



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2021 – 02:15 PM EDT

PDB ID : 2NT6
Title : Crystal structure of pectin methylesterase D178A mutant in complex with hexasaccharide III
Authors : Fries, M.; Brocklehurst, K.; Shevchik, V.E.; Pickersgill, R.W.
Deposited on : 2006-11-07
Resolution : 1.70 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

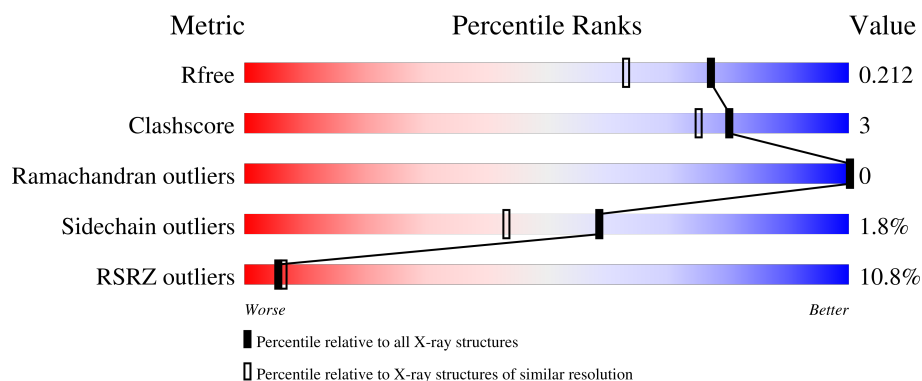
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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	342	<div> <div>11%</div> <div> <div></div> <div>93%</div> <div>6%</div> </div> </div>
1	B	342	<div> <div>11%</div> <div> <div></div> <div>93%</div> <div>6%</div> </div> </div>
2	C	6	<div> <div>50%</div> <div>50%</div> </div>
2	D	6	<div> <div>67%</div> <div>17%</div> <div>17%</div> </div>

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
2	ADA	C	5	X	-	-	X
2	M8C	C	6	-	-	-	X
2	ADA	D	5	-	-	-	X
2	M8C	D	6	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6145 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 Pectinesterase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	0	0
			2606	1628	452	522	4			
1	B	342	Total	C	N	O	S	0	0	0
			2606	1628	452	522	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	178	ALA	ASP	engineered mutation	UNP P0C1A9
B	178	ALA	ASP	engineered mutation	UNP P0C1A9

- Molecule 2 is an oligosaccharide called methyl alpha-D-galactopyranuronate-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-methyl alpha-D-galactopyranuronate.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	6	Total	C	O	0	0	0
			75	38	37			
2	D	6	Total	C	O	0	0	0
			75	38	37			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	395	Total	O	0	0
			395	395		

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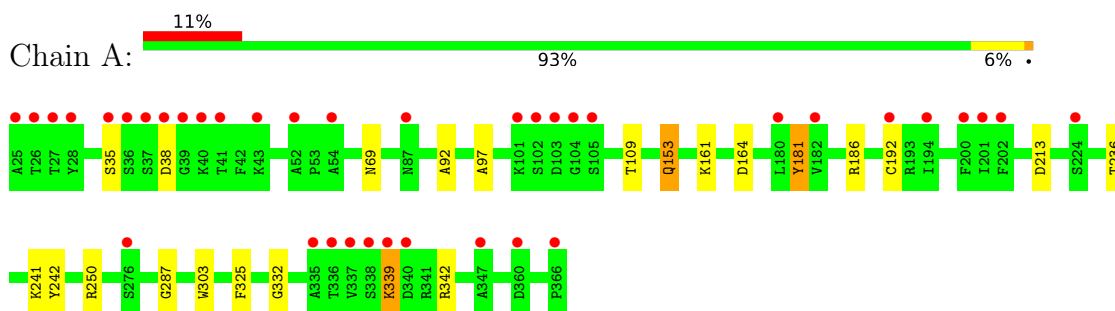
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	388	Total 388	O 388	0	0

