human Fc fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1738	1107	291	332	8			
1	B	217	Total	C	N	O	S	0	0	0
			1726	1101	289	329	7			

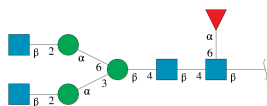
- Molecule 2 is a protein called human Fcg3a receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	169	Total	C	N	O	S	0	0	0
			1365	870	233	258	4			

There are 18 discrepancies between the modelled and reference sequences:

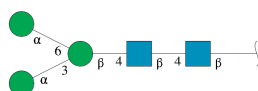
Chain	Residue	Modelled	Actual	Comment	Reference
C	38	GLN	ASN	engineered mutation	UNP P08637
C	74	GLN	ASN	engineered mutation	UNP P08637
C	158	VAL	PHE	SEE REMARK 999	UNP P08637
C	169	GLN	ASN	engineered mutation	UNP P08637
C	191	GLY	-	expression tag	UNP P08637
C	192	LYS	-	expression tag	UNP P08637
C	193	LYS	-	expression tag	UNP P08637
C	194	LYS	-	expression tag	UNP P08637
C	195	LYS	-	expression tag	UNP P08637
C	196	LYS	-	expression tag	UNP P08637
C	197	LYS	-	expression tag	UNP P08637
C	198	GLY	-	expression tag	UNP P08637
C	199	HIS	-	expression tag	UNP P08637
C	200	HIS	-	expression tag	UNP P08637
C	201	HIS	-	expression tag	UNP P08637
C	202	HIS	-	expression tag	UNP P08637
C	203	HIS	-	expression tag	UNP P08637
C	204	HIS	-	expression tag	UNP P08637

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	8	Total	C	N	O	0	0	0
			99	56	4	39			
3	E	8	Total	C	N	O	0	0	0
			99	56	4	39			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



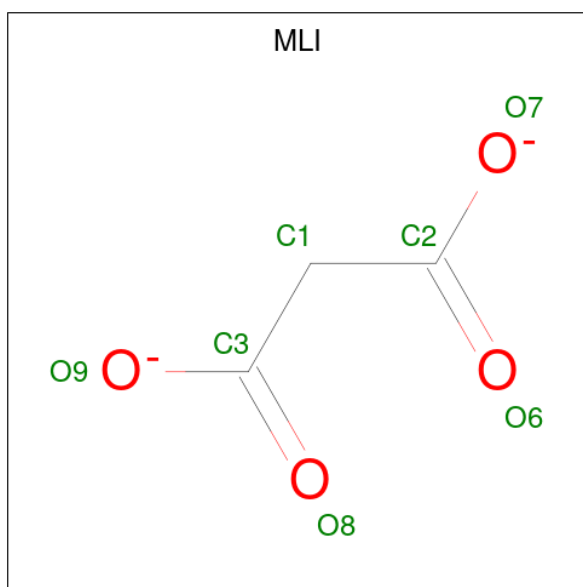
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 is MALONATE ION (CCD ID: MLI) (formula: C₃H₂O₄).



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

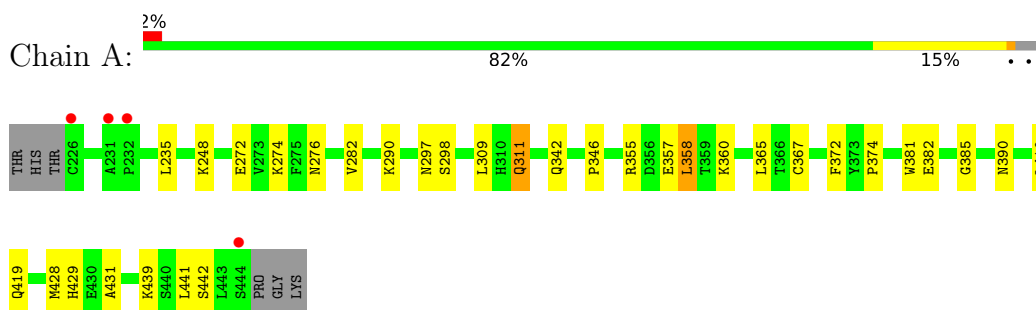
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	122	Total	O	0	0
			122	122		
7	B	111	Total	O	0	0
			111	111		
7	C	87	Total	O	0	0
			87	87		

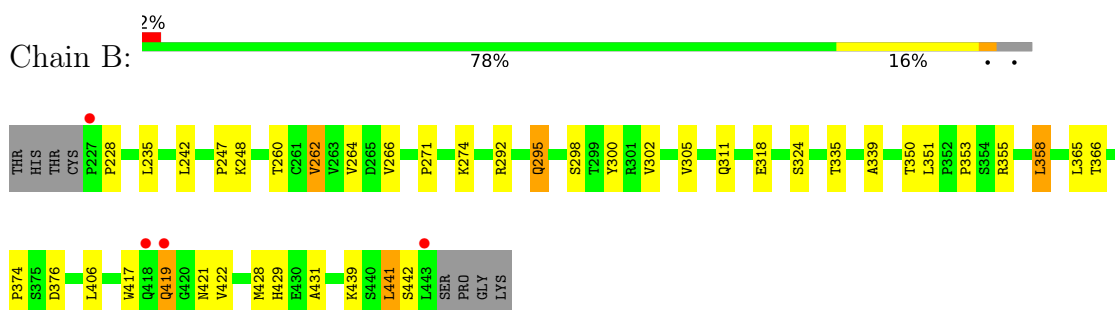
3 Residue-property plots [i](#)

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.

