



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 11:31 AM UTC

PDB ID : 2GJ7 / pdb_00002gj7
Title : Crystal Structure of a gE-gI/Fc complex
Authors : Sprague, E.R.; Wang, C.; Baker, D.; Bjorkman, P.J.
Deposited on : 2006-03-30
Resolution : 5.00 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

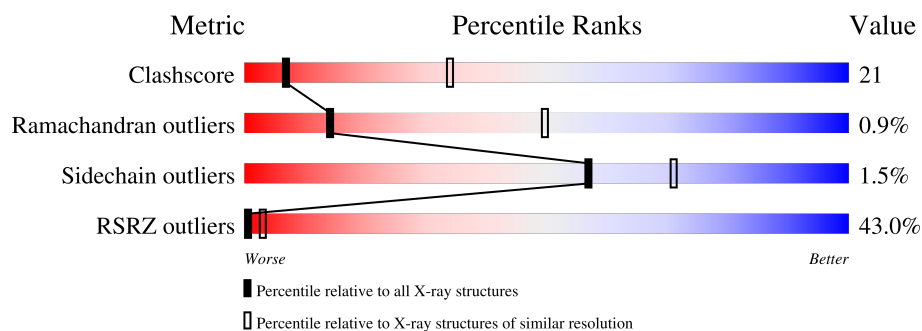
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.00 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))
Clashscore	190562	1040 (6.00-4.00)
Ramachandran outliers	187476	1015 (6.06-3.92)
Sidechain outliers	187428	1136 (6.10-3.90)
RSRZ outliers	180081	1004 (6.02-3.98)

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	227	45% 73% 18% 9%
1	B	227	43% 71% 18% 9%
2	E	401	18% 32% 10% 57%
2	F	401	13% 32% 10% 57%
3	C	8	75% 12% 12%
3	D	8	62% 38%

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	C	8	X	-	-	-
3	FUC	D	8	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6148 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 Ig gamma-1 chain C region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	19	0	0
			1659	1056	279	317	7			
1	B	207	Total	C	N	O	S	11	0	0
			1659	1056	279	317	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	221	LEU	-	cloning artifact	UNP P01857
A	222	GLU	-	cloning artifact	UNP P01857
A	356	GLU	ASP	conflict	UNP P01857
A	358	MET	LEU	conflict	UNP P01857
B	221	LEU	-	cloning artifact	UNP P01857
B	222	GLU	-	cloning artifact	UNP P01857
B	356	GLU	ASP	conflict	UNP P01857
B	358	MET	LEU	conflict	UNP P01857

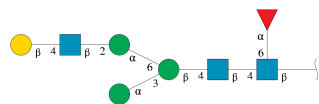
- Molecule 2 is a protein called Glycoprotein E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	171	Total	C	N	O	S	0	0	0
			1319	829	229	250	11			
2	E	171	Total	C	N	O	S	0	0	0
			1319	829	229	250	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	420	ASP	-	cloning artifact	UNP P04488
F	421	ILE	-	cloning artifact	UNP P04488
E	420	ASP	-	cloning artifact	UNP P04488
E	421	ILE	-	cloning artifact	UNP P04488

- Molecule 3 is an oligosaccharide called beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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.