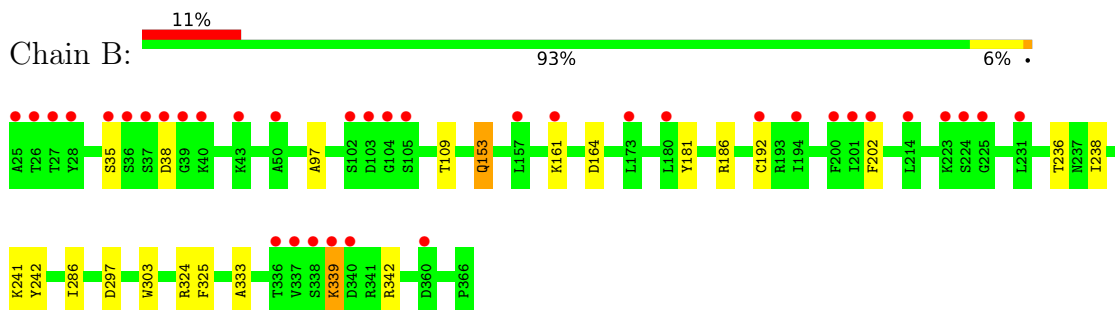
3 Residue-property plots

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: Pectinesterase A



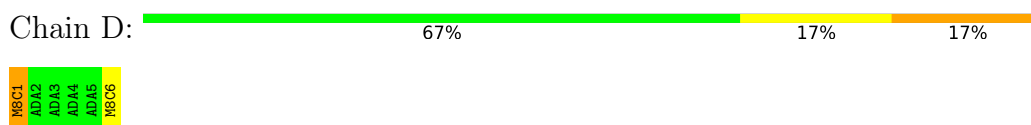
- Molecule 1: Pectinesterase A



- Molecule 2: methyl alpha-D-galactopyranuronate-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-methyl alpha-D-galactopyranuronate



- Molecule 2: methyl alpha-D-galactopyranuronate-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-methyl alpha-D-galactopyranuronate



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.07Å 85.12Å 97.38Å 90.00° 93.37° 90.00°	Depositor
Resolution (Å)	40.66 – 1.70 33.36 – 1.70	Depositor EDS
% Data completeness (in resolution range)	95.1 (40.66-1.70) 95.1 (33.36-1.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.64 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.190 , 0.212 0.189 , 0.212	Depositor DCC
R_{free} test set	4358 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	13.3	Xtriage
Anisotropy	0.137	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 63.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6145	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.15% 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: M8C, ADA

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	0.38	0/2660	0.55	0/3607
1	B	0.38	0/2660	0.54	0/3607
All	All	0.38	0/5320	0.54	0/7214

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2606	0	2525	14	0
1	B	2606	0	2525	15	0
2	C	75	0	45	0	0
2	D	75	0	45	1	0
3	A	395	0	0	4	0
3	B	388	0	0	6	0
All	All	6145	0	5140	28	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 3.