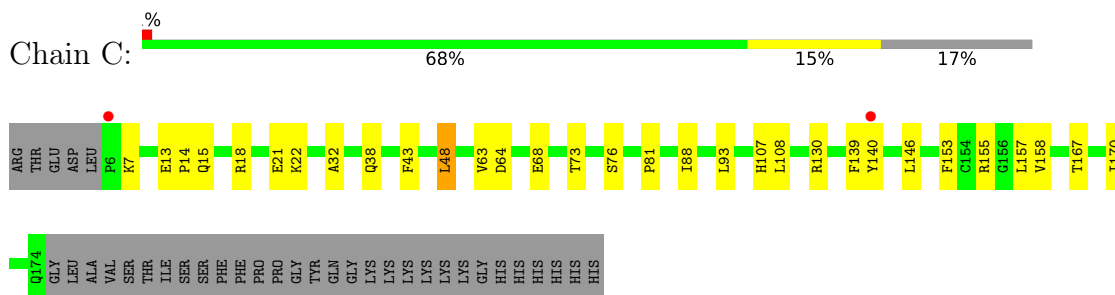
- Molecule 1: human Fc fragment



- Molecule 1: human Fc fragment

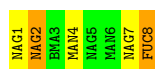


- Molecule 2: human Fcg3a receptor

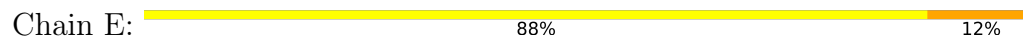


- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 4: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.65Å 88.48Å 140.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	140.29 – 2.20 74.84 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (140.29-2.20) 100.0 (74.84-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.6.0112	Depositor
R, R_{free}	0.195 , 0.252 0.195 , 0.251	Depositor DCC
R_{free} test set	2317 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	39.2	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 39.1	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	0.96	EDS
Total number of atoms	5443	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% 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, BMA, MAN, FUC, NAG

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	0.77	0/1788	0.85	0/2440
1	B	0.76	0/1776	0.86	1/2423 (0.0%)
2	C	0.83	0/1403	0.84	0/1906
All	All	0.78	0/4967	0.85	1/6769 (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	B	419	GLN	N-CA-C	5.46	118.06	111.40

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	1738	0	1703	17	0
1	B	1726	0	1694	21	0
2	C	1365	0	1309	20	0
3	D	99	0	85	2	0
3	E	99	0	85	1	0
4	F	61	0	52	3	0
5	G	28	0	25	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	7	0	2	1	0
7	A	122	0	0	2	0
7	B	111	0	0	0	0
7	C	87	0	0	3	0
All	All	5443	0	4955	63	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 6.

The worst 5 of 63 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:248:LYS:HG3	1:A:428:MET:HE1	1.13	1.10
2:C:63:VAL:HG21	2:C:88:ILE:HD11	1.45	0.95
1:A:248:LYS:HG3	1:A:428:MET:CE	2.00	0.91
1:B:350:THR:HB	1:B:441:LEU:HG	1.53	0.90
1:A:248:LYS:CG	1:A:428:MET:HE1	2.05	0.82

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	217/225 (96%)	214 (99%)	2 (1%)	1 (0%)	24	27
1	B	215/225 (96%)	211 (98%)	4 (2%)	0	100	100
2	C	167/204 (82%)	162 (97%)	5 (3%)	0	100	100
All	All	599/654 (92%)	587 (98%)	11 (2%)	1 (0%)	43	51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	385	GLY

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	203/208 (98%)	190 (94%)	13 (6%)	16	19
1	B	201/208 (97%)	190 (94%)	11 (6%)	19	24
2	C	152/183 (83%)	143 (94%)	9 (6%)	18	22
All	All	556/599 (93%)	523 (94%)	33 (6%)	18	22

5 of 33 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	146	LEU
2	C	157	LEU
2	C	170	ILE
1	A	442	SER
1	A	439	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
2	C	38	GLN
2	C	134	HIS
1	B	295	GLN
1	B	389	ASN
1	B	418	GLN