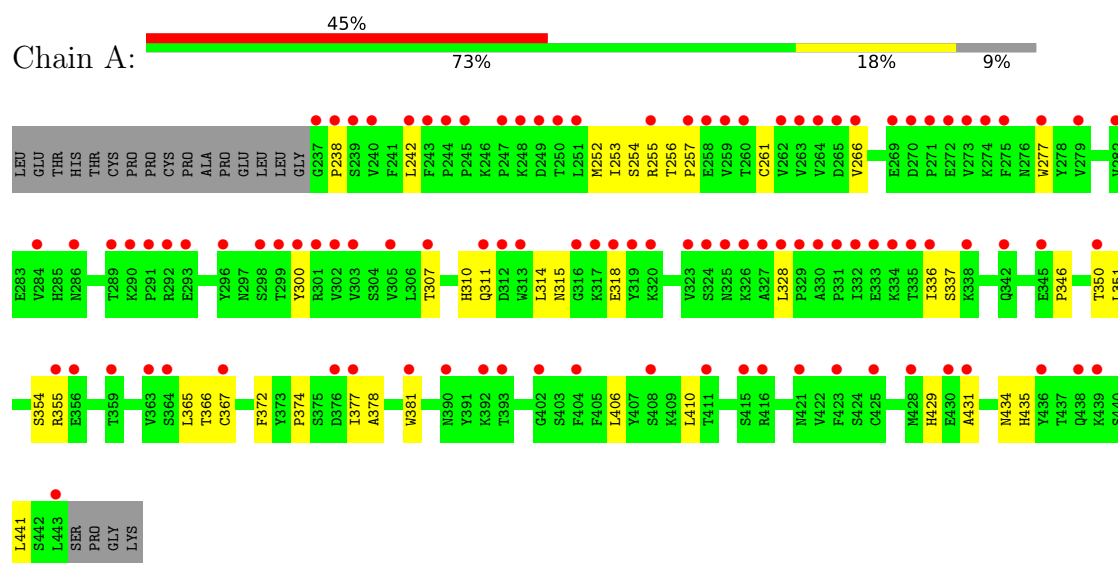


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	0	0	0
			96	54	3	39			
3	D	8	Total	C	N	O	0	0	0
			96	54	3	39			

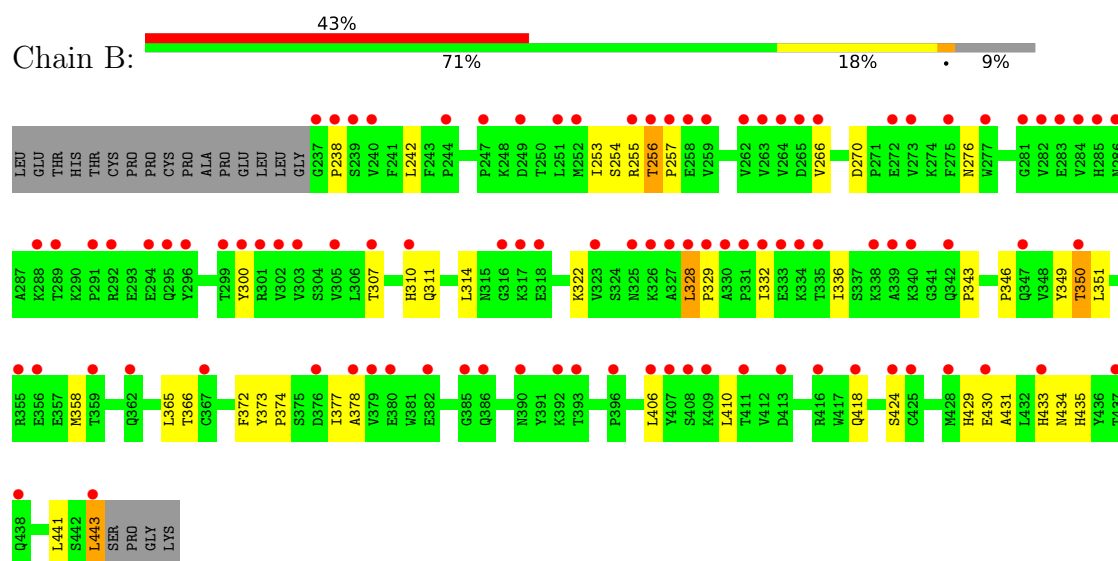
3 Residue-property plots [i](#)

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.

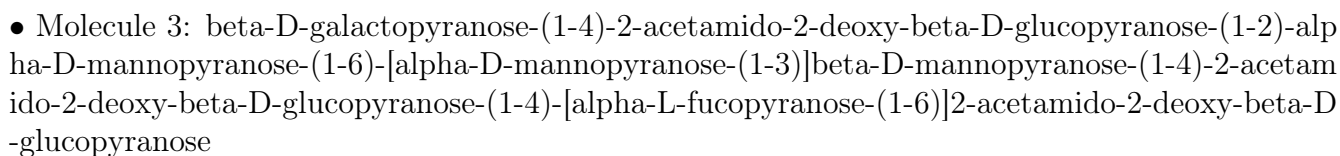
- Molecule 1: Ig gamma-1 chain C region



- Molecule 1: Ig gamma-1 chain C region



- Molecule 2: Glycoprotein E



Chain C:  75% 12% 12%

NAG1	NAG2	MAN3	MAN4	NAG5	GAL6	MAN7	FUC8
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● Molecule 3: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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