All (28) 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:339:LYS:H	1:A:339:LYS:HE2	1.43	0.83
1:B:339:LYS:H	1:B:339:LYS:HE2	1.44	0.82
1:A:109:THR:OG1	1:A:153:GLN:HG3	1.81	0.81
1:B:109:THR:OG1	1:B:153:GLN:HG3	1.87	0.75
1:A:164:ASP:OD1	1:A:186:ARG:HD3	1.99	0.62
1:B:342:ARG:HD3	3:B:633:HOH:O	2.01	0.60
1:B:164:ASP:OD1	1:B:186:ARG:HD3	2.00	0.60
1:A:342:ARG:HD3	3:A:662:HOH:O	2.01	0.59
1:A:97:ALA:HB2	1:A:153:GLN:HG2	1.86	0.57
1:A:181:TYR:HE2	3:A:503:HOH:O	1.88	0.56
1:B:97:ALA:HB2	1:B:153:GLN:HG2	1.89	0.55
1:B:161:LYS:HG3	3:B:532:HOH:O	2.08	0.54
1:A:35:SER:HB2	1:A:38:ASP:HB2	1.89	0.54
1:B:35:SER:HB2	1:B:38:ASP:HB2	1.89	0.54
1:A:241:LYS:HG3	1:A:242:TYR:CE2	2.48	0.48
1:A:161:LYS:HG3	3:A:515:HOH:O	2.14	0.47
1:B:241:LYS:HG3	1:B:242:TYR:CE2	2.51	0.45
1:B:297:ASP:HB2	3:B:817:HOH:O	2.17	0.43
1:A:332:GLY:O	1:B:333:ALA:HA	2.18	0.43
1:A:213:ASP:OD1	1:A:250:ARG:NE	2.46	0.43
1:A:303:TRP:CE2	1:A:325:PHE:HB3	2.54	0.42
1:B:238:ILE:CD1	3:B:550:HOH:O	2.68	0.42
1:A:287:GLY:N	3:A:506:HOH:O	2.53	0.42
1:B:303:TRP:CE2	1:B:325:PHE:HB3	2.55	0.41
1:B:202:PHE:CD2	2:D:1:M8C:HH31	2.56	0.41
1:B:286:ILE:CG2	3:B:550:HOH:O	2.68	0.41
1:A:69:ASN:HA	1:A:92:ALA:O	2.21	0.40
1:B:324:ARG:HG3	3:B:550:HOH:O	2.21	0.40

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	340/342 (99%)	336 (99%)	4 (1%)	0	100	100
1	B	340/342 (99%)	336 (99%)	4 (1%)	0	100	100
All	All	680/684 (99%)	672 (99%)	8 (1%)	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	280/280 (100%)	275 (98%)	5 (2%)	59	43
1	B	280/280 (100%)	275 (98%)	5 (2%)	59	43
All	All	560/560 (100%)	550 (98%)	10 (2%)	59	43

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	GLN
1	A	181	TYR
1	A	192	CYS
1	A	236	THR
1	A	339	LYS
1	B	153	GLN
1	B	181	TYR
1	B	192	CYS
1	B	236	THR
1	B	339	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	87	ASN
1	A	210	ASN

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Mol	Chain	Res	Type
1	A	248	ASN
1	B	69	ASN
1	B	87	ASN
1	B	210	ASN
1	B	248	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 ⓘ

12 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	M8C	C	1	2	14,14,14	1.45	1 (7%)	20,20,20	1.34	2 (10%)
2	ADA	C	2	2	9,12,13	0.49	0	12,17,19	0.72	0
2	ADA	C	3	2	9,12,13	0.33	0	12,17,19	0.62	0
2	ADA	C	4	2	9,12,13	0.42	0	12,17,19	0.98	1 (8%)
2	ADA	C	5	2	9,12,13	0.41	0	12,17,19	0.57	0
2	M8C	C	6	2	13,13,14	1.55	1 (7%)	16,18,20	1.57	3 (18%)
2	M8C	D	1	2	14,14,14	1.45	1 (7%)	20,20,20	1.34	2 (10%)
2	ADA	D	2	2	9,12,13	0.45	0	12,17,19	0.61	0
2	ADA	D	3	2	9,12,13	0.37	0	12,17,19	0.55	0
2	ADA	D	4	2	9,12,13	0.36	0	12,17,19	0.70	0
2	ADA	D	5	2	9,12,13	0.46	0	12,17,19	0.69	0
2	M8C	D	6	2	13,13,14	1.57	1 (7%)	16,18,20	1.60	3 (18%)

