



# Full wwPDB X-ray Structure Validation Report ⓘ

May 3, 2021 – 10:15 pm BST

PDB ID : 5E4T  
Title : Acetylcholinesterase Methylene Blue with PEG  
Authors : Dym, O.  
Deposited on : 2015-10-07  
Resolution : 2.43 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](mailto: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.

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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 : 2.18  
buster-report : 1.1.7 (2018)  
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) : 2.18

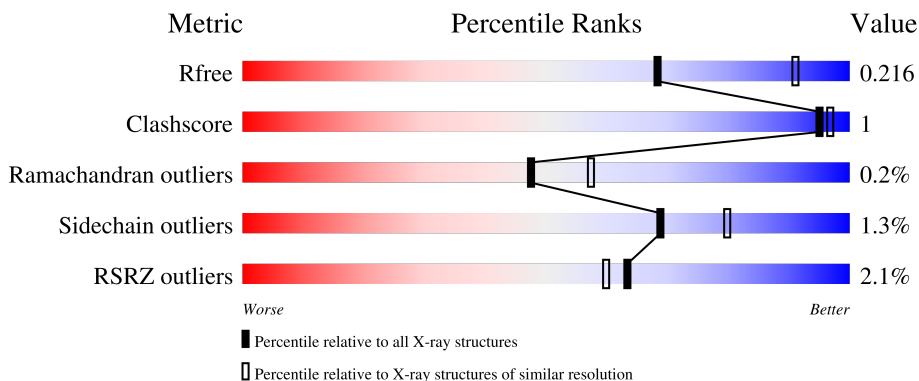
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.43 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	1564 (2.46-2.42)
Clashscore	141614	1631 (2.46-2.42)
Ramachandran outliers	138981	1617 (2.46-2.42)
Sidechain outliers	138945	1617 (2.46-2.42)
RSRZ outliers	127900	1547 (2.46-2.42)

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  $\geq 3$ , 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	543	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 94%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">2% 94% ..</p>
2	B	2	<div style="width: 100%; height: 10px; background-color: orange;"></div> <p style="text-align: center;">100%</p>
3	C	4	<div style="display: flex; align-items: center;"> <div style="width: 75%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 25%; height: 10px; background-color: yellow; margin-right: 5px;"></div> </div> <p style="text-align: center;">75% 25%</p>
4	D	5	<div style="display: flex; align-items: center;"> <div style="width: 20%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 80%; height: 10px; background-color: yellow; margin-right: 5px;"></div> </div> <p style="text-align: center;">20% 80%</p>

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	FUC	B	2	X	-	-	X
3	BMA	C	3	-	-	-	X
3	MAN	C	4	-	-	-	X
4	NAG	D	1	-	-	-	X
4	NAG	D	2	X	-	-	-
4	MAN	D	5	-	-	-	X
8	EDO	A	630	-	-	-	X

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 4707 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 Acetylcholinesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	532	4217	2713	709	772	23	0	2	0

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



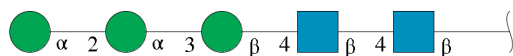
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	2	24	14	1	9	0	0	0

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



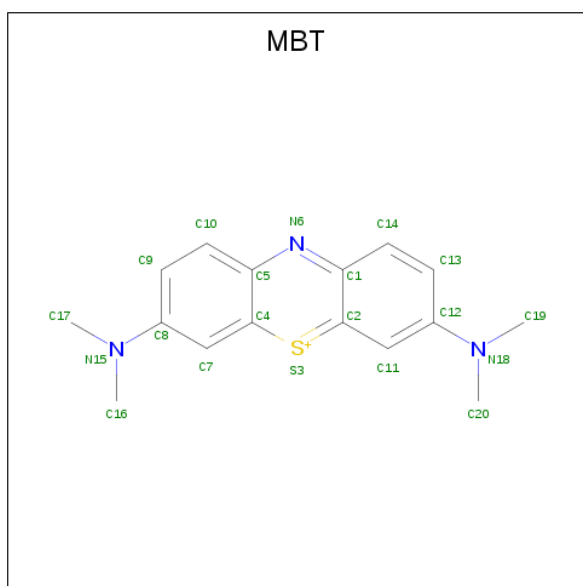
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	4	50	28	2	20	0	0	0

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-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
			Total	C	N	O			
4	D	5	61	34	2	25	0	0	0

- Molecule 5 is 3,7-BIS(DIMETHYLAMINO)PHENOTHIAZIN-5-IUM (three-letter code: MBT) (formula:  $C_{16}H_{18}N_3S$ ).



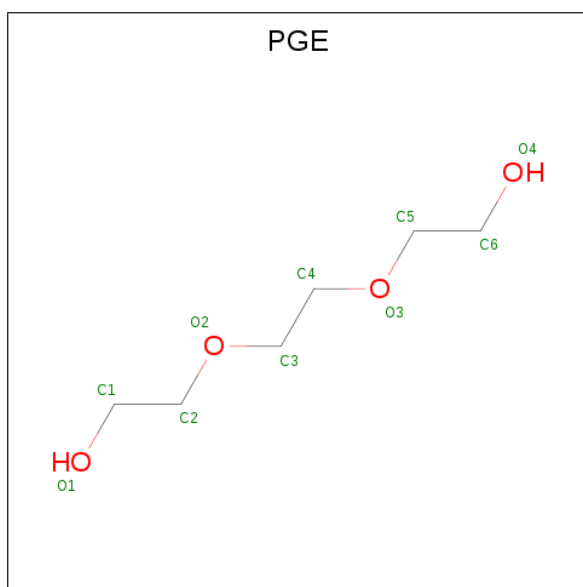
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	S		
5	A	1	20	16	3	1	0	0

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).