Chain D:  62% 38%

NAG1	NAG2	MAN3	MAN4	NAG5	GAL6	MAN7	FUC8
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4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	174.73Å 174.73Å 316.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.00 – 5.00 36.00 – 5.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (36.00-5.00) 99.1 (36.00-5.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.70 (at 5.08Å)	Xtriage
Refinement program		Depositor
R, R_{free}	(Not available) , (Not available) 0.499 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	210.3	Xtriage
Anisotropy	0.477	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 72.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.63	EDS
Total number of atoms	6148	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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: NAG, GAL, MAN, FUC, 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	0.47	0/1705	0.87	0/2322
1	B	0.48	0/1705	0.87	3/2322 (0.1%)
2	E	0.38	0/1359	1.03	12/1866 (0.6%)
2	F	0.38	0/1359	1.02	12/1866 (0.6%)
All	All	0.43	0/6128	0.94	27/8376 (0.3%)

There are no bond length outliers.

The worst 5 of 27 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	319	PRO	CA-C-N	11.73	132.46	120.38
2	F	319	PRO	C-N-CA	11.73	132.46	120.38
2	E	319	PRO	CA-C-N	11.72	132.45	120.38
2	E	319	PRO	C-N-CA	11.72	132.45	120.38
2	E	318	ASN	CA-C-N	7.03	127.62	120.38

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	1659	0	1622	137	0
1	B	1659	0	1623	82	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	1319	0	1239	77	30
2	F	1319	0	1238	138	30
3	C	96	0	82	1	0
3	D	96	0	82	0	0
All	All	6148	0	5886	255	30

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 21.

The worst 5 of 255 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:310:HIS:CD2	2:F:321:PRO:HB3	1.12	1.60
1:A:310:HIS:CG	2:F:321:PRO:CB	1.85	1.60
1:A:310:HIS:CG	2:F:321:PRO:HB3	1.07	1.56
1:A:314:LEU:HB2	2:F:322:ARG:NH1	1.36	1.38
1:A:310:HIS:CD2	2:F:321:PRO:CB	2.01	1.34

The worst 5 of 30 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)
2:F:305:ARG:NH2	2:E:285:GLN:CB[3_444]	0.59	1.61
2:F:285:GLN:CB	2:E:305:ARG:NH2[3_444]	0.81	1.39
2:F:305:ARG:NH1	2:E:285:GLN:O[3_444]	1.28	0.92
2:F:303:THR:CB	2:E:303:THR:CB[3_444]	1.37	0.83
2:F:305:ARG:NH1	2:E:285:GLN:C[3_444]	1.40	0.80

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	205/227 (90%)	203 (99%)	2 (1%)	0	100	100
1	B	205/227 (90%)	199 (97%)	5 (2%)	1 (0%)	24	63
2	E	169/401 (42%)	160 (95%)	6 (4%)	3 (2%)	6	32
2	F	169/401 (42%)	160 (95%)	6 (4%)	3 (2%)	6	32
All	All	748/1256 (60%)	722 (96%)	19 (2%)	7 (1%)	14	50

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	295	ALA
2	F	320	PRO
2	F	339	ALA
2	E	295	ALA
2	E	320	PRO

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	193/210 (92%)	192 (100%)	1 (0%)	81	82
1	B	193/210 (92%)	188 (97%)	5 (3%)	40	61
2	E	144/332 (43%)	142 (99%)	2 (1%)	59	72
2	F	144/332 (43%)	142 (99%)	2 (1%)	59	72
All	All	674/1084 (62%)	664 (98%)	10 (2%)	57	71

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	443	LEU
2	E	288	GLU
2	E	320	PRO
1	B	270	ASP
1	B	328	LEU

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	B	429	HIS
2	E	338	GLN
2	E	383	ASN
2	E	353	HIS
2	E	283	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 ⓘ