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	M8C	C	1	2	-	2/6/26/26	0/1/1/1
2	ADA	C	2	2	-	0/0/21/24	0/1/1/1
2	ADA	C	3	2	-	0/0/21/24	0/1/1/1
2	ADA	C	4	2	-	0/0/21/24	0/1/1/1
2	ADA	C	5	2	1/1/5/6	0/0/21/24	0/1/1/1
2	M8C	C	6	2	-	2/6/23/26	0/1/1/1
2	M8C	D	1	2	-	2/6/26/26	0/1/1/1
2	ADA	D	2	2	-	0/0/21/24	0/1/1/1
2	ADA	D	3	2	-	0/0/21/24	0/1/1/1
2	ADA	D	4	2	-	0/0/21/24	0/1/1/1
2	ADA	D	5	2	-	0/0/21/24	0/1/1/1
2	M8C	D	6	2	-	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)
2	D	6	M8C	O6B-C6	5.31	1.46	1.33
2	C	6	M8C	O6B-C6	5.26	1.46	1.33
2	C	1	M8C	O6B-C6	4.95	1.45	1.33
2	D	1	M8C	O6B-C6	4.89	1.45	1.33

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	6	M8C	O6B-C6-C5	4.83	120.10	110.58
2	C	6	M8C	O6B-C6-C5	4.82	120.08	110.58
2	D	1	M8C	O6B-C6-C5	4.74	119.91	110.58
2	C	1	M8C	O6B-C6-C5	4.65	119.74	110.58
2	C	4	ADA	C1-C2-C3	2.83	113.15	109.67
2	D	1	M8C	O6B-C6-O6A	-2.53	118.89	123.84
2	D	6	M8C	O6B-C6-O6A	-2.50	118.95	123.84
2	C	1	M8C	O6B-C6-O6A	-2.46	119.04	123.84
2	C	6	M8C	O6B-C6-O6A	-2.36	119.22	123.84
2	C	6	M8C	CH3-O6B-C6	2.08	120.65	115.94
2	D	6	M8C	CH3-O6B-C6	2.03	120.53	115.94

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	5	ADA	C1

All (8) torsion outliers are listed below:

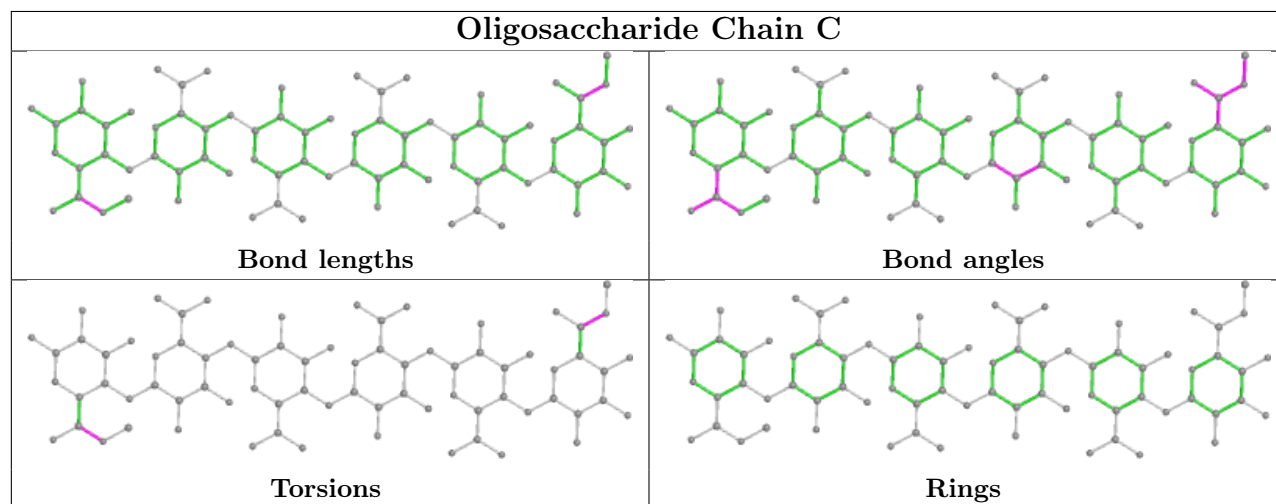
Mol	Chain	Res	Type	Atoms
2	C	1	M8C	C5-C6-O6B-CH3
2	C	6	M8C	C5-C6-O6B-CH3
2	D	6	M8C	O6A-C6-O6B-CH3
2	C	6	M8C	O6A-C6-O6B-CH3
2	D	6	M8C	C5-C6-O6B-CH3
2	D	1	M8C	C5-C6-O6B-CH3
2	C	1	M8C	O6A-C6-O6B-CH3
2	D	1	M8C	O6A-C6-O6B-CH3