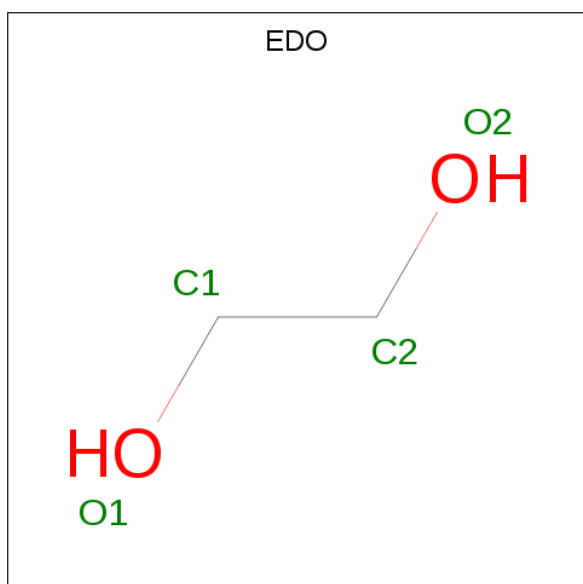
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0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

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 10 6 4	0	0
7	A	1	Total C O 10 6 4	0	0
7	A	1	Total C O 10 6 4	0	0
7	A	1	Total C O 10 6 4	0	0
7	A	1	Total C O 10 6 4	0	0

- Molecule 8 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0

- Molecule 9 is water.

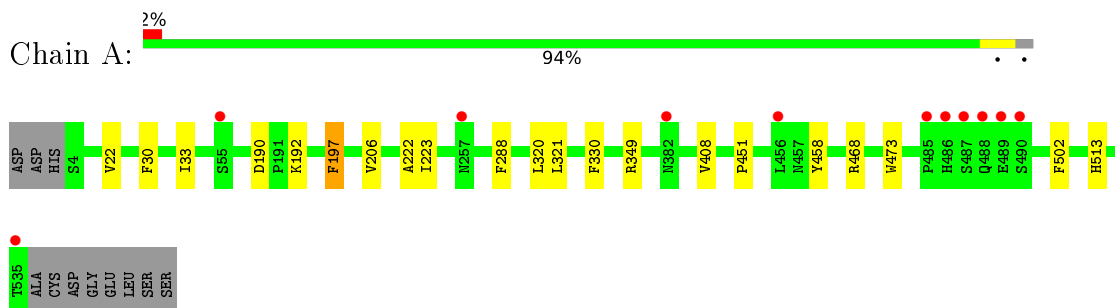
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	165	Total O 165 165	0	0



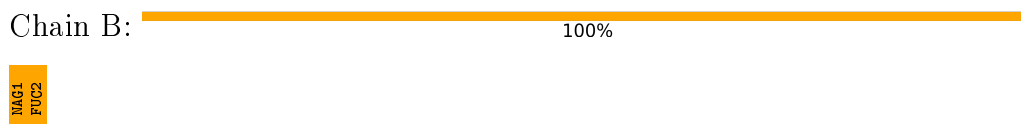
### 3 Residue-property plots [i](#)

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.

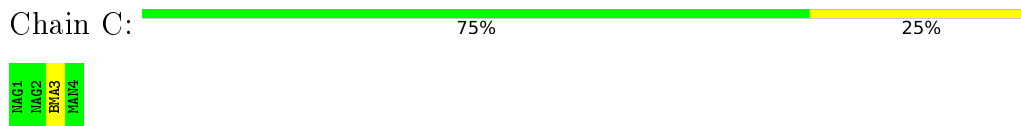
- Molecule 1: Acetylcholinesterase



- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



- Molecule 4: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-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 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.11Å 111.11Å 137.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	96.23 – 2.43 29.05 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.1 (96.23-2.43) 99.2 (29.05-2.43)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.90 (at 2.42Å)	Xtrriage
Refinement program	REFMAC 5.8.0073	Depositor
R, $R_{free}$	0.173 , 0.206 0.184 , 0.216	Depositor DCC
$R_{free}$ test set	1909 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.6	Xtrriage
Anisotropy	0.061	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 40.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4707	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MBT, FUC, PEG, EDO, PGE, BMA, MAN, 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	0.31	0/4345	0.51	0/5903

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	4217	0	4047	10	0
2	B	24	0	22	1	0
3	C	50	0	43	0	0
4	D	61	0	52	0	0
5	A	20	0	18	0	0
6	A	84	0	120	0	0
7	A	50	0	70	0	0
8	A	36	0	54	0	0
9	A	165	0	0	0	0
All	All	4707	0	4426	11	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 1.

All (11) 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:321:LEU:HD11	1:A:408:VAL:HG22	1.84	0.60
1:A:321:LEU:HD11	1:A:408:VAL:CG2	2.44	0.47
1:A:223:ILE:HA	1:A:320:LEU:O	2.15	0.46
1:A:502:PHE:CZ	1:A:513:HIS:HB2	2.52	0.45
1:A:197:PHE:CB	1:A:223:ILE:HB	2.48	0.44
1:A:451:PRO:HA	1:A:458:TYR:CD1	2.53	0.44
1:A:30:PHE:HB3	1:A:33:ILE:HD11	2.02	0.41
1:A:190:ASP:OD1	1:A:192:LYS:HG2	2.21	0.41
1:A:197:PHE:HB3	1:A:223:ILE:HB	2.03	0.41
2:B:1:NAG:H61	2:B:2:FUC:O2	2.21	0.40
1:A:206:VAL:CG1	1:A:222:ALA:HB1	2.51	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	532/543 (98%)	512 (96%)	19 (4%)	1 (0%)	47 57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	VAL

### 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	454/474 (96%)	448 (99%)	6 (1%)	69 80

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	197	PHE
1	A	288	PHE
1	A	330	PHE
1	A	349	ARG
1	A	468	ARG
1	A	473	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	374	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