16 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	C	1	1,3	14,14,15	0.48	0	17,19,21	0.75	0
3	NAG	C	2	3	14,14,15	0.46	0	17,19,21	0.76	0
3	BMA	C	3	3	11,11,12	0.63	0	15,15,17	0.54	0
3	MAN	C	4	3	11,11,12	0.51	0	15,15,17	0.56	0
3	NAG	C	5	3	14,14,15	0.49	0	17,19,21	0.78	0
3	GAL	C	6	3	11,11,12	0.74	0	15,15,17	0.70	0
3	MAN	C	7	3	11,11,12	0.60	0	15,15,17	0.74	0
3	FUC	C	8	3	10,10,11	0.73	0	14,14,16	1.08	2 (14%)
3	NAG	D	1	1,3	14,14,15	0.69	0	17,19,21	0.89	0
3	NAG	D	2	3	14,14,15	0.59	0	17,19,21	0.81	1 (5%)
3	BMA	D	3	3	11,11,12	0.78	0	15,15,17	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MAN	D	4	3	11,11,12	0.56	0	15,15,17	0.85	1 (6%)
3	NAG	D	5	3	14,14,15	0.48	0	17,19,21	0.82	0
3	GAL	D	6	3	11,11,12	0.55	0	15,15,17	0.48	0
3	MAN	D	7	3	11,11,12	0.53	0	15,15,17	0.52	0
3	FUC	D	8	3	10,10,11	0.75	0	14,14,16	0.89	1 (7%)

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

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	MAN	C1-O5-C5	2.63	115.71	112.19
3	C	8	FUC	C3-C4-C5	2.53	113.65	109.81
3	D	2	NAG	C2-N2-C7	-2.23	119.91	122.90
3	D	8	FUC	C1-C2-C3	2.18	112.83	109.64
3	C	8	FUC	C1-C2-C3	2.06	112.65	109.64

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	C	8	FUC	C1
3	D	8	FUC	C1

5 of 11 torsion outliers are listed below:

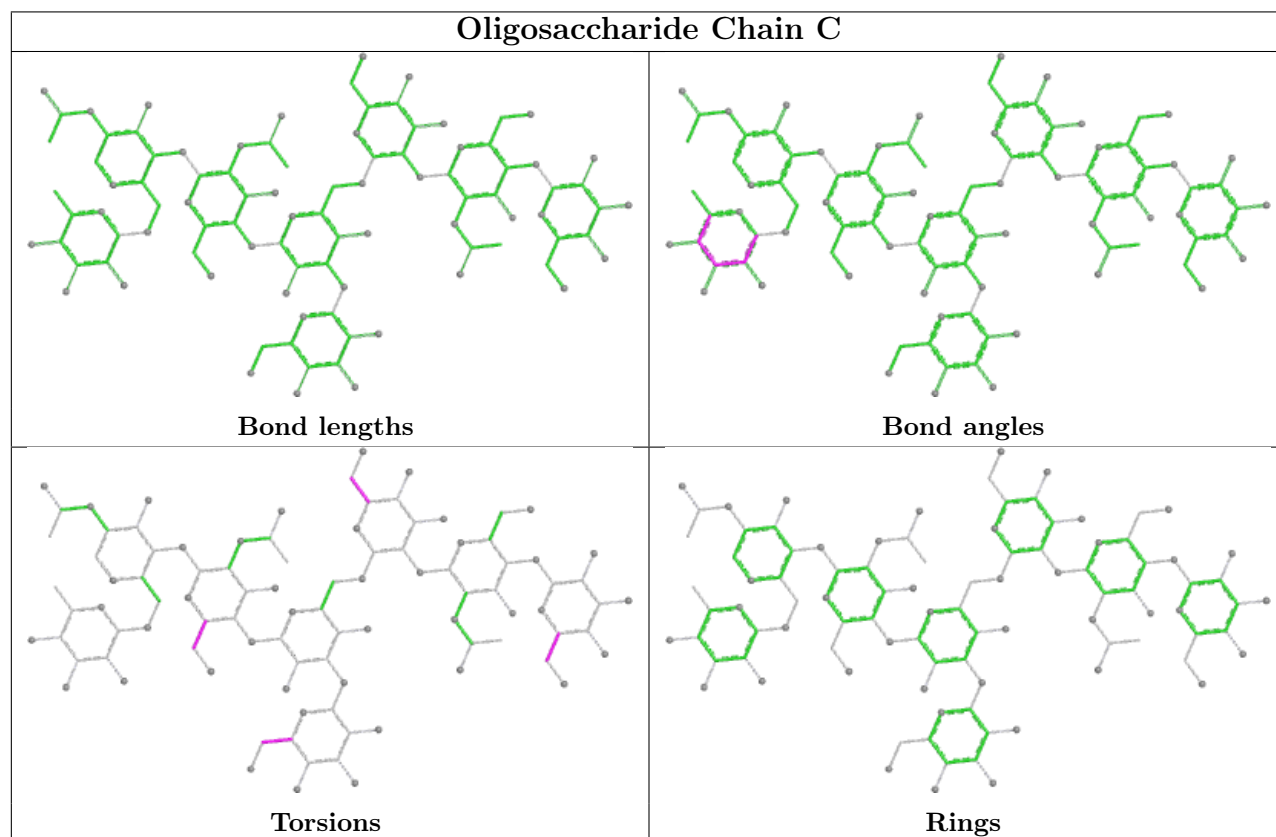
Mol	Chain	Res	Type	Atoms
3	D	4	MAN	O5-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6
3	C	2	NAG	C4-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6

