



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 02:07 AM UTC

PDB ID : 9D06 / pdb_00009d06
Title : Crystal structure of IgG1 FC M252R at pH 5.6
Authors : Reddem, E.R.; Shapiro, L.
Deposited on : 2024-08-06
Resolution : 2.33 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

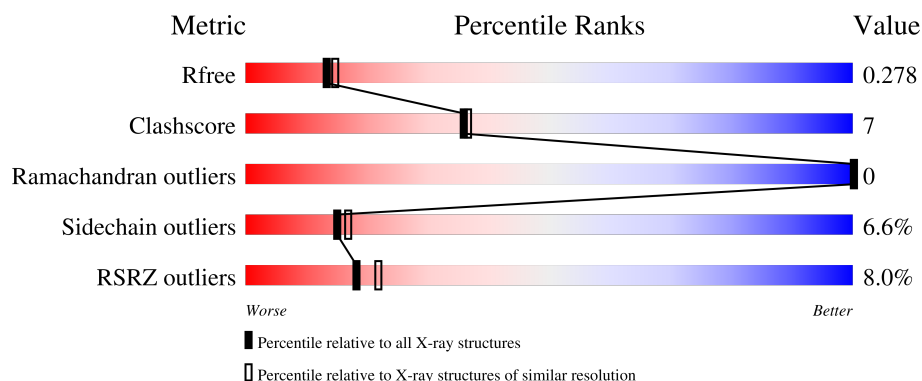
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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	231	<div> <div>3%</div> <div>72%</div> <div>17%</div> <div>10%</div> </div>
1	B	231	<div> <div>8%</div> <div>32%</div> <div>13%</div> <div>55%</div> </div>
2	C	7	<div> <div>14%</div> <div>71%</div> <div>14%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2604 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 Immunoglobulin gamma-1 heavy chain Fc fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1667	1060	283	319	5			
1	B	104	Total	C	N	O	S	0	0	0
			826	521	139	163	3			

There are 40 discrepancies between the modelled and reference sequences:

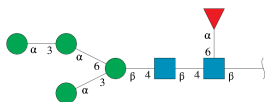
Chain	Residue	Modelled	Actual	Comment	Reference
A	217	MET	-	initiating methionine	UNP P0DOX5
A	218	GLY	-	expression tag	UNP P0DOX5
A	219	TRP	-	expression tag	UNP P0DOX5
A	220	SER	-	expression tag	UNP P0DOX5
A	221	CYS	-	expression tag	UNP P0DOX5
A	222	ILE	-	expression tag	UNP P0DOX5
A	223	ILE	-	expression tag	UNP P0DOX5
A	224	LEU	-	expression tag	UNP P0DOX5
A	225	PHE	-	expression tag	UNP P0DOX5
A	226	LEU	-	expression tag	UNP P0DOX5
A	227	VAL	-	expression tag	UNP P0DOX5
A	228	ALA	-	expression tag	UNP P0DOX5
A	229	THR	-	expression tag	UNP P0DOX5
A	230	ALA	-	expression tag	UNP P0DOX5
A	231	THR	-	expression tag	UNP P0DOX5
A	232	GLY	-	expression tag	UNP P0DOX5
A	233	VAL	-	expression tag	UNP P0DOX5
A	234	HIS	-	expression tag	UNP P0DOX5
A	235	SER	-	expression tag	UNP P0DOX5
A	252	ARG	MET	engineered mutation	UNP P0DOX5
B	217	MET	-	initiating methionine	UNP P0DOX5
B	218	GLY	-	expression tag	UNP P0DOX5
B	219	TRP	-	expression tag	UNP P0DOX5
B	220	SER	-	expression tag	UNP P0DOX5
B	221	CYS	-	expression tag	UNP P0DOX5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	222	ILE	-	expression tag	UNP P0DOX5
B	223	ILE	-	expression tag	UNP P0DOX5
B	224	LEU	-	expression tag	UNP P0DOX5
B	225	PHE	-	expression tag	UNP P0DOX5
B	226	LEU	-	expression tag	UNP P0DOX5
B	227	VAL	-	expression tag	UNP P0DOX5
B	228	ALA	-	expression tag	UNP P0DOX5
B	229	THR	-	expression tag	UNP P0DOX5
B	230	ALA	-	expression tag	UNP P0DOX5
B	231	THR	-	expression tag	UNP P0DOX5
B	232	GLY	-	expression tag	UNP P0DOX5
B	233	VAL	-	expression tag	UNP P0DOX5
B	234	HIS	-	expression tag	UNP P0DOX5
B	235	SER	-	expression tag	UNP P0DOX5
B	252	ARG	MET	engineered mutation	UNP P0DOX5