11 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	B	1	2,1	14,14,15	0.46	0	17,19,21	1.19	1 (5%)
2	FUC	B	2	2	10,10,11	0.86	0	14,14,16	1.67	2 (14%)
3	NAG	C	1	1,3	14,14,15	0.55	0	17,19,21	0.77	0
3	NAG	C	2	3	14,14,15	0.52	0	17,19,21	0.82	0
3	BMA	C	3	3	11,11,12	0.40	0	15,15,17	0.84	1 (6%)
3	MAN	C	4	3	11,11,12	0.59	0	15,15,17	0.84	0
4	NAG	D	1	4,1	14,14,15	0.50	0	17,19,21	1.07	1 (5%)
4	NAG	D	2	4	14,14,15	0.61	0	17,19,21	1.86	2 (11%)
4	BMA	D	3	4	11,11,12	0.45	0	15,15,17	0.93	1 (6%)
4	MAN	D	4	4	11,11,12	0.56	0	15,15,17	0.92	0
4	MAN	D	5	4	11,11,12	0.57	0	15,15,17	1.45	2 (13%)

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

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	2	NAG	O5-C1-C2	6.00	120.75	111.29
2	B	2	FUC	O5-C1-C2	5.07	118.60	110.77
4	D	2	NAG	C1-O5-C5	4.12	117.77	112.19
4	D	5	MAN	C1-O5-C5	4.06	117.70	112.19
2	B	1	NAG	C1-O5-C5	3.34	116.72	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	5	MAN	C1-C2-C3	3.05	113.41	109.67
2	B	2	FUC	O2-C2-C1	2.57	114.40	109.15
4	D	1	NAG	C1-O5-C5	2.30	115.31	112.19
3	C	3	BMA	O5-C5-C6	2.22	110.68	107.20
4	D	3	BMA	C1-C2-C3	2.03	112.16	109.67

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	2	FUC	C1
4	D	2	NAG	C1

All (11) torsion outliers are listed below:

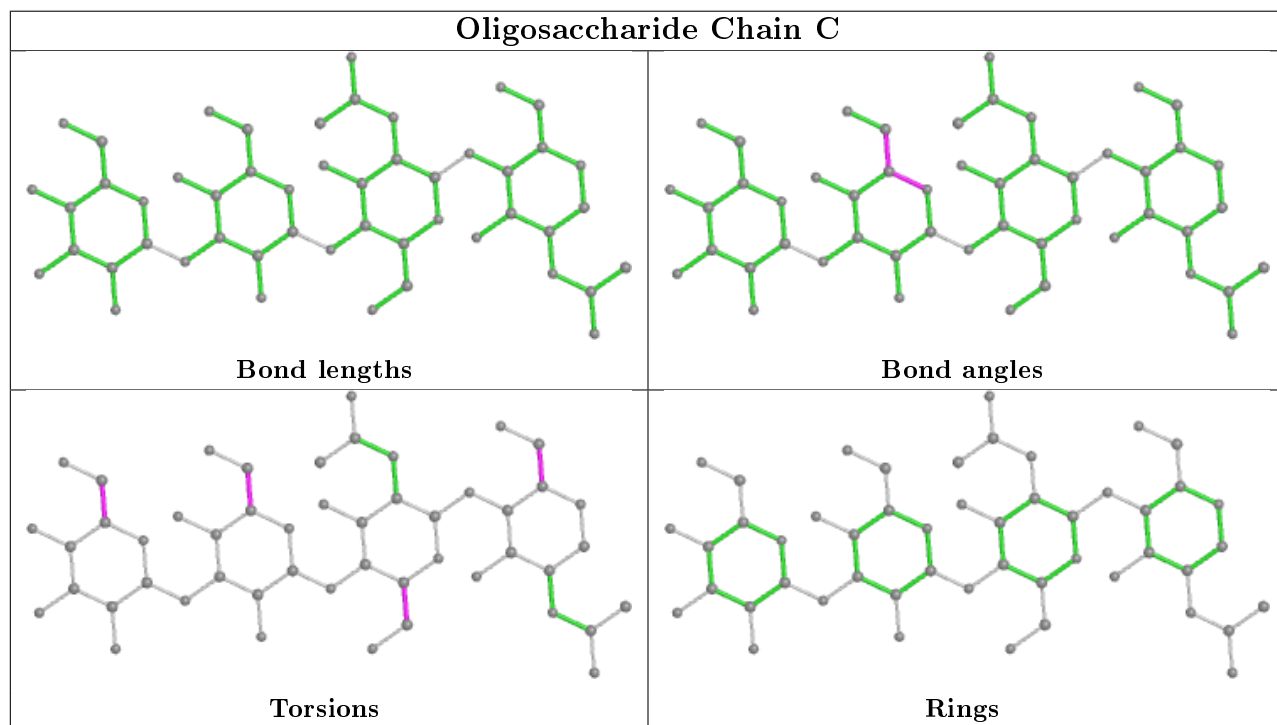
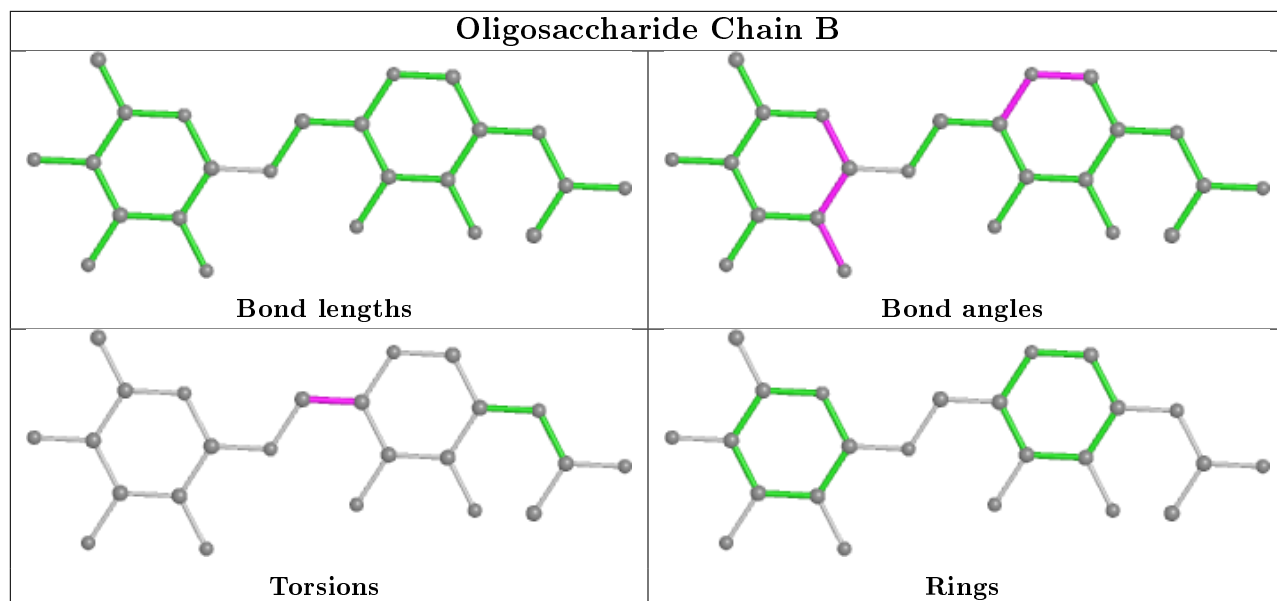
Mol	Chain	Res	Type	Atoms
2	B	1	NAG	O5-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6
3	C	2	NAG	C4-C5-C6-O6
4	D	2	NAG	C4-C5-C6-O6
4	D	2	NAG	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
3	C	3	BMA	C4-C5-C6-O6
3	C	3	BMA	O5-C5-C6-O6
3	C	1	NAG	C4-C5-C6-O6
3	C	4	MAN	C4-C5-C6-O6
3	C	1	NAG	O5-C5-C6-O6