5.3.3 RNA [i](#)

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 ⓘ

23 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	NAG	D	1	3,1	14,14,15	0.52	0	17,19,21	1.15	2 (11%)
3	NAG	D	2	3	14,14,15	0.69	0	17,19,21	1.05	2 (11%)
3	BMA	D	3	3	11,11,12	0.32	0	15,15,17	1.04	0
3	MAN	D	4	3	11,11,12	0.61	0	15,15,17	1.30	1 (6%)
3	NAG	D	5	3	14,14,15	0.47	0	17,19,21	0.93	0
3	MAN	D	6	3	11,11,12	0.59	0	15,15,17	0.94	0
3	NAG	D	7	3	14,14,15	0.53	0	17,19,21	1.06	1 (5%)
3	FUC	D	8	3	10,10,11	0.81	0	14,14,16	1.75	3 (21%)
3	NAG	E	1	3,1	14,14,15	0.67	0	17,19,21	1.27	2 (11%)
3	NAG	E	2	3	14,14,15	0.65	0	17,19,21	1.39	3 (17%)
3	BMA	E	3	3	11,11,12	0.49	0	15,15,17	1.57	3 (20%)
3	MAN	E	4	3	11,11,12	1.36	1 (9%)	15,15,17	1.65	4 (26%)
3	NAG	E	5	3	14,14,15	0.59	0	17,19,21	1.90	5 (29%)
3	MAN	E	6	3	11,11,12	0.61	0	15,15,17	1.15	2 (13%)
3	NAG	E	7	3	14,14,15	0.53	0	17,19,21	1.99	4 (23%)
3	FUC	E	8	3	10,10,11	0.87	0	14,14,16	2.05	5 (35%)
4	NAG	F	1	4,2	14,14,15	1.81	1 (7%)	17,19,21	1.93	4 (23%)
4	NAG	F	2	4	14,14,15	0.73	1 (7%)	17,19,21	1.51	4 (23%)
4	BMA	F	3	4	11,11,12	0.46	0	15,15,17	1.07	1 (6%)
4	MAN	F	4	4	11,11,12	0.69	0	15,15,17	1.87	3 (20%)
4	MAN	F	5	4	11,11,12	0.68	0	15,15,17	1.47	3 (20%)
5	NAG	G	1	5,2	14,14,15	1.94	1 (7%)	17,19,21	2.17	6 (35%)
5	NAG	G	2	5	14,14,15	0.58	0	17,19,21	1.44	5 (29%)

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	NAG	D	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
3	BMA	D	3	3	-	0/2/19/22	0/1/1/1
3	MAN	D	4	3	-	2/2/19/22	0/1/1/1
3	NAG	D	5	3	-	4/6/23/26	0/1/1/1
3	MAN	D	6	3	-	2/2/19/22	0/1/1/1
3	NAG	D	7	3	-	2/6/23/26	0/1/1/1
3	FUC	D	8	3	1/1/4/5	-	0/1/1/1
3	NAG	E	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
3	MAN	E	4	3	1/1/4/5	2/2/19/22	0/1/1/1
3	NAG	E	5	3	-	2/6/23/26	0/1/1/1
3	MAN	E	6	3	-	2/2/19/22	0/1/1/1
3	NAG	E	7	3	-	3/6/23/26	0/1/1/1
3	FUC	E	8	3	-	-	0/1/1/1
4	NAG	F	1	4,2	-	3/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1
4	BMA	F	3	4	-	1/2/19/22	0/1/1/1
4	MAN	F	4	4	1/1/4/5	2/2/19/22	0/1/1/1
4	MAN	F	5	4	1/1/4/5	2/2/19/22	0/1/1/1
5	NAG	G	1	5,2	-	2/6/23/26	0/1/1/1
5	NAG	G	2	5	-	2/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	G	1	NAG	O5-C1	-6.63	1.32	1.43
4	F	1	NAG	O5-C1	-6.45	1.32	1.43
3	E	4	MAN	O5-C1	-4.19	1.36	1.43
4	F	2	NAG	O5-C1	-2.17	1.40	1.43

The worst 5 of 63 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	1	NAG	C1-O5-C5	-6.29	103.76	112.19
4	F	1	NAG	C1-O5-C5	-4.97	105.53	112.19
3	E	7	NAG	C4-C3-C2	-4.71	104.11	111.02
4	F	4	MAN	C1-O5-C5	4.50	118.22	112.19
4	F	4	MAN	C1-C2-C3	4.44	116.11	109.64

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	D	8	FUC	C1
3	E	4	MAN	C1
4	F	4	MAN	C1
4	F	5	MAN	C1

5 of 33 torsion outliers are listed below:

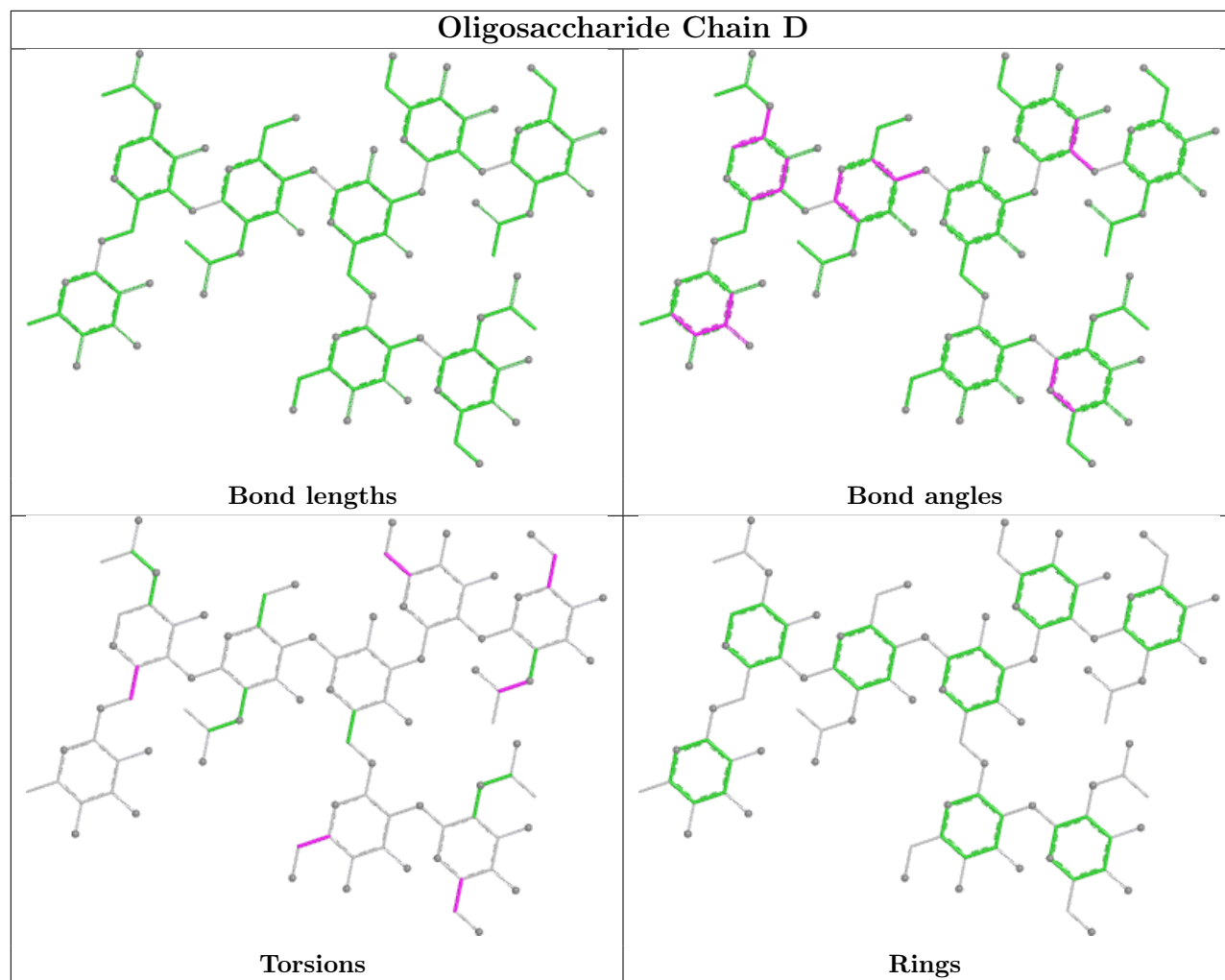
Mol	Chain	Res	Type	Atoms
4	F	1	NAG	C1-C2-N2-C7
3	E	6	MAN	O5-C5-C6-O6
3	E	7	NAG	O5-C5-C6-O6
3	D	7	NAG	O5-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6

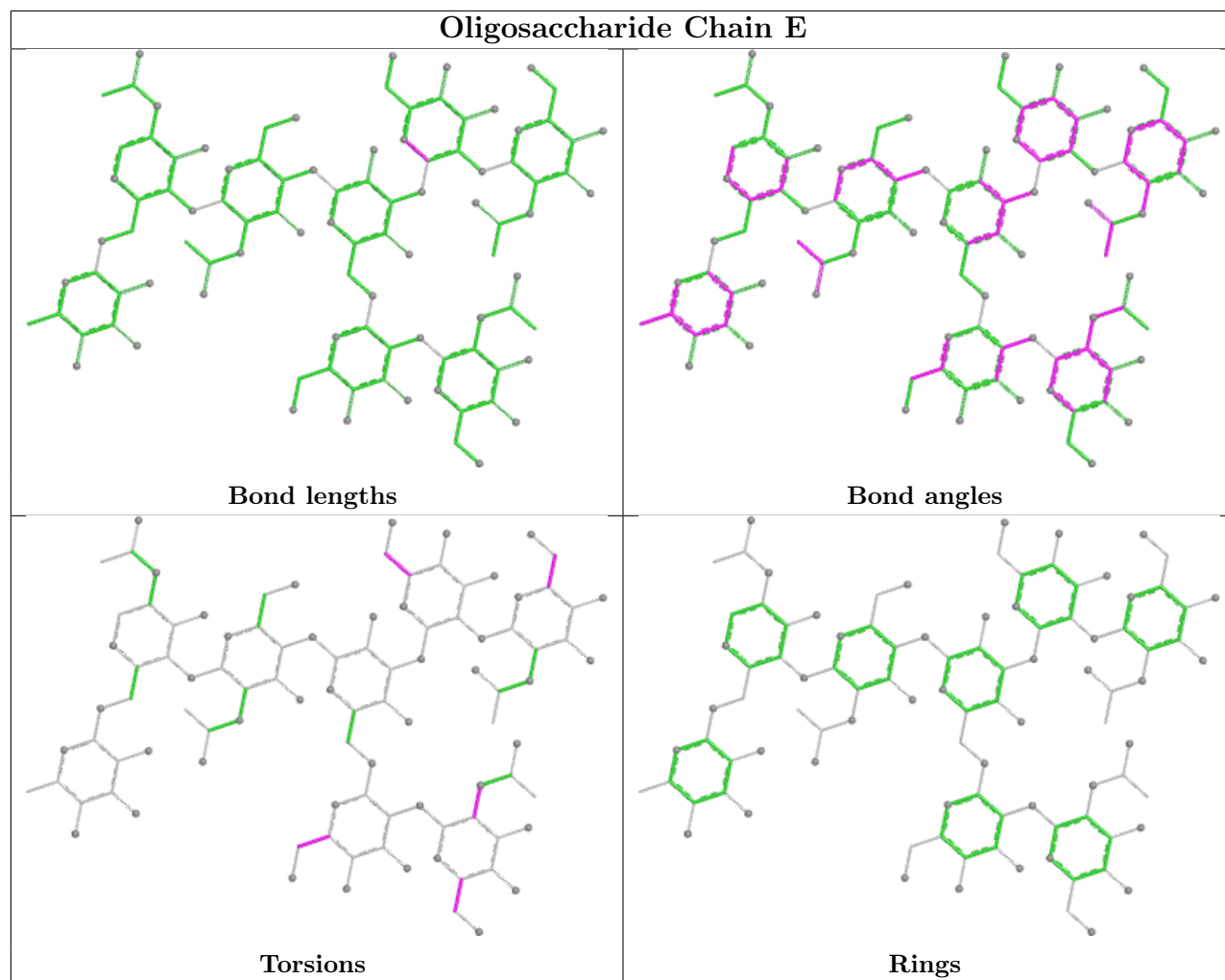
There are no ring outliers.

