

wwPDB X-ray Structure Validation Summary Report (i)

Feb 17, 2024 – 11:33 AM EST

PDB ID : 3SET

Title: Ni-mediated Dimer of Maltose-binding Protein A216H/K220H by Synthetic

Symmetrization (Form I)

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Deposited on : 2011-06-11

Resolution : 1.90 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 (1)) were used in the production of this report:

 $Mol Probity \quad : \quad 4.02b\text{--}467$

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

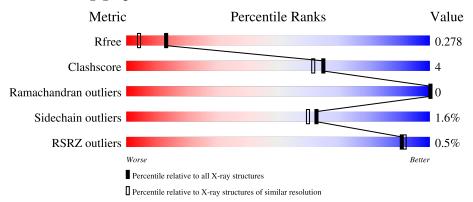
Validation Pipeline (wwPDB-VP) : 2.36

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 1.90 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	Similar resolution $(\# \text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	372	87%	12% •
1	С	372	88%	11% •
2	В	2	100%	
2	D	2	50% 50%	



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5994 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 Maltose-binding periplasmic protein.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	370	Total 2842	C 1829	11	O 544	S 6	0	0	0
1	С	369	Total 2818		N 461	O 537	S 6	0	1	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P0AEX9
A	216	HIS	ALA	engineered mutation	UNP POAEX9
A	220	HIS	LYS	engineered mutation	UNP POAEX9
A	360	ALA	-	expression tag	UNP P0AEX9
A	361	ALA	-	expression tag	UNP P0AEX9
A	362	LEU	-	expression tag	UNP P0AEX9
A	363	ALA	-	expression tag	UNP P0AEX9
A	364	ALA	-	expression tag	UNP P0AEX9
A	365	ALA	-	expression tag	UNP P0AEX9
A	366	GLN	-	expression tag	UNP P0AEX9
A	367	THR	-	expression tag	UNP P0AEX9
A	368	ASN	-	expression tag	UNP P0AEX9
A	369	ALA	-	expression tag	UNP P0AEX9
A	370	ALA	-	expression tag	UNP P0AEX9
A	371	ALA	-	expression tag	UNP P0AEX9
A	372	ALA	-	expression tag	UNP P0AEX9
С	1	MET	-	initiating methionine	UNP P0AEX9
С	216	HIS	ALA	engineered mutation	UNP P0AEX9
С	220	HIS	LYS	engineered mutation	UNP P0AEX9
С	360	ALA	-	expression tag	UNP P0AEX9
С	361	ALA	-	expression tag	UNP P0AEX9
С	362	LEU	-	expression tag	UNP P0AEX9
С	363	ALA	-	expression tag	UNP P0AEX9
С	364	ALA		expression tag	UNP P0AEX9
С	365	ALA	_	expression tag	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
С	366	GLN	-	expression tag	UNP P0AEX9
С	367	THR	-	expression tag	UNP P0AEX9
С	368	ASN	-	expression tag	UNP P0AEX9
С	369	ALA	-	expression tag	UNP P0AEX9
С	370	ALA	-	expression tag	UNP P0AEX9
С	371	ALA	-	expression tag	UNP P0AEX9
С	372	ALA	=	expression tag	UNP P0AEX9

• Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace
2	В	2	Total C 23 12		0	2	0
2	D	2	Total C 23 12	O 11	0	2	0

• Molecule 3 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Ni 1 1	0	0
3	С	1	Total Ni 1 1	0	0

• Molecule 4 is water.