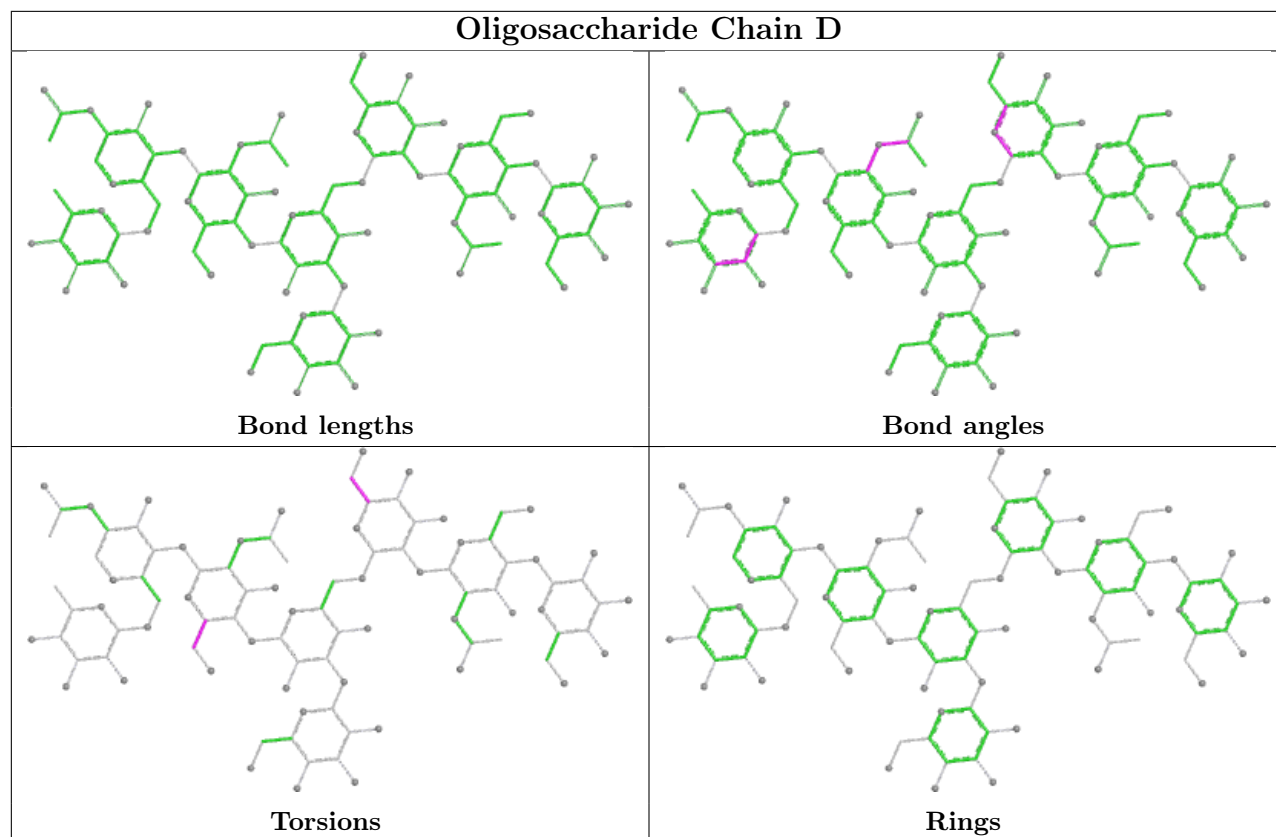
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	8	FUC	1	0
3	C	1	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](#)

