wwPDB X-ray Structure Validation Summary Report (i)

Nov 23, 2022 – 03:28 pm GMT

PDB ID	:	8A47
Title	:	IdeS in complex with IgG1 Fc
Authors	:	Sudol, A.S.L.; Tews, I.; Crispin, M.
Deposited on	:	2022-06-10
Resolution	:	2.34 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 (i) 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 (i)) were used in the production of this report:

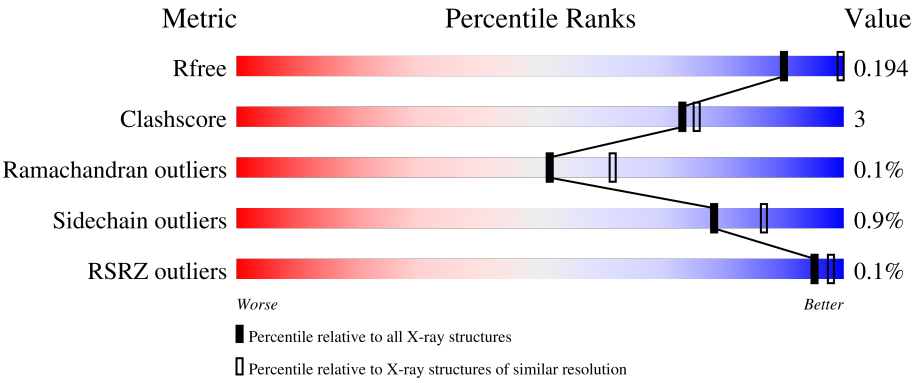
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.3
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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}	130704	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (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 <=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	90% 6% ·
1	B	227	89% 5% 5%
2	C	308	82% 14% · ·
3	D	7	71% 29%
4	E	8	100%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 12063 atoms, of which 5878 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 IgG1 Fc.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	217	Total	C	H	N	O	S	107	1	0
			3435	1102	1704	290	331	8			
1	B	215	Total	C	H	N	O	S	106	0	0
			3396	1089	1687	287	326	7			

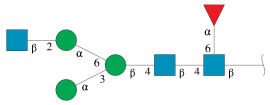
- Molecule 2 is a protein called IgG-degrading protease.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	297	Total	C	H	N	O	S	145	0	0
			4674	1501	2314	402	452	5			

There are 10 discrepancies between the modelled and reference sequences:

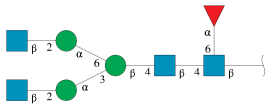
Chain	Residue	Modelled	Actual	Comment	Reference
C	40	MET	-	initiating methionine	UNP A0A8B6IYA1
C	94	ALA	CYS	engineered mutation	UNP A0A8B6IYA1
C	350	LEU	-	expression tag	UNP A0A8B6IYA1
C	351	GLU	-	expression tag	UNP A0A8B6IYA1
C	352	HIS	-	expression tag	UNP A0A8B6IYA1
C	353	HIS	-	expression tag	UNP A0A8B6IYA1
C	354	HIS	-	expression tag	UNP A0A8B6IYA1
C	355	HIS	-	expression tag	UNP A0A8B6IYA1
C	356	HIS	-	expression tag	UNP A0A8B6IYA1
C	357	HIS	-	expression tag	UNP A0A8B6IYA1

- Molecule 3 is an oligosaccharide called 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	D	7	Total	C	H	N	O	18	0	0
			165	48	80	3	34			

- Molecule 4 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
4	E	8	Total	C	H	N	O	20	0	0
			192	56	93	4	39			

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	Na	0	0
			1	1		

- Molecule 6 is water.

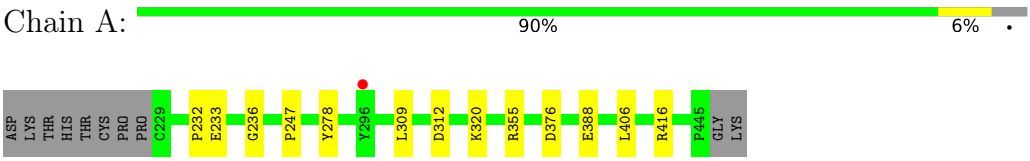
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	45	Total	O	0	0
			45	45		
6	B	65	Total	O	0	0
			65	65		
6	C	90	Total	O	0	0
			90	90		

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.