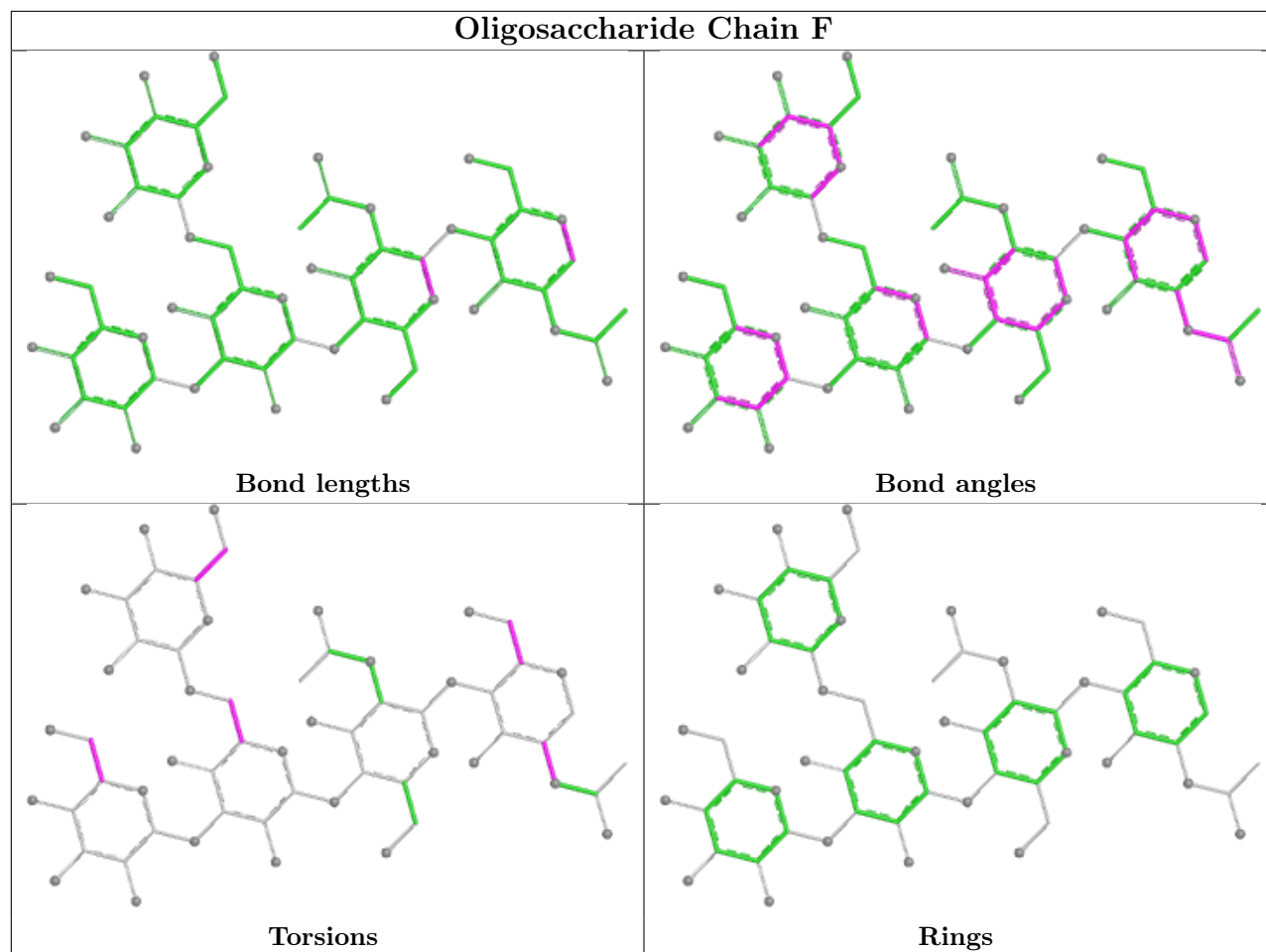
8 monomers are involved in 6 short contacts:

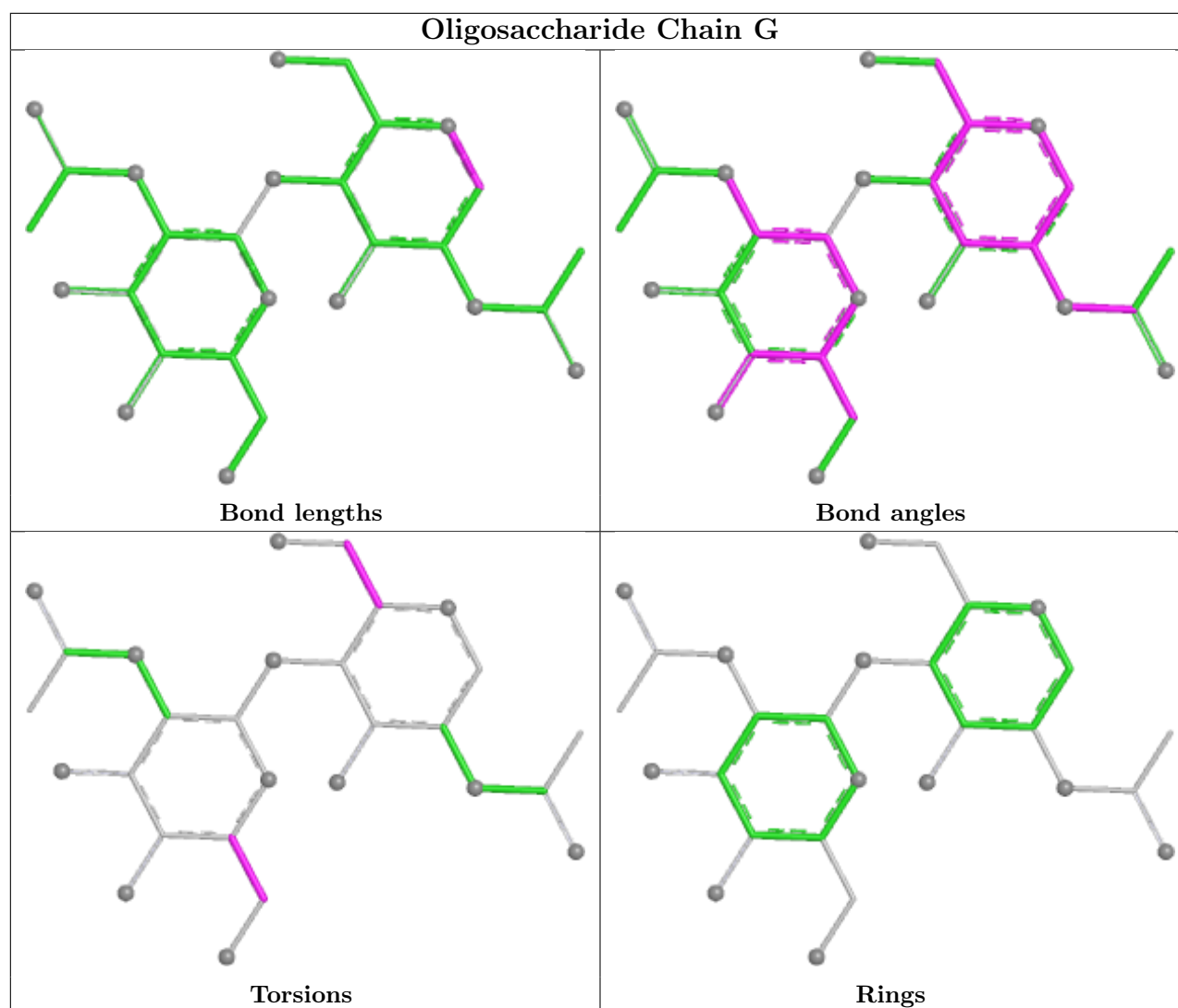
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	8	FUC	2	0
3	E	1	NAG	1	0
4	F	1	NAG	1	0
4	F	3	BMA	1	0
4	F	2	NAG	1	0
4	F	5	MAN	1	0
3	D	2	NAG	1	0
5	G	2	NAG	1	0

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](#)

1 ligand is 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	$\# Z > 2$
6	MLI	A	509	-	6,6,6	1.26	0	7,7,7	1.14	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
6	MLI	A	509	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

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
6	A	509	MLI	1	0

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	219/225 (97%)	-0.17	4 (1%) 67 64	24, 37, 66, 81	0
1	B	217/225 (96%)	-0.11	4 (1%) 67 64	27, 41, 64, 88	0
2	C	169/204 (82%)	-0.24	2 (1%) 76 74	24, 34, 56, 87	0
All	All	605/654 (92%)	-0.17	10 (1%) 69 66	24, 38, 64, 88	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	231	ALA	3.5
1	B	443	LEU	3.3
1	A	232	PRO	3.3
2	C	140	TYR	3.0
1	A	444	SER	2.9

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
4	MAN	F	4	11/12	0.42	0.16	88,97,99,102	0
3	NAG	E	5	14/15	0.58	0.16	65,78,87,89	0
4	MAN	F	5	11/12	0.64	0.12	97,100,103,104	0