There are no ring outliers.

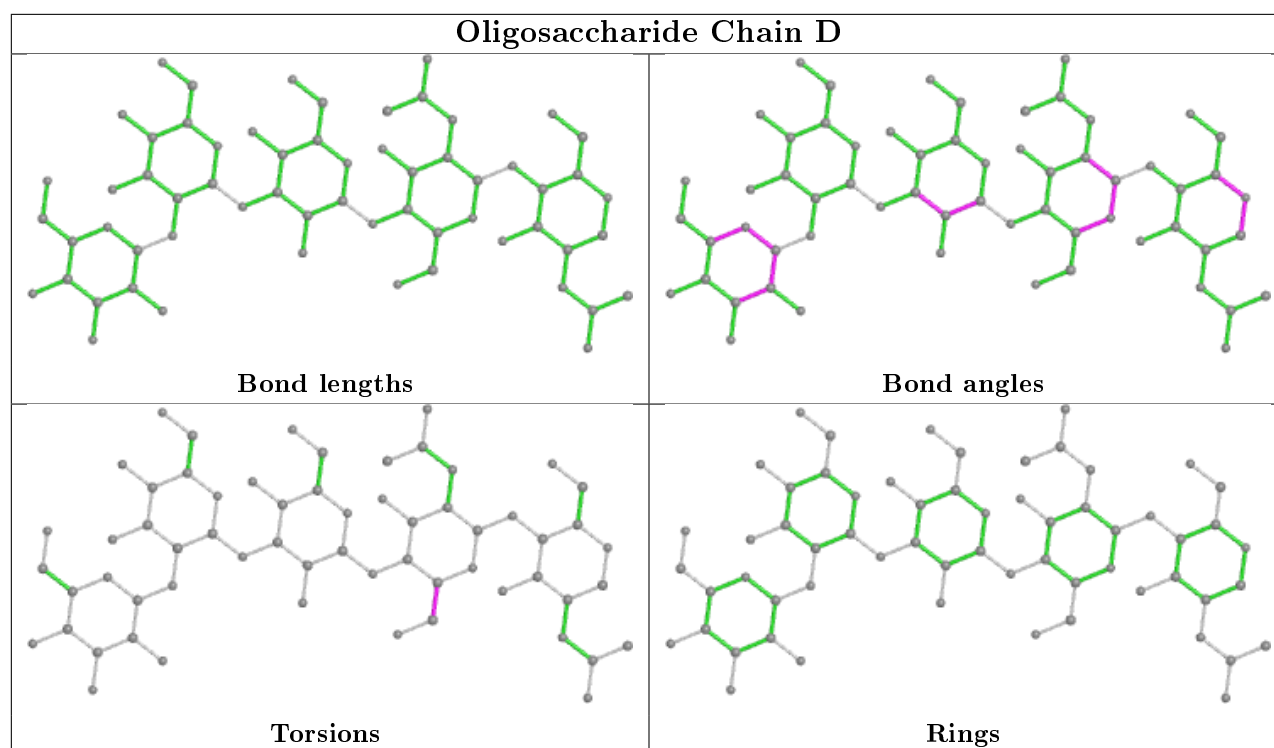
2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	NAG	1	0
2	B	2	FUC	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](#)