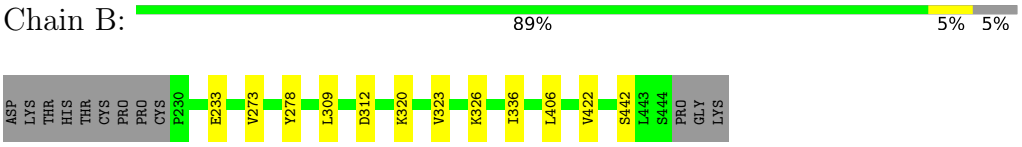
• Molecule 1: IgG1 Fc

Chain A:



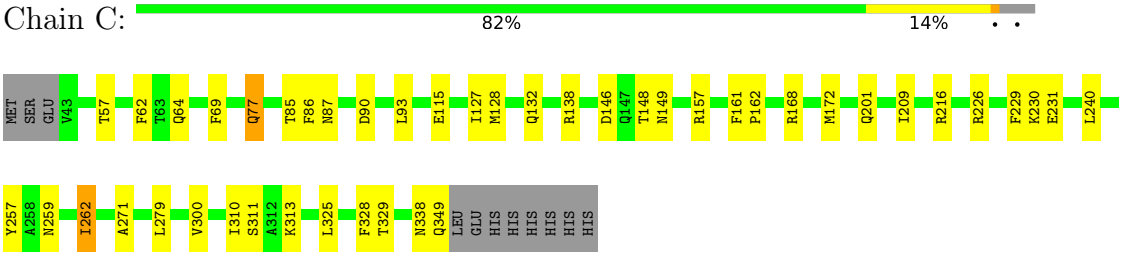
• Molecule 1: IgG1 Fc

Chain B:



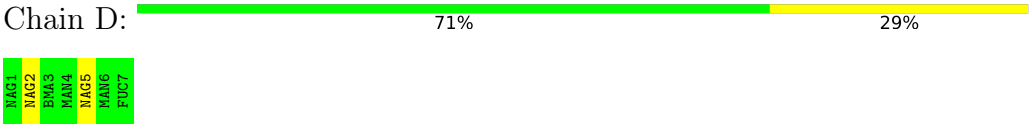
• Molecule 2: IgG-degrading protease

Chain C:



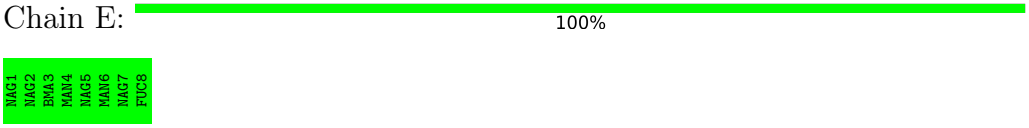
• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:



• Molecule 4: 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

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	217.56Å 108.45Å 63.09Å 90.00° 90.06° 90.00°	Depositor
Resolution (Å)	48.53 – 2.34 48.53 – 2.34	Depositor EDS
% Data completeness (in resolution range)	97.4 (48.53-2.34) 95.2 (48.53-2.34)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.99 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.171 , 0.203 0.163 , 0.194	Depositor DCC
R_{free} test set	3049 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	34.8	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.37$, $< L^2 > = 0.20$	Xtriage
Estimated twinning fraction	0.427 for -h,-k,l	Xtriage
Reported twinning fraction	0.507 for H, K, L 0.493 for -h,-k,l	Depositor
Outliers	0 of 60351 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	12063	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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, NAG, MAN, NA, 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	0.40	0/1780	0.70	0/2427
1	B	0.43	0/1757	0.73	0/2394
2	C	0.40	0/2412	0.77	0/3263
All	All	0.41	0/5949	0.74	0/8084

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	1731	1704	1696	8	1
1	B	1709	1687	1680	6	1
2	C	2360	2314	2303	26	0
3	D	85	80	73	0	0
4	E	99	93	85	0	0
5	C	1	0	0	0	0
6	A	45	0	0	2	0
6	B	65	0	0	2	0
6	C	90	0	0	3	0
All	All	6185	5878	5837	39	1