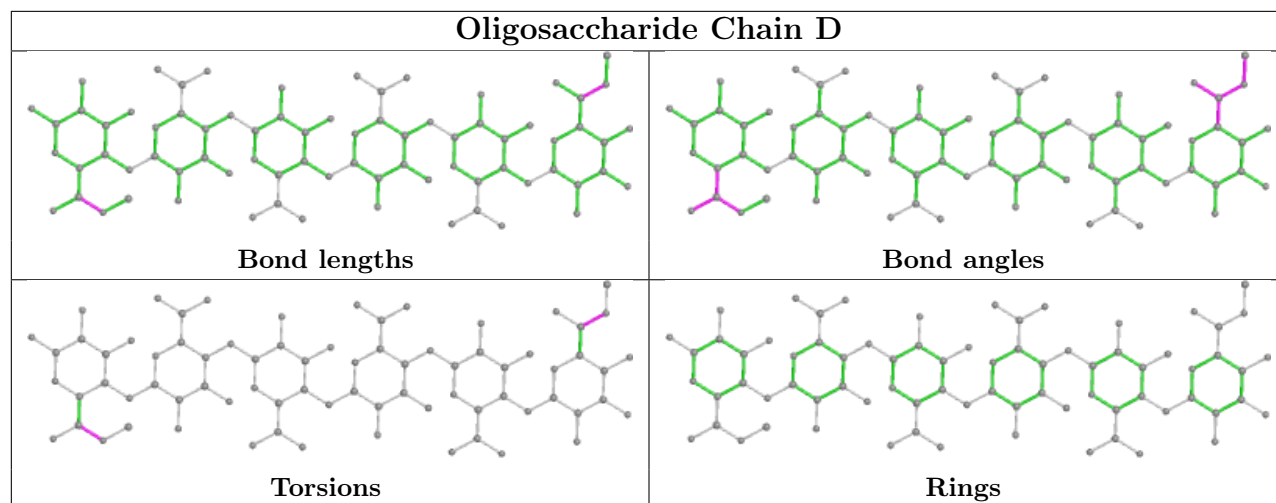
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	M8C	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 ⓘ

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	342/342 (100%)	0.74	38 (11%) 5 6	5, 12, 26, 43	1 (0%)
1	B	342/342 (100%)	0.70	36 (10%) 6 7	5, 12, 26, 43	1 (0%)
All	All	684/684 (100%)	0.72	74 (10%) 5 6	5, 12, 27, 43	2 (0%)