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]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
2	C	7	Total	C	N	O	0	0	0
			82	46	2	34			

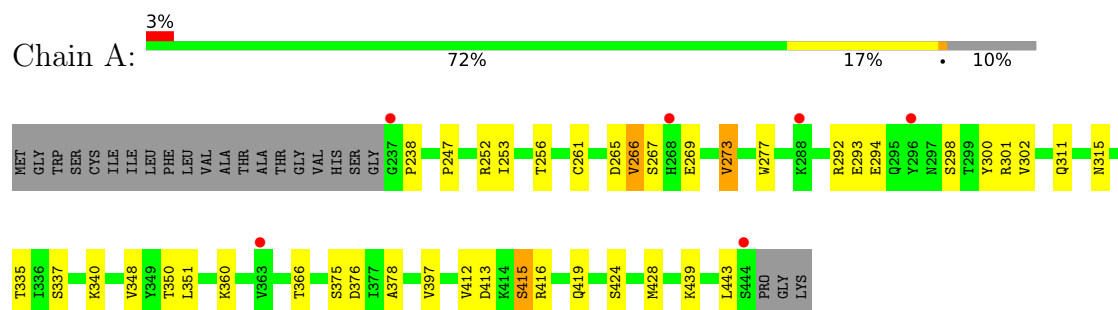
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	22	Total	O	0	0
			22	22		
3	B	7	Total	O	0	0
			7	7		

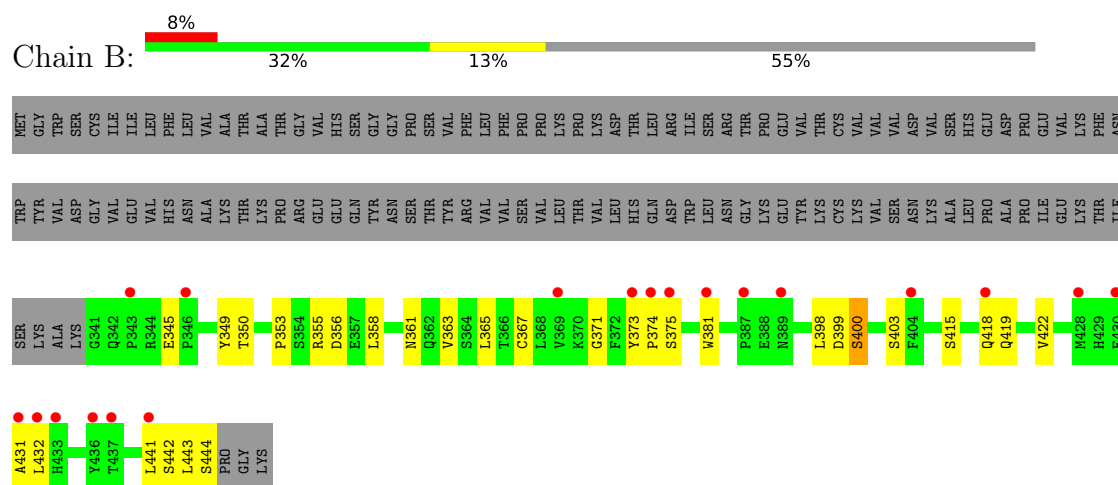
3 Residue-property plots

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: Immunoglobulin gamma-1 heavy chain Fc fragment



- Molecule 1: Immunoglobulin gamma-1 heavy chain Fc fragment



- Molecule 2: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]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