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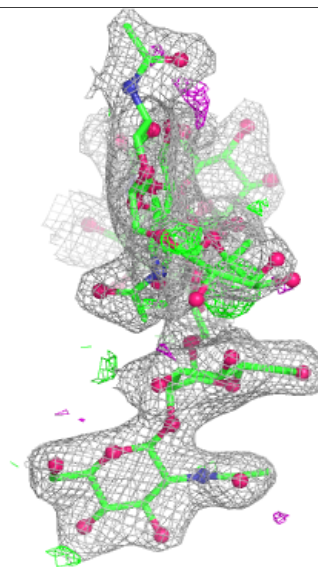
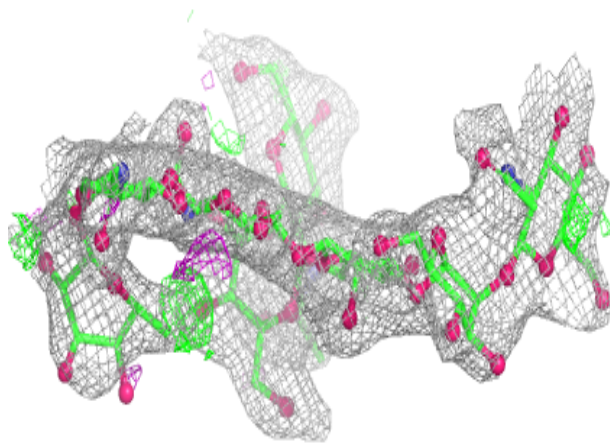
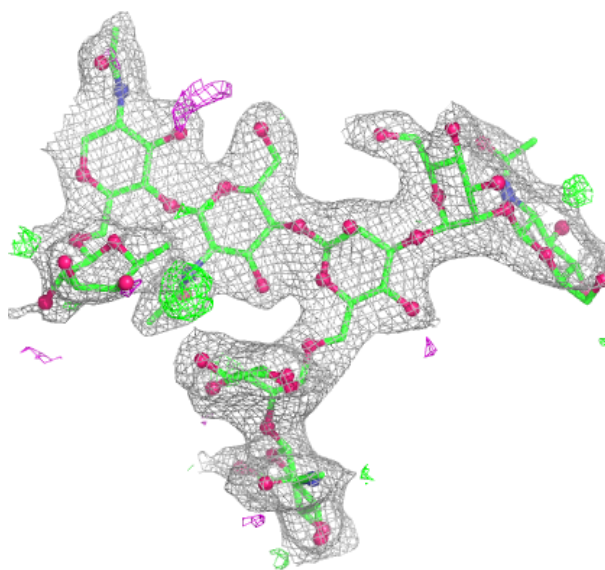
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	D	5	14/15	0.67	0.15	78,83,88,92	0
3	FUC	D	8	10/11	0.68	0.17	64,73,77,86	0
3	FUC	E	8	10/11	0.72	0.15	54,60,63,65	0
4	BMA	F	3	11/12	0.78	0.11	77,85,92,93	0
4	NAG	F	2	14/15	0.81	0.12	52,58,69,71	0
3	MAN	E	4	11/12	0.84	0.11	56,59,61,69	0
3	MAN	D	4	11/12	0.87	0.09	50,58,66,69	0
4	NAG	F	1	14/15	0.87	0.11	37,43,51,53	0
3	NAG	E	7	14/15	0.88	0.11	44,48,54,56	0
5	NAG	G	2	14/15	0.90	0.09	36,38,40,40	0
5	NAG	G	1	14/15	0.92	0.08	28,31,36,36	0
3	MAN	D	6	11/12	0.93	0.07	37,38,39,41	0
3	BMA	E	3	11/12	0.93	0.07	35,40,43,51	0
3	MAN	E	6	11/12	0.93	0.07	37,41,43,44	0
3	NAG	D	1	14/15	0.94	0.07	28,31,44,53	0
3	BMA	D	3	11/12	0.95	0.06	31,34,37,41	0
3	NAG	E	2	14/15	0.95	0.07	25,28,30,33	0
3	NAG	D	7	14/15	0.95	0.07	37,42,46,47	0
3	NAG	E	1	14/15	0.96	0.07	30,33,38,41	0
3	NAG	D	2	14/15	0.97	0.05	29,30,31,31	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