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	207/227 (91%)	2.38	103 (49%) 0 2	9, 23, 36, 50	5 (2%)
1	B	207/227 (91%)	2.10	97 (46%) 0 2	10, 24, 45, 56	4 (1%)
2	E	171/401 (42%)	2.05	72 (42%) 0 3	11, 18, 28, 36	0
2	F	171/401 (42%)	1.67	53 (30%) 1 5	0, 0, 0, 0	0
All	All	756/1256 (60%)	2.07	325 (42%) 0 3	0, 19, 37, 56	9 (1%)

The worst 5 of 325 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	265	ASP	11.8
2	E	390	LEU	11.7
1	A	238	PRO	10.8
2	F	390	LEU	9.2
1	A	237	GLY	8.5

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	NAG	C	1	14/15	-	-	31,33,43,54	0
3	NAG	C	2	14/15	-	-	23,27,30,31	0

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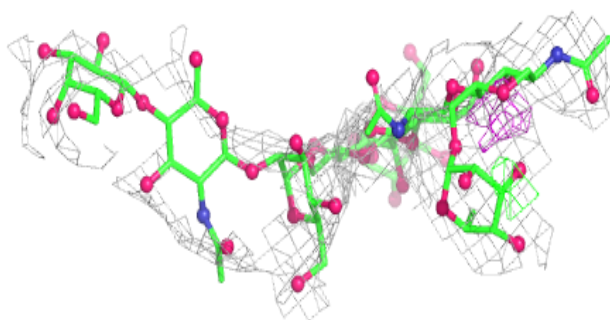
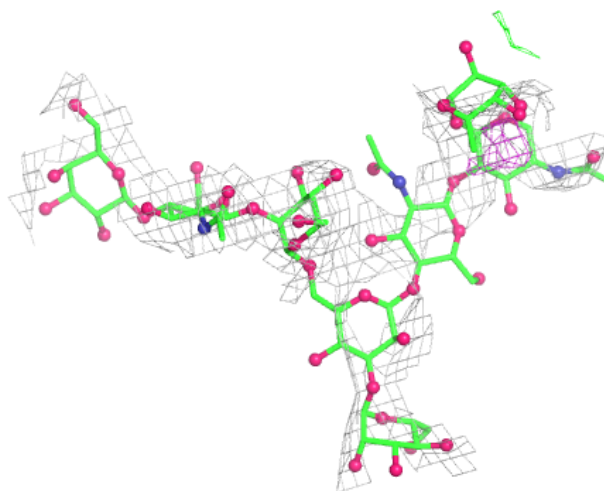