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	67.39Å 111.45Å 63.66Å 90.00° 95.74° 90.00°	Depositor
Resolution (Å)	32.50 – 2.33 32.50 – 2.33	Depositor EDS
% Data completeness (in resolution range)	99.0 (32.50-2.33) 99.0 (32.50-2.33)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.225 , 0.279 0.226 , 0.278	Depositor DCC
R_{free} test set	1988 reflections (9.92%)	wwPDB-VP
Wilson B-factor (Å ²)	59.2	Xtriage
Anisotropy	0.414	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2604	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FUC, MAN, NAG, BMA

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	0.44	0/1713	0.64	0/2334
1	B	0.39	0/849	0.56	0/1156
All	All	0.42	0/2562	0.61	0/3490

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	1667	0	1634	19	0
1	B	826	0	784	18	0
2	C	82	0	70	1	0
3	A	22	0	0	1	0
3	B	7	0	0	0	0
All	All	2604	0	2488	37	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 7.

All (37) 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:351:LEU:HB2	1:A:366:THR:HB	1.75	0.69
1:A:413:ASP:HB2	1:A:416:ARG:HG3	1.79	0.65
1:B:353:PRO:HB3	1:B:363:VAL:HG13	1.79	0.63
1:B:353:PRO:HG3	1:B:363:VAL:HG11	1.81	0.62
1:A:252:ARG:HH11	1:A:252:ARG:HG2	1.65	0.60
1:B:355:ARG:HA	1:B:358:LEU:HD12	1.83	0.60
1:A:238:PRO:HA	1:A:265:ASP:HB2	1.83	0.60
1:B:415:SER:O	1:B:419:GLN:HG3	2.04	0.58
1:A:266:VAL:CG2	1:A:300:TYR:HB2	2.35	0.56
2:C:2:NAG:H83	2:C:4:MAN:H61	1.88	0.56
1:B:367:CYS:HB2	1:B:381:TRP:CZ2	2.43	0.54
1:B:345:GLU:HG2	1:B:432:LEU:HD23	1.89	0.54
1:A:301:ARG:NH1	3:A:501:HOH:O	2.29	0.53
1:B:353:PRO:HG3	1:B:363:VAL:CG1	2.38	0.52
1:B:418:GLN:HA	1:B:443:LEU:HD12	1.91	0.51
1:B:350:THR:HB	1:B:441:LEU:HB2	1.92	0.51
1:A:273:VAL:HG13	1:A:302:VAL:HG11	1.93	0.51
1:A:292:ARG:HB3	1:A:302:VAL:HG22	1.93	0.51
1:A:311:GLN:O	1:A:315:ASN:OD1	2.28	0.51
1:A:348:VAL:HB	1:A:439:LYS:HD3	1.93	0.50
1:A:415:SER:O	1:A:419:GLN:HG3	2.13	0.48
1:B:353:PRO:HD3	1:B:365:LEU:HD23	1.94	0.48
1:A:294:GLU:OE1	1:A:298:SER:HA	2.14	0.48
1:A:267:SER:HB2	1:A:269:GLU:OE1	2.13	0.47
1:A:378:ALA:HB3	1:A:428:MET:HE3	1.97	0.46
1:B:398:LEU:HD13	1:B:399:ASP:O	2.14	0.46
1:B:422:VAL:HA	1:B:442:SER:HB3	1.98	0.46
1:B:399:ASP:OD1	1:B:400:SER:N	2.47	0.45
1:A:261:CYS:HB2	1:A:277:TRP:CH2	2.53	0.44
1:A:350:THR:C	1:A:351:LEU:HD23	2.43	0.43
1:A:247:PRO:HG3	1:A:376:ASP:OD2	2.19	0.42
1:B:353:PRO:HD3	1:B:365:LEU:CD2	2.48	0.42
1:B:373:TYR:CD2	1:B:374:PRO:HA	2.54	0.42
1:A:360:LYS:HE3	1:B:349:TYR:OH	2.19	0.42
1:B:345:GLU:HG3	1:B:431:ALA:O	2.20	0.41
1:B:371:GLY:HA2	1:B:403:SER:OG	2.21	0.41
1:A:350:THR:O	1:A:351:LEU:HD23	2.21	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	206/231 (89%)	201 (98%)	5 (2%)	0	100	100
1	B	102/231 (44%)	99 (97%)	3 (3%)	0	100	100
All	All	308/462 (67%)	300 (97%)	8 (3%)	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	194/211 (92%)	180 (93%)	14 (7%)	13	14
1	B	96/211 (46%)	91 (95%)	5 (5%)	21	25
All	All	290/422 (69%)	271 (93%)	19 (7%)	15	17