27 ligands 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
6	PEG	A	623	-	6,6,6	0.46	0	5,5,5	0.21	0
8	EDO	A	633	-	3,3,3	0.44	0	2,2,2	0.31	0
8	EDO	A	632	-	3,3,3	0.54	0	2,2,2	0.17	0
6	PEG	A	620	-	6,6,6	0.46	0	5,5,5	0.21	0
8	EDO	A	638	-	3,3,3	0.49	0	2,2,2	0.20	0
8	EDO	A	635	-	3,3,3	0.46	0	2,2,2	0.31	0
7	PGE	A	628	-	9,9,9	0.48	0	8,8,8	0.31	0
6	PEG	A	614	-	6,6,6	0.44	0	5,5,5	0.21	0
7	PGE	A	625	-	9,9,9	0.47	0	8,8,8	0.23	0
8	EDO	A	631	-	3,3,3	0.52	0	2,2,2	0.16	0
6	PEG	A	617	-	6,6,6	0.46	0	5,5,5	0.18	0
8	EDO	A	630	-	3,3,3	0.53	0	2,2,2	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	PEG	A	619	-	6,6,6	0.49	0	5,5,5	0.17	0
7	PGE	A	626	-	9,9,9	0.49	0	8,8,8	0.19	0
6	PEG	A	615	-	6,6,6	0.46	0	5,5,5	0.19	0
6	PEG	A	622	-	6,6,6	0.49	0	5,5,5	0.16	0
6	PEG	A	616	-	6,6,6	0.45	0	5,5,5	0.26	0
5	MBT	A	612	-	22,22,22	1.98	6 (27%)	32,32,32	2.34	8 (25%)
6	PEG	A	621	-	6,6,6	0.48	0	5,5,5	0.19	0
6	PEG	A	624	-	6,6,6	0.43	0	5,5,5	0.25	0
6	PEG	A	613	-	6,6,6	0.51	0	5,5,5	0.31	0
8	EDO	A	637	-	3,3,3	0.44	0	2,2,2	0.35	0
7	PGE	A	629	-	9,9,9	0.47	0	8,8,8	0.26	0
8	EDO	A	636	-	3,3,3	0.47	0	2,2,2	0.27	0
8	EDO	A	634	-	3,3,3	0.50	0	2,2,2	0.08	0
6	PEG	A	618	-	6,6,6	0.46	0	5,5,5	0.35	0
7	PGE	A	627	-	9,9,9	0.50	0	8,8,8	0.19	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	PEG	A	623	-	-	1/4/4/4	-
8	EDO	A	633	-	-	0/1/1/1	-
8	EDO	A	632	-	-	1/1/1/1	-
6	PEG	A	620	-	-	1/4/4/4	-
8	EDO	A	638	-	-	1/1/1/1	-
8	EDO	A	635	-	-	1/1/1/1	-
7	PGE	A	628	-	-	4/7/7/7	-
6	PEG	A	614	-	-	2/4/4/4	-
7	PGE	A	625	-	-	3/7/7/7	-
8	EDO	A	631	-	-	1/1/1/1	-
6	PEG	A	617	-	-	2/4/4/4	-
8	EDO	A	630	-	-	0/1/1/1	-
6	PEG	A	619	-	-	1/4/4/4	-
7	PGE	A	626	-	-	4/7/7/7	-
6	PEG	A	615	-	-	1/4/4/4	-
6	PEG	A	622	-	-	1/4/4/4	-
6	PEG	A	616	-	-	1/4/4/4	-
5	MBT	A	612	-	-	0/8/8/8	0/3/3/3
6	PEG	A	621	-	-	3/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PEG	A	624	-	-	1/4/4/4	-
6	PEG	A	613	-	-	4/4/4/4	-
8	EDO	A	637	-	-	0/1/1/1	-
7	PGE	A	629	-	-	4/7/7/7	-
8	EDO	A	636	-	-	0/1/1/1	-
8	EDO	A	634	-	-	1/1/1/1	-
6	PEG	A	618	-	-	2/4/4/4	-
7	PGE	A	627	-	-	2/7/7/7	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	612	MBT	C1-C2	3.78	1.49	1.43
5	A	612	MBT	C2-S3	-3.77	1.60	1.72
5	A	612	MBT	C4-S3	-3.67	1.60	1.72
5	A	612	MBT	C5-C4	3.31	1.48	1.43
5	A	612	MBT	C8-N15	2.81	1.44	1.37
5	A	612	MBT	C12-N18	2.51	1.43	1.37

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	612	MBT	C4-S3-C2	7.80	111.35	103.79
5	A	612	MBT	C2-C1-N6	-4.98	121.76	125.87
5	A	612	MBT	C5-N6-C1	4.83	124.68	117.86
5	A	612	MBT	C4-C5-N6	-4.25	122.36	125.87
5	A	612	MBT	C7-C4-S3	3.38	121.31	117.35
5	A	612	MBT	C11-C2-S3	2.34	120.09	117.35
5	A	612	MBT	C7-C4-C5	-2.12	118.98	121.74
5	A	612	MBT	C8-C7-C4	2.11	123.35	120.95

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	621	PEG	O1-C1-C2-O2
7	A	625	PGE	O2-C3-C4-O3
6	A	613	PEG	O1-C1-C2-O2
6	A	617	PEG	O2-C3-C4-O4
6	A	614	PEG	O2-C3-C4-O4