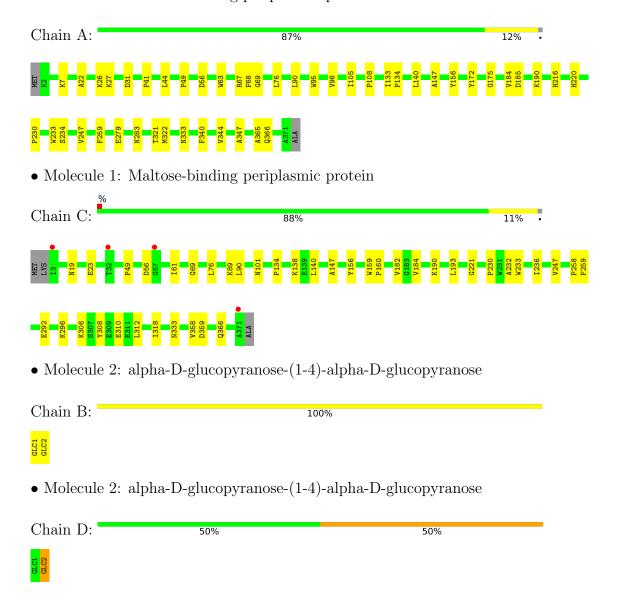
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	150	Total O 150 150	0	0
4	С	136	Total O 136 136	0	0



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: Maltose-binding periplasmic protein





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	72.94Å 62.27Å 89.71Å	Depositor
a, b, c, α , β , γ	90.00° 109.03° 90.00°	Depositor
Resolution (Å)	42.90 - 1.90	Depositor
rtesolution (A)	50.19 - 1.90	EDS
% Data completeness	99.7 (42.90-1.90)	Depositor
(in resolution range)	99.7 (50.19-1.90)	EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.70 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
D D.	0.232 , 0.282	Depositor
R, R_{free}	0.228 , 0.278	DCC
R_{free} test set	3000 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	14.7	Xtriage
Anisotropy	0.612	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 41.5	EDS
L-test for twinning ²	$ < L >=0.44, < L^2>=0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5994	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.88% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: GLC, NI

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		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.27	0/2913	0.43	0/3964	
1	С	0.25	0/2892	0.41	0/3941	
All	All	0.26	0/5805	0.42	0/7905	

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	2842	0	2775	23	0
1	С	2818	0	2736	22	0
2	В	23	0	21	2	0
2	D	23	0	21	1	0
3	A	1	0	0	0	0
3	С	1	0	0	0	0
4	A	150	0	0	2	0
4	С	136	0	0	2	0
All	All	5994	0	5553	46	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 4.



The worst 5 of 46 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} & ext{Interatomic} \ & ext{distance} \ & ext{(Å)} \end{aligned}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:C:69:GLY:HA3	1:C:333:ASN:O	1.89	0.72
1:A:69:GLY:HA3	1:A:333:ASN:O	1.91	0.71
1:C:184:VAL:HG22	1:C:366:GLN:HE21	1.67	0.57
1:A:98:VAL:HG21	1:A:108:PRO:HD3	1.89	0.56
1:C:306:LYS:O	1:C:310:GLU:HG3	2.06	0.55

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	$368/372 \ (99\%)$	361 (98%)	7 (2%)	0	100	100
1	C	$368/372 \ (99\%)$	363 (99%)	5 (1%)	0	100	100
All	All	736/744 (99%)	724 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

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.

\mathbf{Mol}	Chain	Analysed	lysed Rotameric		Percentiles	
1	A	287/293 (98%)	282 (98%)	5 (2%)	60 57	

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
1	С	282/293 (96%)	278 (99%)	4 (1%)	67 65		
All	All	569/586 (97%)	560 (98%)	9 (2%)	62 60		

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	101	ASN
1	С	259	PHE
1	A	247	VAL
1	A	259	PHE
1	С	56	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	368	ASN
1	С	50	GLN
1	С	368	ASN
1	С	219	ASN
1	С	366	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)

4 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	Mol Type	Chain	Peg	Res Link	Bond lengths			Bond angles		
MIOI		Chain	rtes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GLC	В	1[A]	2	12,12,12	0.59	0	17,17,17	0.58	0
2	GLC	В	2[A]	2	11,11,12	0.63	0	15,15,17	0.68	0
2	GLC	D	1[A]	2	12,12,12	0.51	0	17,17,17	0.64	0
2	GLC	D	2[A]	2	11,11,12	0.56	0	15,15,17	0.98	1 (6%)