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	253	ILE
1	A	256	THR
1	A	266	VAL
1	A	273	VAL
1	A	293	GLU
1	A	335	THR
1	A	337	SER
1	A	340	LYS

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Mol	Chain	Res	Type
1	A	375	SER
1	A	397	VAL
1	A	412	VAL
1	A	415	SER
1	A	424	SER
1	A	443	LEU
1	B	356	ASP
1	B	361	ASN
1	B	375	SER
1	B	400	SER
1	B	444	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	HIS
1	A	276	ASN
1	A	389	ASN
1	B	347	GLN
1	B	434	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 ⓘ

7 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	NAG	C	1	2,1	14,14,15	0.32	0	17,19,21	0.52	0
2	NAG	C	2	2	14,14,15	0.33	0	17,19,21	0.44	0
2	BMA	C	3	2	11,11,12	0.71	0	15,15,17	1.24	1 (6%)
2	MAN	C	4	2	11,11,12	1.62	2 (18%)	15,15,17	1.29	4 (26%)
2	MAN	C	5	2	11,11,12	0.85	0	15,15,17	1.58	2 (13%)
2	MAN	C	6	2	11,11,12	1.05	0	15,15,17	1.02	1 (6%)
2	FUC	C	7	2	10,10,11	1.20	1 (10%)	14,14,16	0.98	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	NAG	C	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
2	MAN	C	4	2	-	0/2/19/22	0/1/1/1
2	MAN	C	5	2	-	2/2/19/22	0/1/1/1
2	MAN	C	6	2	-	2/2/19/22	0/1/1/1
2	FUC	C	7	2	-	-	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	4	MAN	O5-C1	-3.66	1.37	1.43
2	C	7	FUC	C4-C5	2.92	1.59	1.52
2	C	4	MAN	O2-C2	2.61	1.48	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	5	MAN	C1-O5-C5	3.74	117.19	112.19
2	C	3	BMA	C1-O5-C5	3.66	117.09	112.19
2	C	6	MAN	C1-O5-C5	2.67	115.77	112.19
2	C	5	MAN	O2-C2-C3	-2.65	104.66	110.15
2	C	4	MAN	O3-C3-C4	-2.19	105.21	110.38
2	C	4	MAN	O2-C2-C1	2.17	114.19	109.22
2	C	4	MAN	C2-C3-C4	2.07	114.50	110.86
2	C	4	MAN	C3-C4-C5	2.03	113.90	110.23

There are no chirality outliers.

All (6) torsion outliers are listed below:

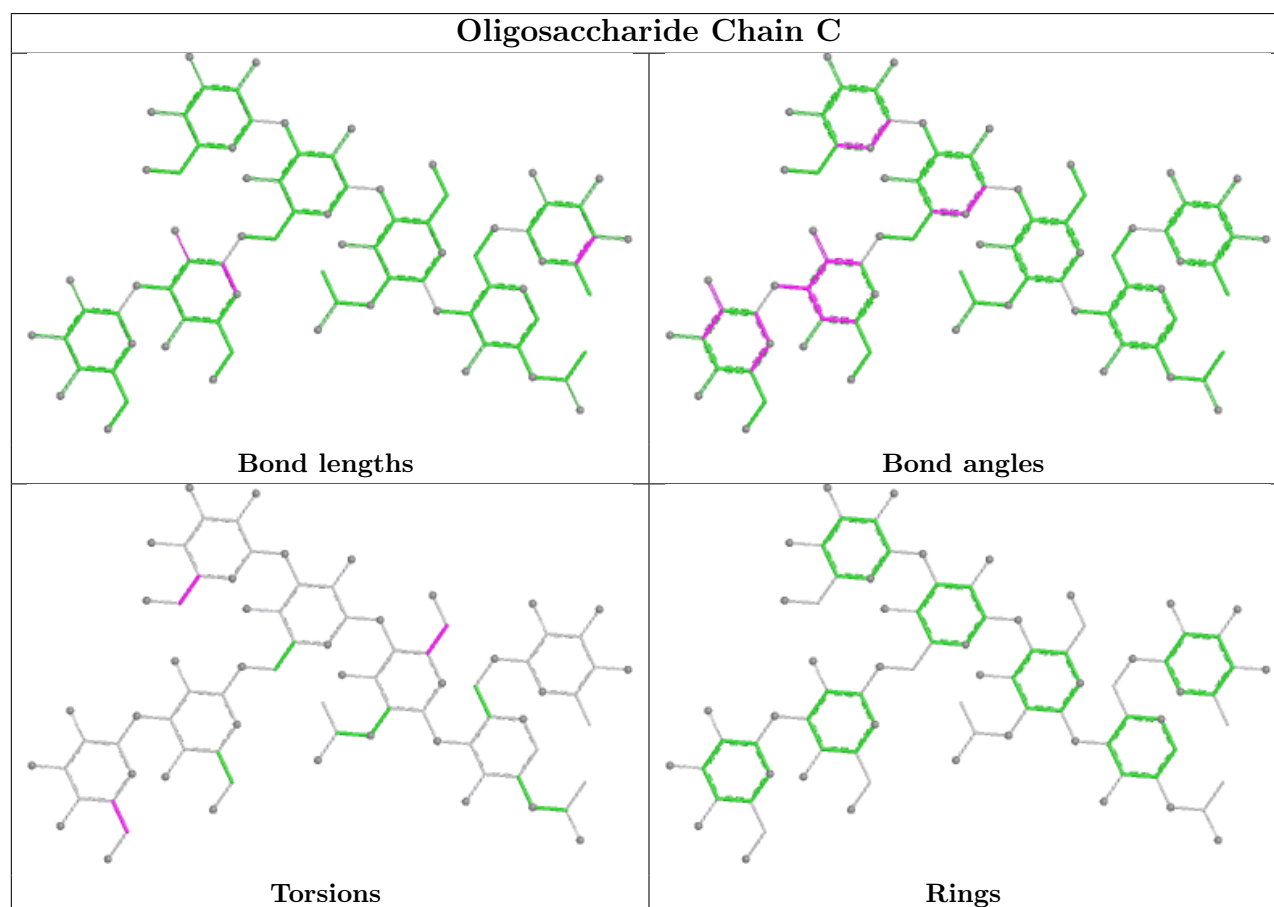
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	O5-C5-C6-O6
2	C	5	MAN	O5-C5-C6-O6
2	C	5	MAN	C4-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	C	6	MAN	C4-C5-C6-O6
2	C	6	MAN	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	4	MAN	1	0
2	C	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

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	208/231 (90%)	0.71	6 (2%) 53 60	46, 61, 82, 102	0
1	B	104/231 (45%)	1.14	19 (18%) 3 4	53, 72, 95, 103	0
All	All	312/462 (67%)	0.85	25 (8%) 18 22	46, 63, 90, 103	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	436	TYR	3.5
1	B	431	ALA	3.0
1	B	433	HIS	2.8
1	A	296	TYR	2.7
1	A	268	HIS	2.7
1	B	389	ASN	2.6
1	A	363	VAL	2.6
1	A	444	SER	2.6
1	B	418	GLN	2.6
1	B	381	TRP	2.4
1	A	237	GLY	2.4
1	B	373	TYR	2.3
1	B	375	SER	2.3
1	B	430	GLU	2.3
1	B	437	THR	2.3
1	B	441	LEU	2.2
1	B	404	PHE	2.2
1	B	369	VAL	2.1
1	B	346	PRO	2.1
1	B	387	PRO	2.1
1	B	428	MET	2.0
1	B	432	LEU	2.0
1	B	343	PRO	2.0
1	B	374	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	288	LYS	2.0