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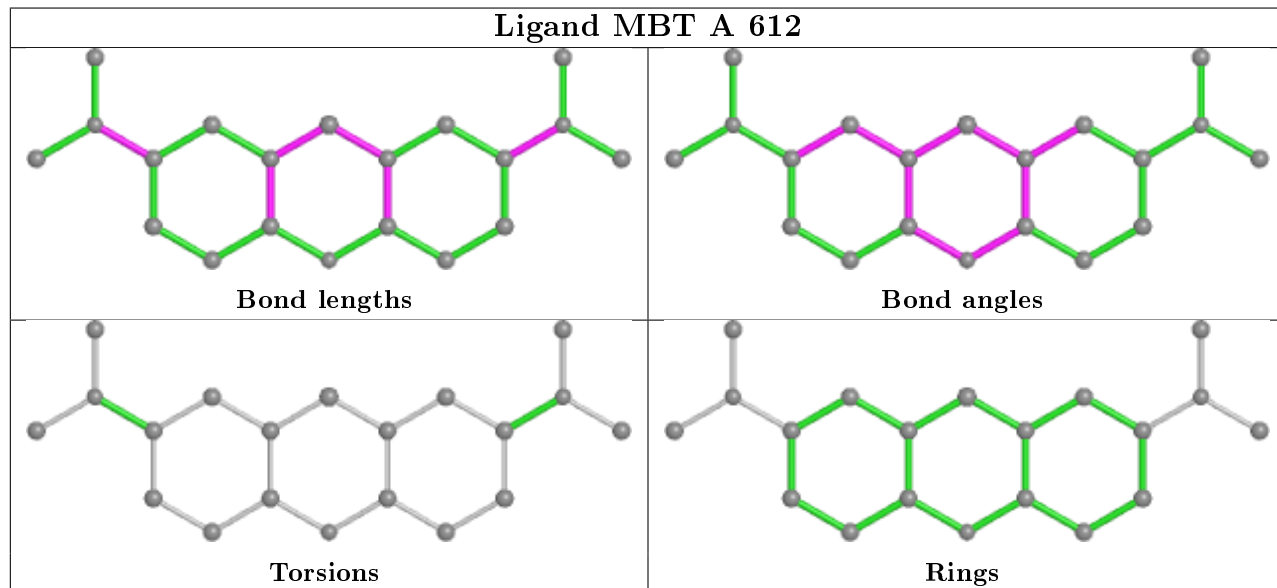
Mol	Chain	Res	Type	Atoms
6	A	615	PEG	O1-C1-C2-O2
7	A	626	PGE	O3-C5-C6-O4
7	A	628	PGE	O3-C5-C6-O4
6	A	613	PEG	O2-C3-C4-O4
6	A	617	PEG	O1-C1-C2-O2
6	A	620	PEG	O1-C1-C2-O2
7	A	629	PGE	O1-C1-C2-O2
8	A	635	EDO	O1-C1-C2-O2
8	A	638	EDO	O1-C1-C2-O2
6	A	621	PEG	O2-C3-C4-O4
7	A	627	PGE	O2-C3-C4-O3
6	A	614	PEG	O1-C1-C2-O2
6	A	624	PEG	O2-C3-C4-O4
8	A	631	EDO	O1-C1-C2-O2
8	A	632	EDO	O1-C1-C2-O2
7	A	628	PGE	O2-C3-C4-O3
7	A	626	PGE	O1-C1-C2-O2
7	A	625	PGE	C1-C2-O2-C3
7	A	628	PGE	C6-C5-O3-C4
6	A	619	PEG	C1-C2-O2-C3
6	A	618	PEG	C4-C3-O2-C2
7	A	628	PGE	C1-C2-O2-C3
7	A	629	PGE	O2-C3-C4-O3
7	A	625	PGE	C3-C4-O3-C5
6	A	613	PEG	C1-C2-O2-C3
6	A	621	PEG	C1-C2-O2-C3
7	A	629	PGE	C3-C4-O3-C5
6	A	613	PEG	C4-C3-O2-C2
7	A	627	PGE	C3-C4-O3-C5
6	A	618	PEG	C1-C2-O2-C3
7	A	626	PGE	C4-C3-O2-C2
7	A	626	PGE	O2-C3-C4-O3
6	A	623	PEG	C4-C3-O2-C2
8	A	634	EDO	O1-C1-C2-O2
6	A	616	PEG	C1-C2-O2-C3
6	A	622	PEG	C4-C3-O2-C2
7	A	629	PGE	C4-C3-O2-C2

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 all instances of the Ligand of Interest. In

addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	532/543 (97%)	-0.27	11 (2%) 63 60	33, 44, 62, 107	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	488	GLN	5.8
1	A	485	PRO	5.0
1	A	486	HIS	3.5
1	A	535	THR	3.3
1	A	489	GLU	2.9
1	A	257	ASN	2.8
1	A	487	SER	2.6
1	A	382	ASN	2.6
1	A	55	SER	2.5
1	A	490	SER	2.4
1	A	456	LEU	2.2

### 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, 95<sup>th</sup> 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.

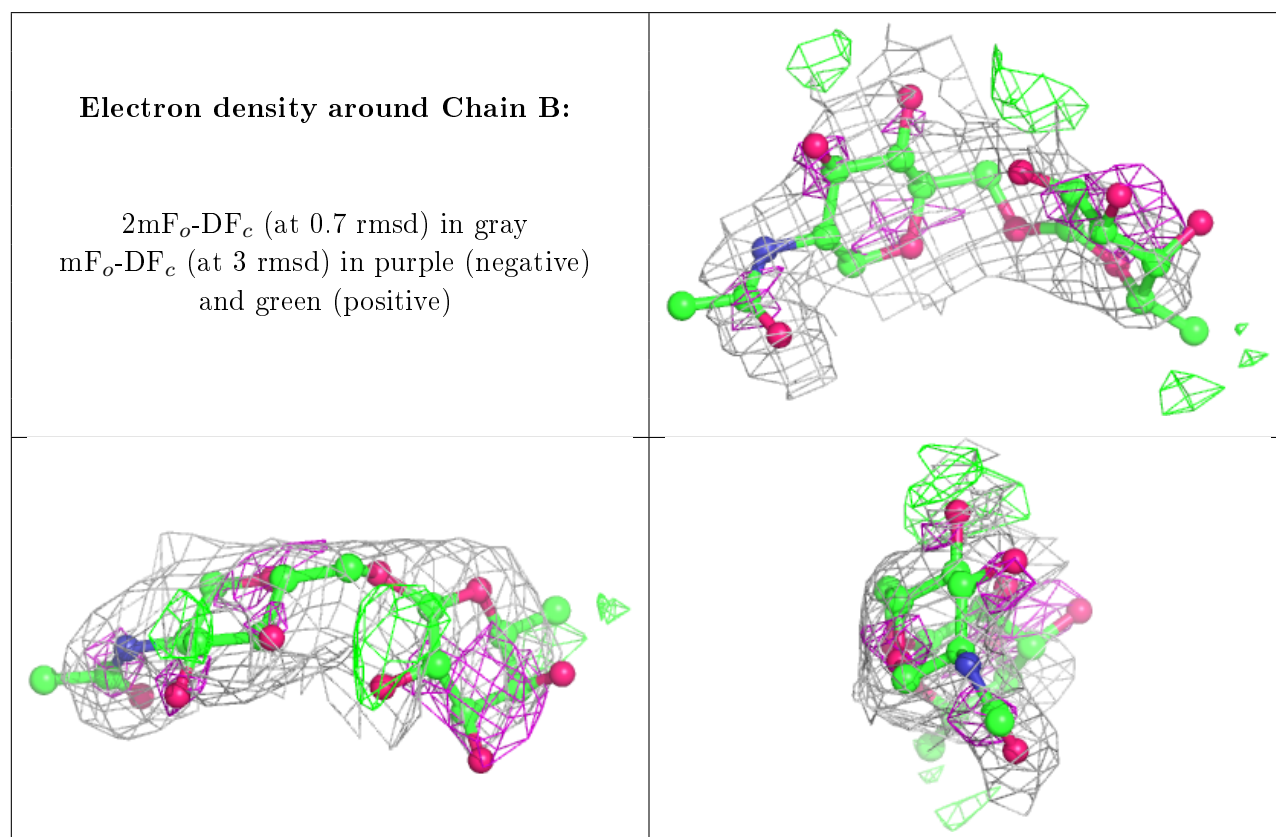
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	FUC	B	2	10/11	0.65	0.48	76,79,80,81	0