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

The worst 5 of 39 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)
2:C:62:PHE:HB3	2:C:69:PHE:CZ	2.31	0.65
2:C:64:GLN:NE2	6:C:501:HOH:O	2.20	0.64
1:B:422:VAL:HG22	1:B:442:SER:HB2	1.81	0.63
2:C:271:ALA:HB1	2:C:279:LEU:HD21	1.79	0.63
2:C:57:THR:HB	2:C:77:GLN:HG3	1.82	0.61

All (1) 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:A:416:ARG:HH11	1:B:233:GLU:OE2[4_555]	1.58	0.02

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	216/227 (95%)	213 (99%)	2 (1%)	1 (0%)	29 31
1	B	213/227 (94%)	211 (99%)	2 (1%)	0	100 100
2	C	295/308 (96%)	280 (95%)	15 (5%)	0	100 100
All	All	724/762 (95%)	704 (97%)	19 (3%)	1 (0%)	51 62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	GLY

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	201/209 (96%)	201 (100%)	0	100 100
1	B	198/209 (95%)	197 (100%)	1 (0%)	88 93
2	C	252/264 (96%)	247 (98%)	5 (2%)	55 66
All	All	651/682 (96%)	645 (99%)	6 (1%)	78 87

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	168	ARG
2	C	257	TYR
2	C	262	ILE
2	C	77	GLN
1	B	336	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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

15 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	1,3	14,14,15	0.86	0	17,19,21	1.03	0
3	NAG	D	2	3	14,14,15	1.19	2 (14%)	17,19,21	0.59	0
3	BMA	D	3	3	11,11,12	0.96	0	15,15,17	0.82	0
3	MAN	D	4	3	11,11,12	0.96	0	15,15,17	1.01	0
3	NAG	D	5	3	14,14,15	0.84	0	17,19,21	1.03	1 (5%)
3	MAN	D	6	3	11,11,12	0.63	0	15,15,17	0.58	0
3	FUC	D	7	3	10,10,11	1.03	0	14,14,16	1.06	0
4	NAG	E	1	1,4	14,14,15	0.88	0	17,19,21	0.72	0
4	NAG	E	2	4	14,14,15	0.58	0	17,19,21	0.62	0
4	BMA	E	3	4	11,11,12	0.76	0	15,15,17	0.88	0
4	MAN	E	4	4	11,11,12	0.96	0	15,15,17	0.79	0
4	NAG	E	5	4	14,14,15	0.62	0	17,19,21	0.75	0
4	MAN	E	6	4	11,11,12	1.09	0	15,15,17	1.04	0
4	NAG	E	7	4	14,14,15	0.70	0	17,19,21	0.81	0
4	FUC	E	8	4	10,10,11	0.68	0	14,14,16	0.96	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	D	1	1,3	-	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	-	0/2/19/22	0/1/1/1
3	NAG	D	5	3	-	1/6/23/26	0/1/1/1
3	MAN	D	6	3	-	1/2/19/22	0/1/1/1
3	FUC	D	7	3	-	-	0/1/1/1
4	NAG	E	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	0/2/19/22	0/1/1/1
4	MAN	E	4	4	-	2/2/19/22	0/1/1/1
4	NAG	E	5	4	-	0/6/23/26	0/1/1/1
4	MAN	E	6	4	-	1/2/19/22	0/1/1/1
4	NAG	E	7	4	-	0/6/23/26	0/1/1/1
4	FUC	E	8	4	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	2	NAG	C4-C5	2.22	1.57	1.53
3	D	2	NAG	O4-C4	2.03	1.47	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	5	NAG	C6-C5-C4	2.38	118.57	113.00

There are no chirality outliers.