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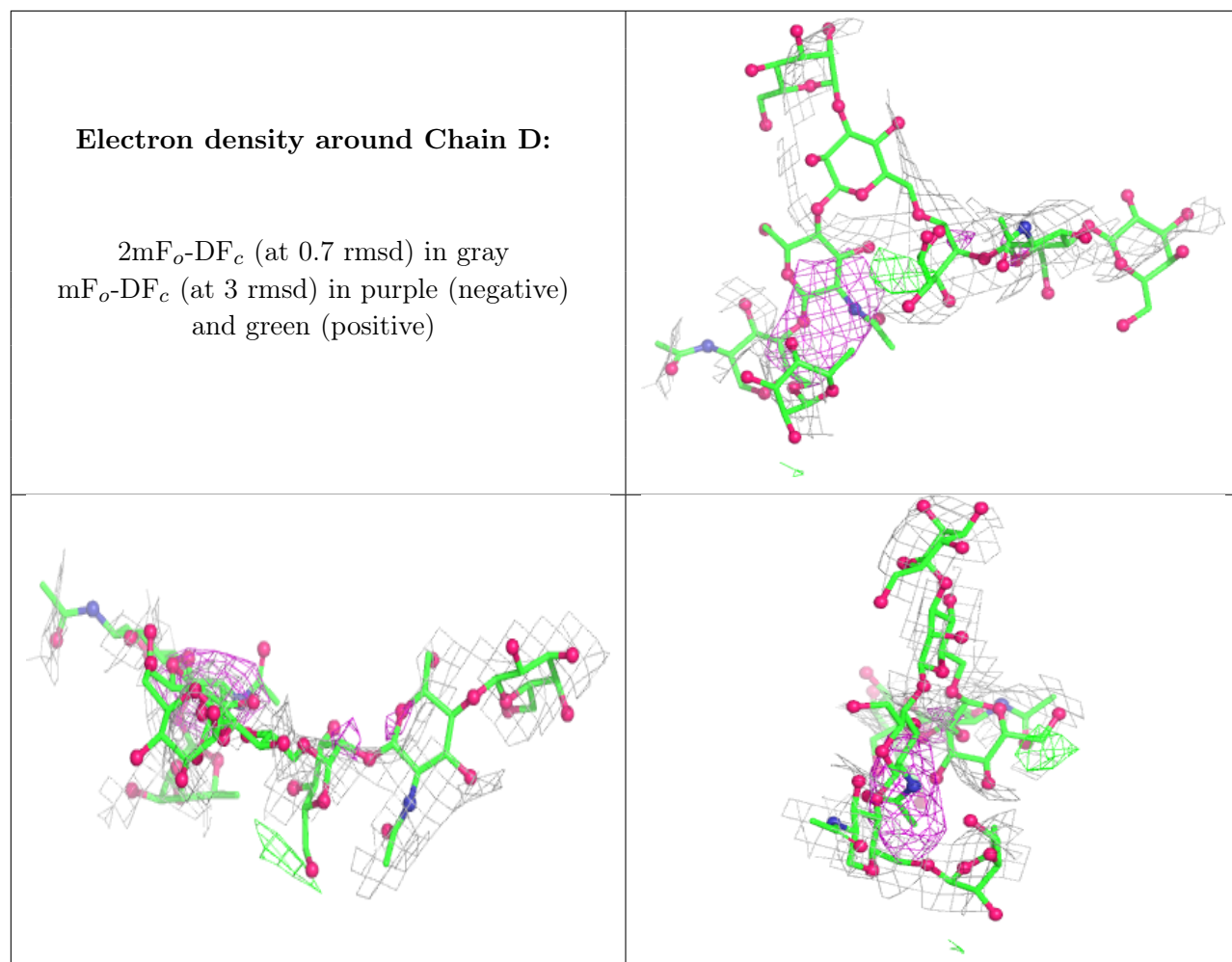
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	BMA	C	3	11/12	-	-	30,33,36,40	0
3	MAN	C	4	11/12	-	-	36,37,40,40	0
3	NAG	C	5	14/15	-	-	33,36,39,42	0
3	GAL	C	6	11/12	-	-	47,53,56,61	0
3	MAN	C	7	11/12	-	-	39,44,46,47	0
3	FUC	C	8	10/11	-	-	62,68,74,74	0
3	NAG	D	1	14/15	-	-	40,42,45,54	0
3	NAG	D	2	14/15	-	-	29,34,36,38	0
3	BMA	D	3	11/12	-	-	25,28,33,34	0
3	MAN	D	4	11/12	-	-	27,29,32,36	0
3	NAG	D	5	14/15	-	-	29,33,37,43	0
3	GAL	D	6	11/12	-	-	48,52,55,58	0
3	MAN	D	7	11/12	-	-	40,42,43,44	0
3	FUC	D	8	10/11	-	-	57,60,63,64	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 C:

$2mF_o-DF_c$ (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.