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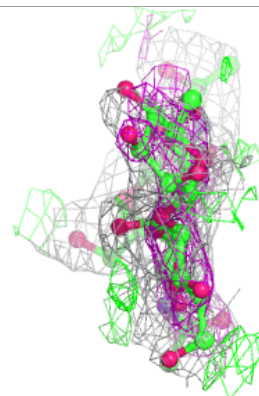
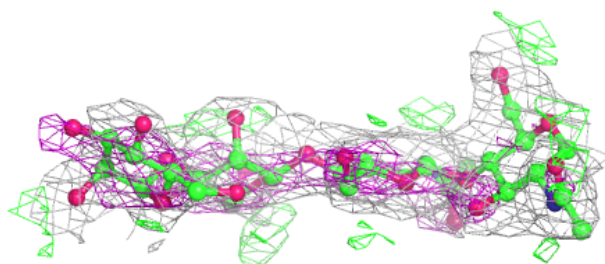
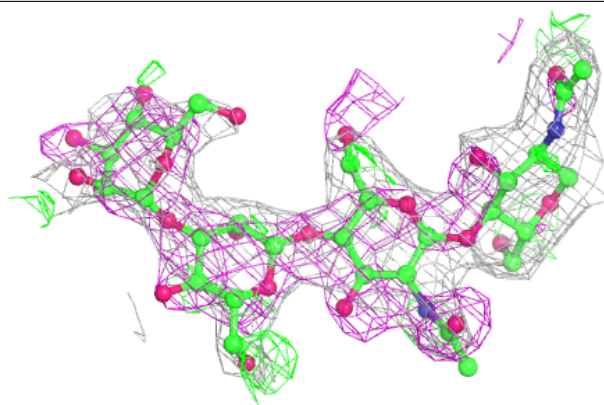
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	MAN	D	5	11/12	0.67	0.47	59,63,65,68	0
3	BMA	C	3	11/12	0.72	0.53	60,63,64,64	0
4	NAG	D	1	14/15	0.75	0.46	63,66,69,73	0
3	MAN	C	4	11/12	0.80	0.58	60,63,64,65	0
4	BMA	D	3	11/12	0.81	0.51	66,68,70,71	0
4	NAG	D	2	14/15	0.83	0.60	72,74,76,76	0
4	MAN	D	4	11/12	0.84	0.45	56,59,61,62	0
2	NAG	B	1	14/15	0.84	0.38	56,62,66,72	0
3	NAG	C	2	14/15	0.86	0.51	48,51,54,56	0
3	NAG	C	1	14/15	0.94	0.27	37,38,43,44	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.

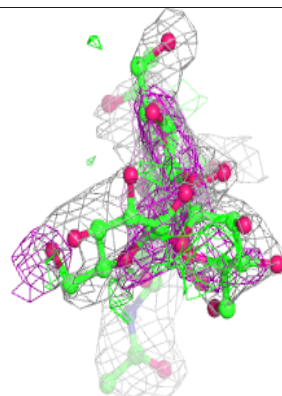
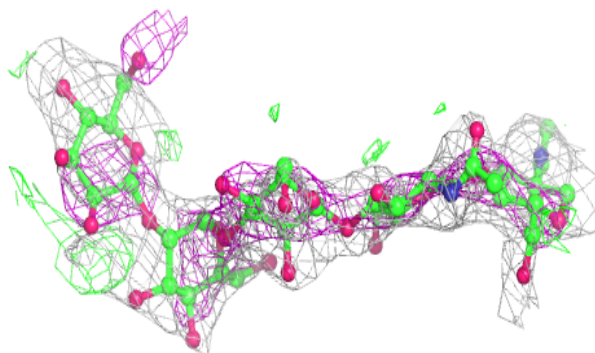
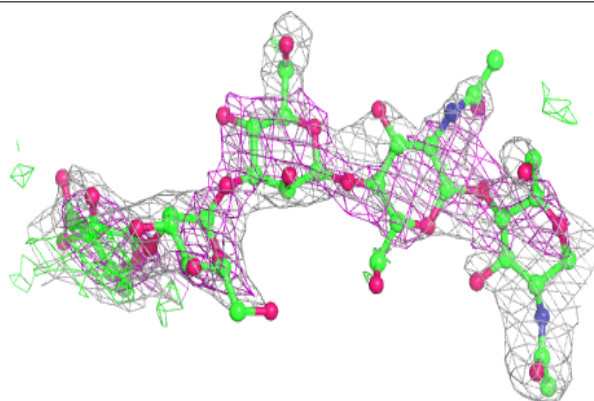