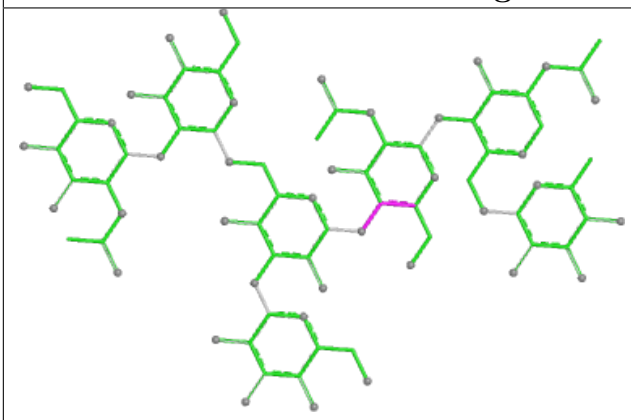
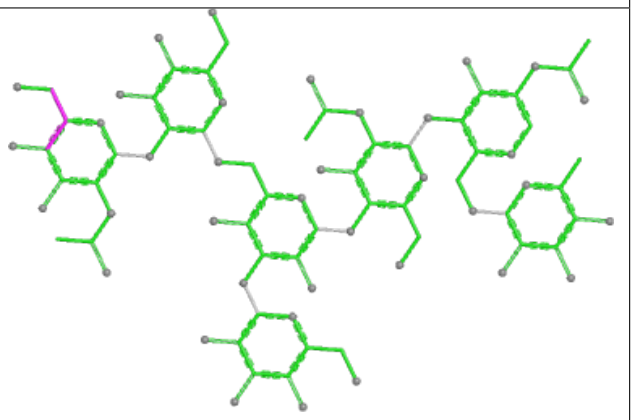
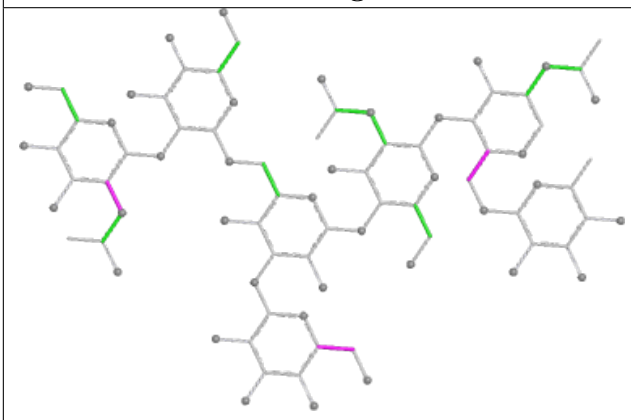
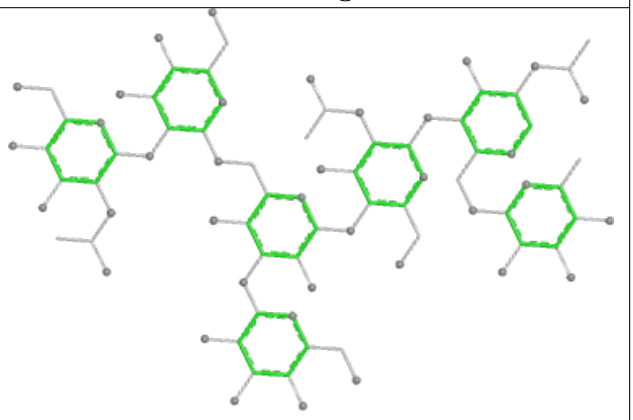
5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	4	MAN	O5-C5-C6-O6
4	E	4	MAN	C4-C5-C6-O6
4	E	1	NAG	C4-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6

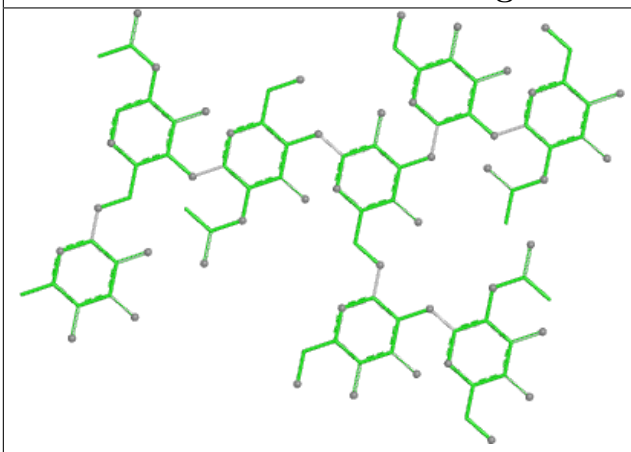
There are no ring outliers.

No monomer is involved in short contacts.