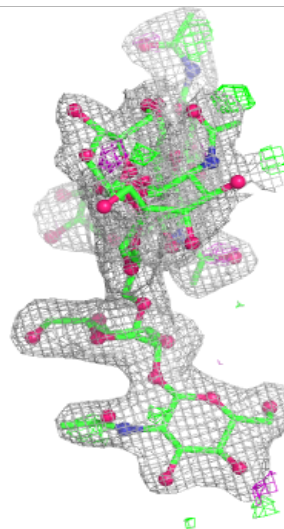
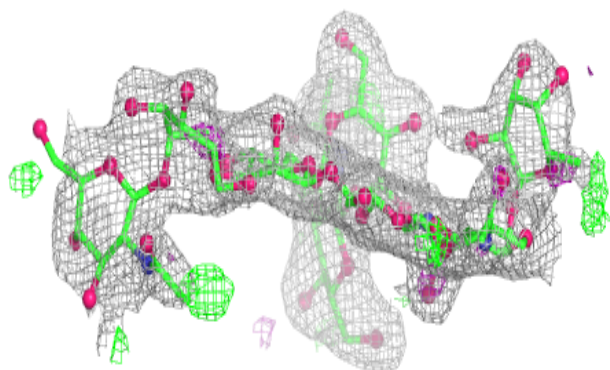
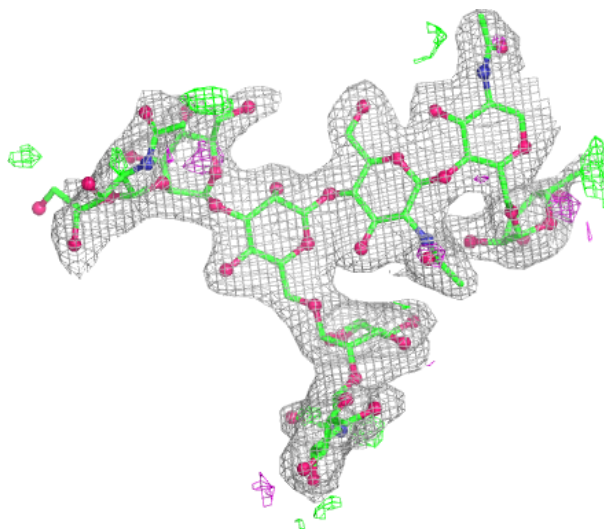
Electron density around Chain D:

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



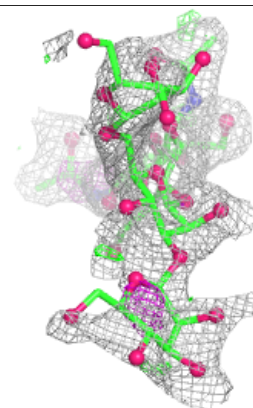
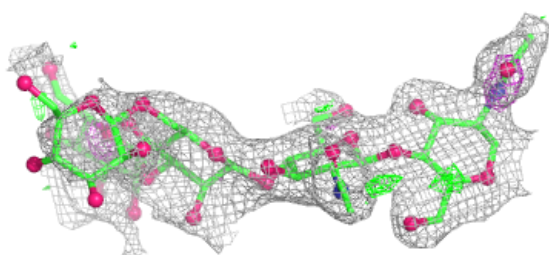
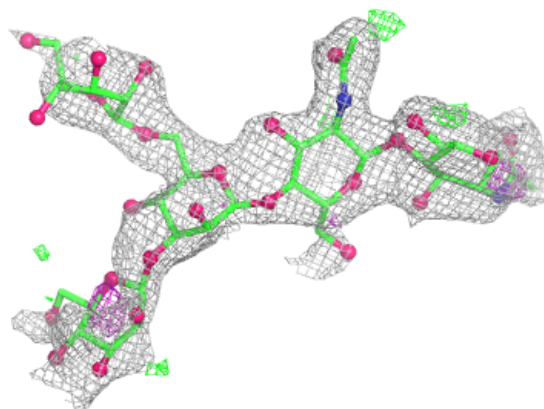
Electron density around Chain E:

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

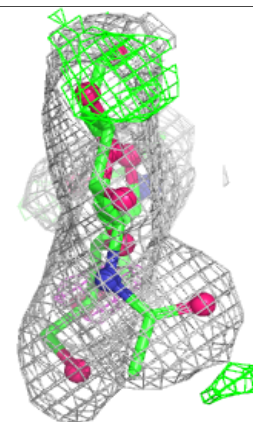
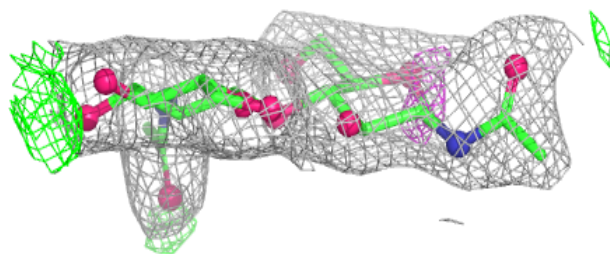
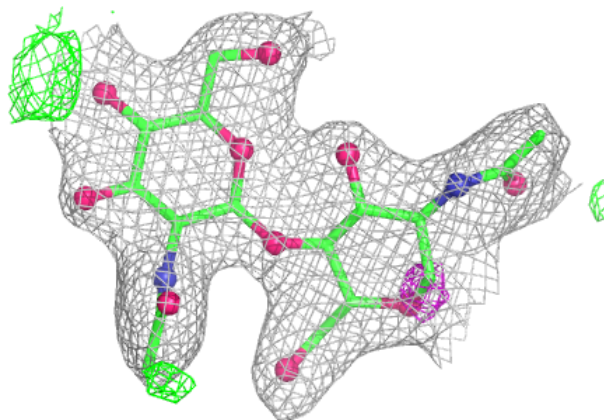


Electron density around Chain F:

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

**Electron density around Chain G:**

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



6.4 Ligands

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	MLI	A	509	7/7	0.92	0.14	40,43,46,48	0

6.5 Other polymers

There are no such residues in this entry.