**Electron density around Chain C:**

$2mF_o-DF_c$  (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)



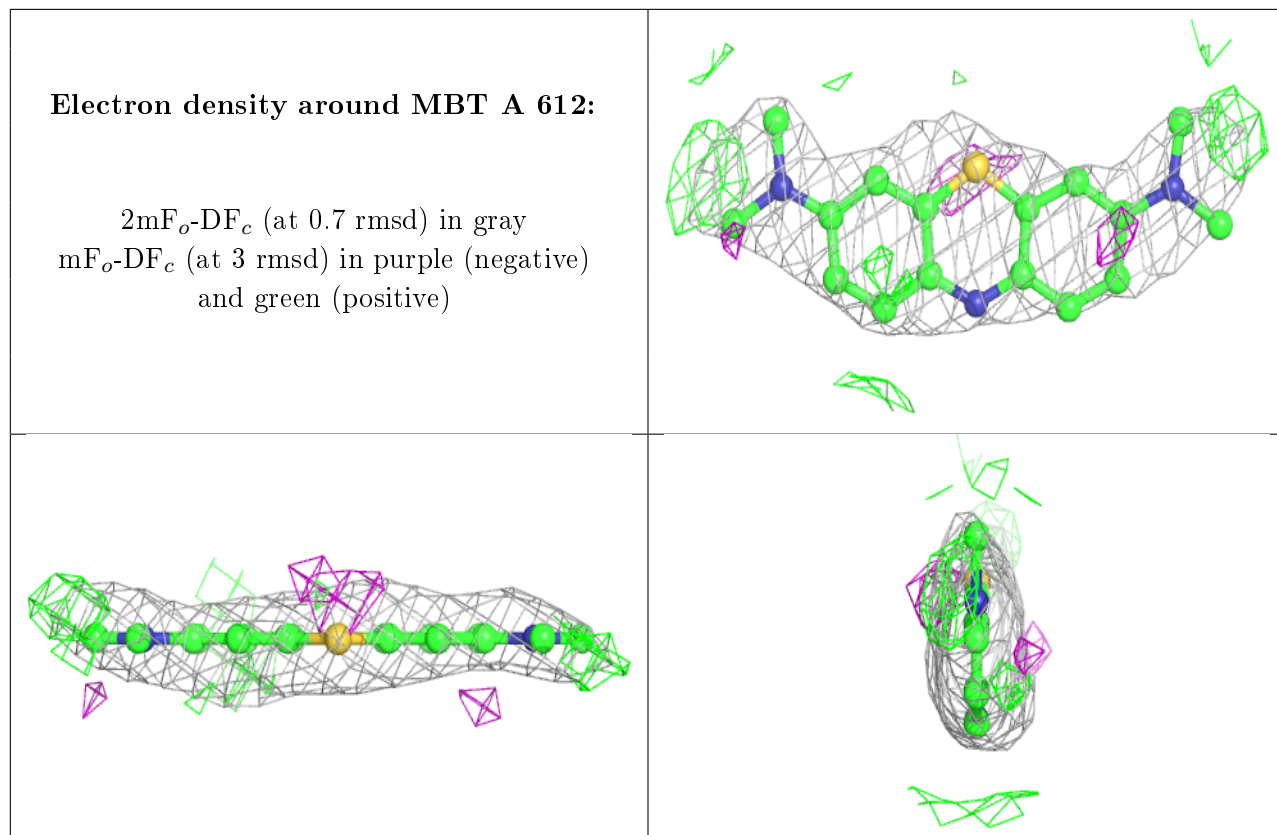


## 6.4 Ligands i

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
6	PEG	A	621	7/7	0.66	0.28	86,87,91,91	0
8	EDO	A	638	4/4	0.69	0.22	77,79,79,80	0
8	EDO	A	632	4/4	0.73	0.23	57,59,59,60	0
6	PEG	A	622	7/7	0.73	0.30	85,87,88,89	0
7	PGE	A	628	10/10	0.75	0.26	59,65,69,69	0
8	EDO	A	630	4/4	0.76	0.41	64,67,69,69	0
8	EDO	A	634	4/4	0.77	0.21	64,64,65,67	0
6	PEG	A	624	7/7	0.80	0.27	86,88,91,91	0
6	PEG	A	620	7/7	0.81	0.25	82,86,92,93	0
7	PGE	A	627	10/10	0.82	0.31	81,86,90,90	0
6	PEG	A	623	7/7	0.84	0.35	81,88,91,91	0
6	PEG	A	619	7/7	0.84	0.28	60,66,77,78	0
7	PGE	A	626	10/10	0.84	0.35	78,79,84,85	0
6	PEG	A	617	7/7	0.85	0.23	66,68,69,70	0
7	PGE	A	625	10/10	0.85	0.30	65,73,74,76	0
8	EDO	A	635	4/4	0.85	0.29	81,81,82,83	0
8	EDO	A	631	4/4	0.85	0.21	57,62,64,64	0
6	PEG	A	616	7/7	0.86	0.21	68,70,71,72	0
6	PEG	A	614	7/7	0.86	0.26	78,79,82,82	0
8	EDO	A	636	4/4	0.88	0.27	76,78,78,78	0
6	PEG	A	618	7/7	0.89	0.22	65,67,76,76	0
5	MBT	A	612	20/20	0.89	0.21	70,74,76,84	0
8	EDO	A	637	4/4	0.91	0.15	73,74,74,74	0
7	PGE	A	629	10/10	0.92	0.17	61,72,77,78	0
6	PEG	A	615	7/7	0.92	0.30	66,66,69,71	0
6	PEG	A	613	7/7	0.92	0.15	64,64,67,67	0
8	EDO	A	633	4/4	0.94	0.09	69,69,70,71	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



## 6.5 Other polymers [i](#)

There are no such residues in this entry.