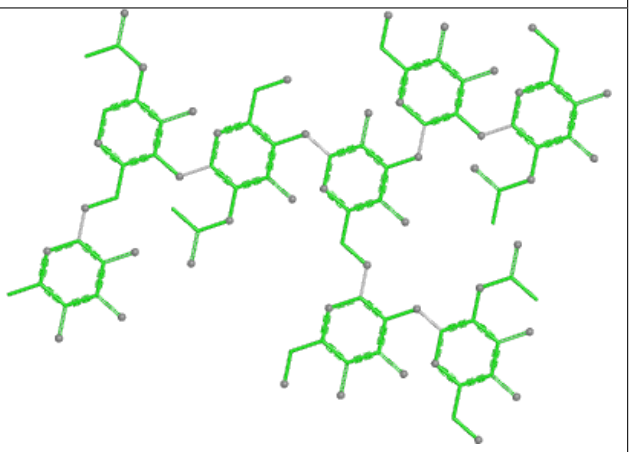
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain D	
	
Bond lengths	**Bond angles**
	
Torsions	**Rings**

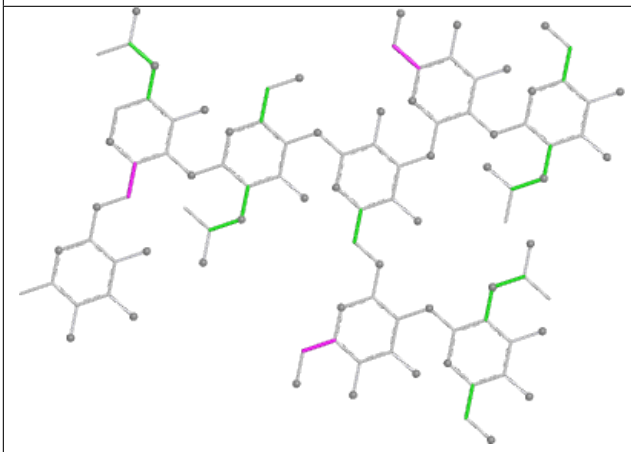
Oligosaccharide Chain E



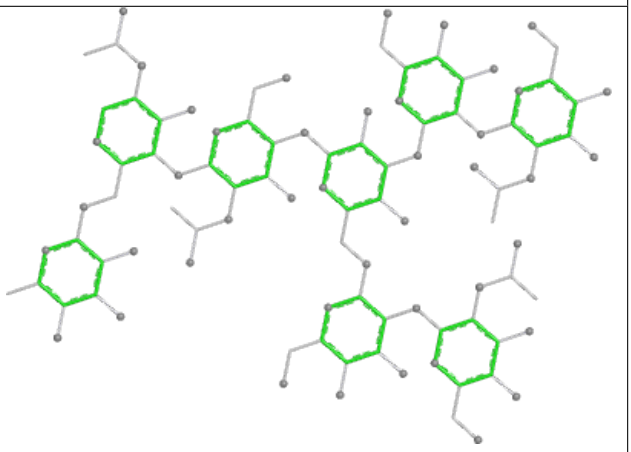
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

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, 95^{th} 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	217/227 (95%)	-0.04	1 (0%) 91 95	21, 43, 68, 87	0
1	B	215/227 (94%)	-0.12	0 100 100	21, 32, 52, 75	0
2	C	297/308 (96%)	-0.15	0 100 100	18, 36, 56, 92	0
All	All	729/762 (95%)	-0.11	1 (0%) 95 98	18, 36, 63, 92	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	296	TYR	2.6

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

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, 95^{th} 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	FUC	D	7	10/11	0.92	0.14	43,54,57,60	3
3	MAN	D	4	11/12	0.95	0.18	42,50,58,63	3
3	NAG	D	2	14/15	0.95	0.17	40,56,63,68	2
4	NAG	E	5	14/15	0.95	0.16	39,55,67,70	3
3	NAG	D	5	14/15	0.96	0.21	46,56,60,60	3
3	MAN	D	6	11/12	0.97	0.14	41,52,58,59	4
3	NAG	D	1	14/15	0.97	0.15	39,48,53,56	1