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	THR	15.7
1	A	37	SER	14.7
1	A	192	CYS	14.1
1	B	192	CYS	13.5
1	A	26	THR	13.3
1	B	25	ALA	12.8
1	B	37	SER	11.8
1	A	25	ALA	11.5
1	B	38	ASP	10.8
1	A	38	ASP	9.8
1	A	36	SER	8.7
1	A	39	GLY	8.6
1	A	40	LYS	7.9
1	B	36	SER	7.8
1	B	27	THR	7.5
1	B	39	GLY	7.4
1	B	40	LYS	6.3
1	B	35	SER	6.1
1	A	27	THR	6.0
1	A	35	SER	6.0
1	A	337	VAL	5.6
1	A	339	LYS	4.7
1	B	339	LYS	4.4
1	B	102	SER	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	224	SER	4.1
1	A	41	THR	4.0
1	A	102	SER	3.8
1	A	336	THR	3.5
1	A	28	TYR	3.4
1	B	360	ASP	3.4
1	B	103	ASP	3.3
1	A	224	SER	3.3
1	B	105	SER	3.2
1	B	201	ILE	3.1
1	B	43	LYS	3.1
1	A	180	LEU	3.0
1	A	338	SER	2.8
1	A	105	SER	2.8
1	B	28	TYR	2.8
1	A	103	ASP	2.8
1	A	360	ASP	2.8
1	A	201	ILE	2.7
1	A	43	LYS	2.7
1	A	182	VAL	2.7
1	B	104	GLY	2.6
1	B	200	PHE	2.6
1	B	340	ASP	2.6
1	B	336	THR	2.5
1	A	104	GLY	2.4
1	A	200	PHE	2.3
1	A	52	ALA	2.3
1	A	335	ALA	2.3
1	B	223	LYS	2.2
1	B	180	LEU	2.2
1	B	161	LYS	2.2
1	A	340	ASP	2.2
1	B	338	SER	2.2
1	B	337	VAL	2.2
1	A	194	ILE	2.2
1	B	194	ILE	2.2
1	B	231	LEU	2.1
1	A	54	ALA	2.1
1	A	347	ALA	2.1
1	B	157	LEU	2.1
1	A	202	PHE	2.1
1	B	202	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	101	LYS	2.1
1	B	50	ALA	2.1
1	B	173	LEU	2.1
1	B	214	LEU	2.1
1	A	276	SER	2.0
1	A	366	PRO	2.0
1	B	225	GLY	2.0
1	A	87	ASN	2.0

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 ⓘ

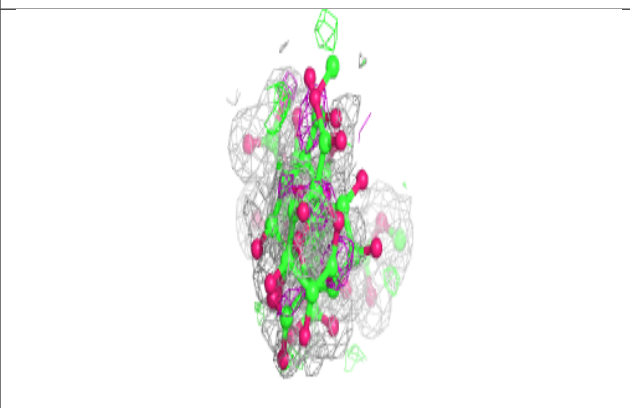
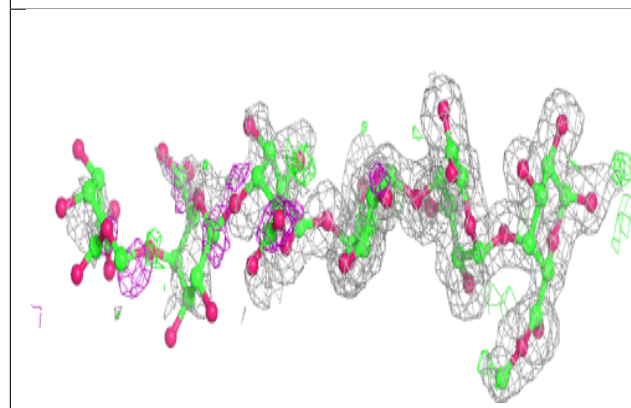
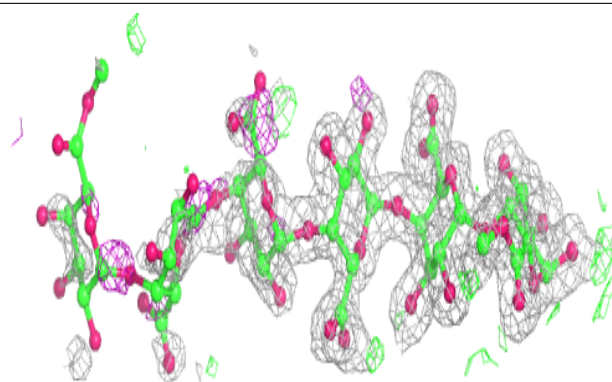
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	M8C	D	6	13/14	0.01	0.69	64,65,65,65	0
2	ADA	D	5	12/13	0.32	0.69	54,57,58,60	0
2	ADA	C	5	12/13	0.40	0.66	57,59,60,61	0
2	M8C	C	6	13/14	0.44	0.72	64,65,66,66	0
2	ADA	C	4	12/13	0.54	0.40	40,45,46,50	0
2	ADA	D	4	12/13	0.63	0.34	37,41,42,46	0
2	ADA	D	3	12/13	0.89	0.13	18,20,22,26	0
2	ADA	C	3	12/13	0.92	0.11	16,20,23,27	0
2	M8C	C	1	14/14	0.93	0.10	11,14,15,16	0
2	M8C	D	1	14/14	0.94	0.09	10,12,15,15	0
2	ADA	D	2	12/13	0.96	0.08	10,11,12,13	0
2	ADA	C	2	12/13	0.98	0.06	9,10,11,13	0