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

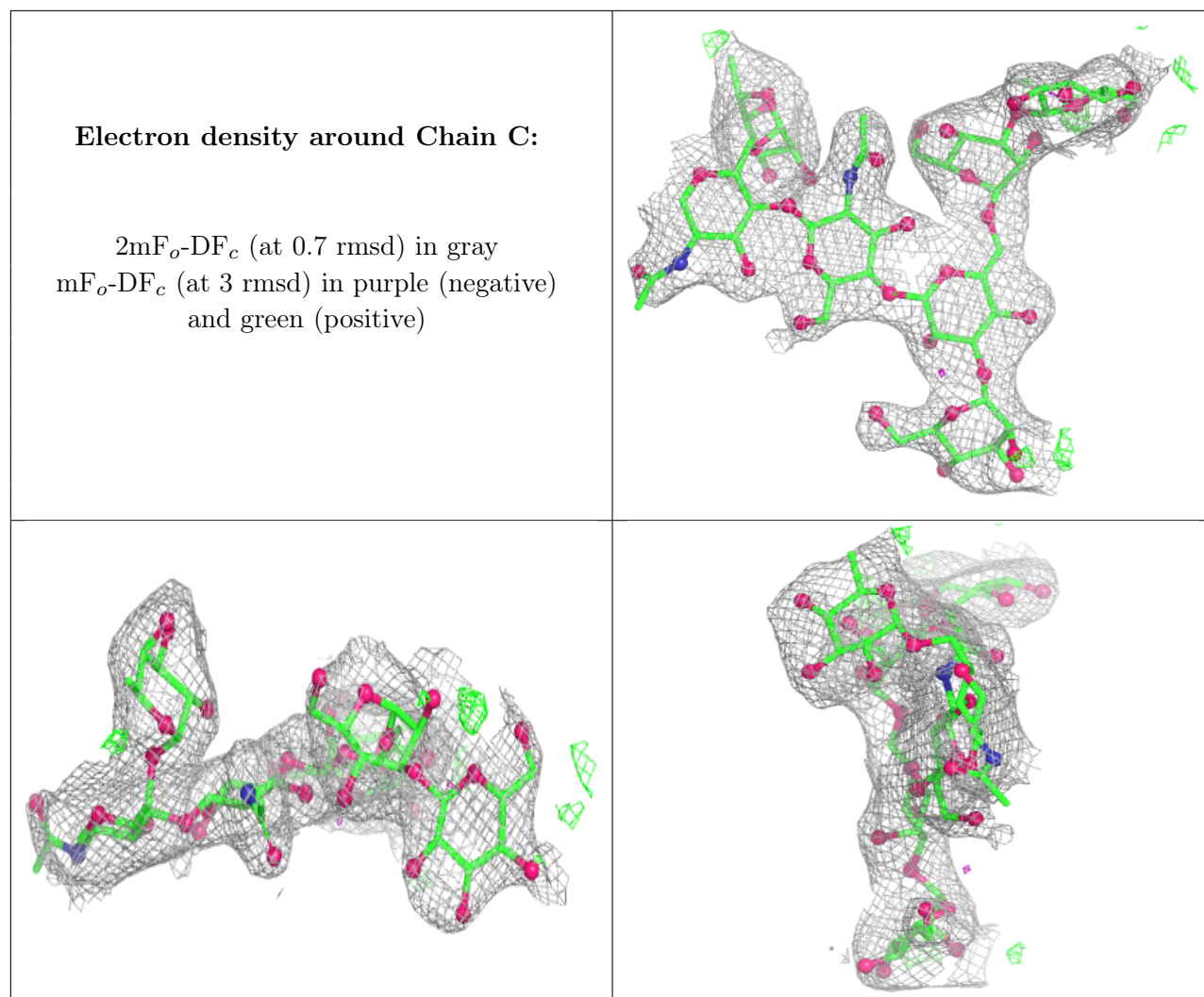
There are no non-standard protein/DNA/RNA residues in this entry.

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	MAN	C	6	11/12	0.69	0.14	81,89,91,93	0
2	FUC	C	7	10/11	0.69	0.14	82,90,93,93	0
2	MAN	C	5	11/12	0.76	0.16	75,84,93,99	0
2	MAN	C	4	11/12	0.77	0.11	81,87,91,92	0
2	NAG	C	1	14/15	0.86	0.12	79,83,88,89	0
2	BMA	C	3	11/12	0.87	0.09	76,78,81,82	0
2	NAG	C	2	14/15	0.91	0.10	73,81,86,87	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.



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.