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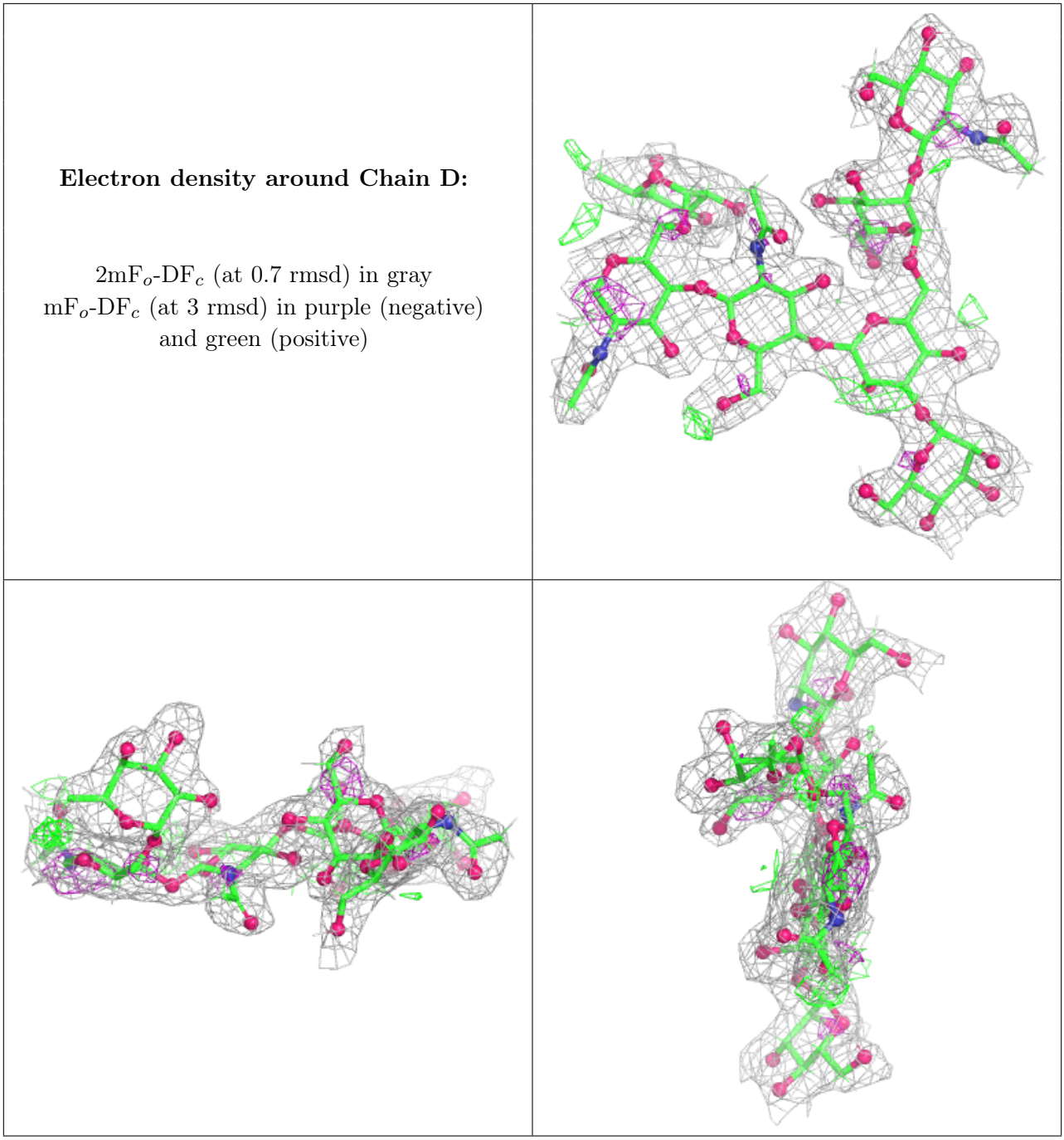
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	BMA	D	3	11/12	0.97	0.09	40,47,50,50	2
4	MAN	E	4	11/12	0.98	0.11	36,43,48,50	3
4	MAN	E	6	11/12	0.98	0.11	31,34,37,37	3
4	NAG	E	7	14/15	0.98	0.11	31,36,43,45	3
4	FUC	E	8	10/11	0.98	0.13	35,43,48,48	3
4	BMA	E	3	11/12	0.99	0.10	29,34,37,39	2
4	NAG	E	1	14/15	0.99	0.13	27,29,31,35	1
4	NAG	E	2	14/15	0.99	0.12	29,31,33,34	2