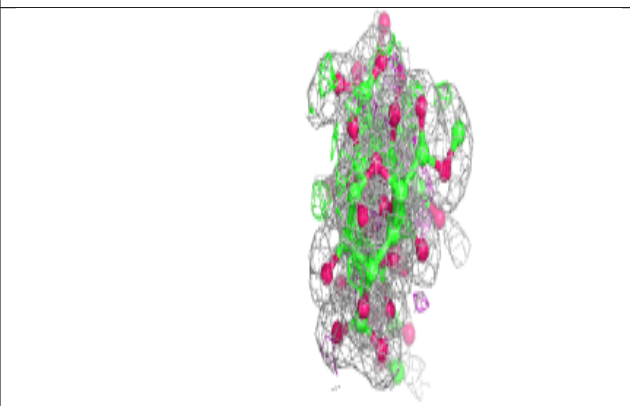
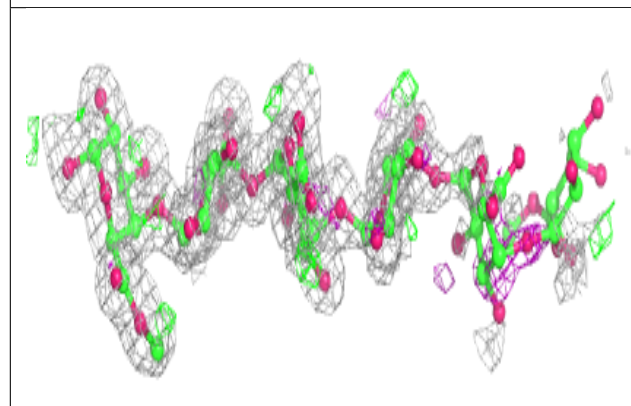
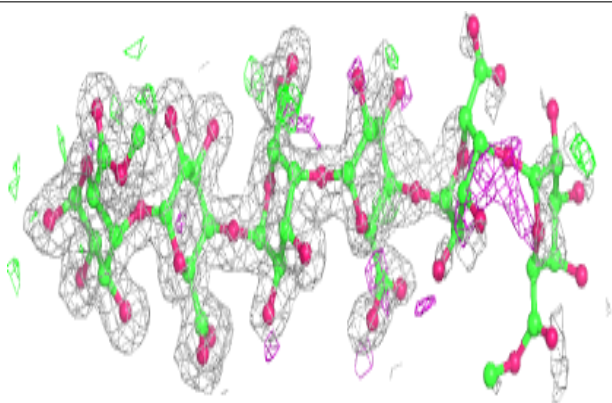
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

**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)



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.