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

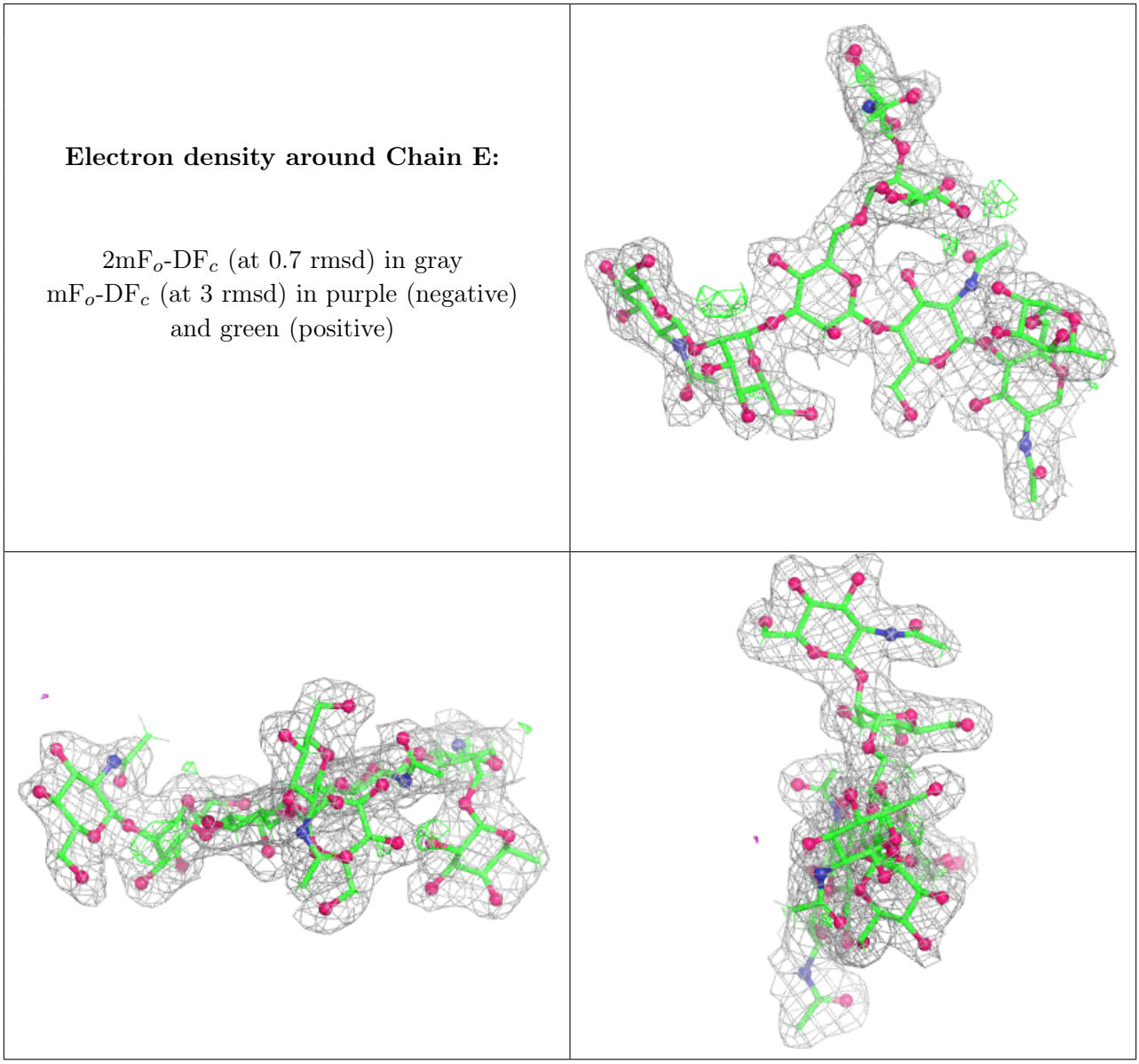
Electron density around Chain D:

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



Electron density around Chain E:

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



6.4 Ligands

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, 95^{th} 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
5	NA	C	401	1/1	0.98	0.13	29,29,29,29	0

6.5 Other polymers

There are no such residues in this entry.