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	GLC	В	1[A]	2	-	0/2/22/22	0/1/1/1
2	GLC	В	2[A]	2	-	0/2/19/22	0/1/1/1
2	GLC	D	1[A]	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2[A]	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

\mathbf{Mol}	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	D	2[A]	GLC	O5-C5-C6	2.33	110.86	107.20

There are no chirality outliers.

There are no torsion outliers.

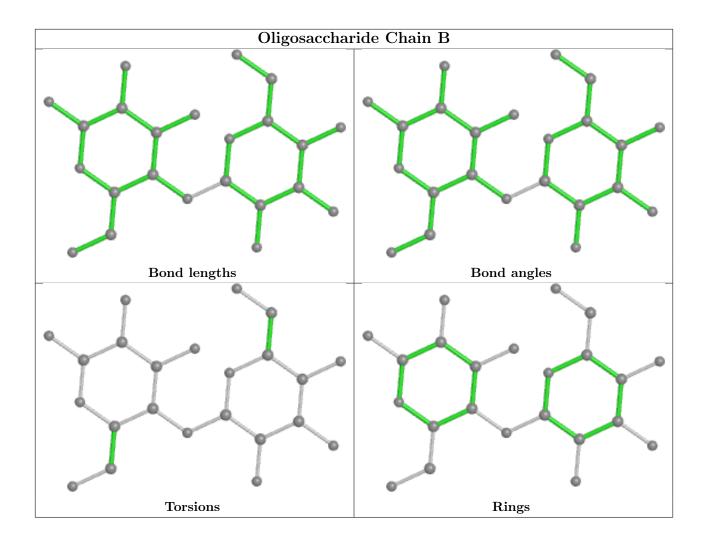
There are no ring outliers.

3 monomers are involved in 3 short contacts:

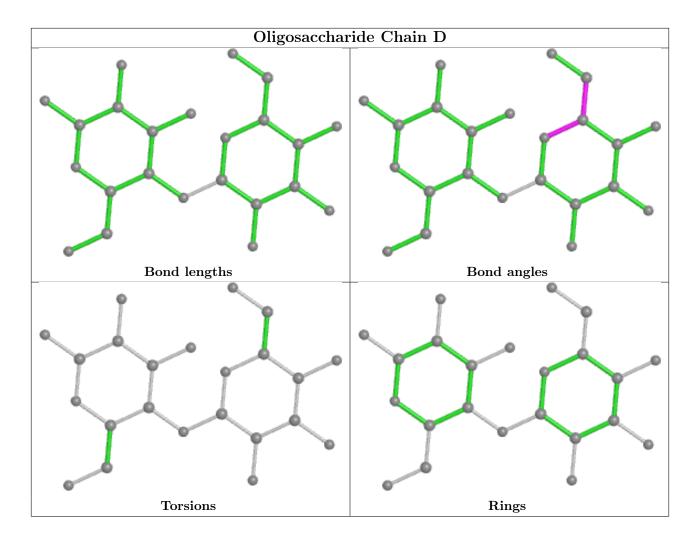
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	1[A]	GLC	1	0
2	D	2[A]	GLC	1	0
2	В	2[A]	GLC	1	0

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)

Of 2 ligands modelled in this entry, 2 are 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 (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^{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	$OWAB(A^2)$	Q < 0.9
1	A	370/372 (99%)	-0.00	0 100 100	5, 15, 29, 47	0
1	С	369/372~(99%)	0.01	4 (1%) 80 82	6, 16, 36, 48	0
All	All	739/744 (99%)	0.00	4 (0%) 91 92	5, 16, 32, 48	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ	
1	С	371	ALA	4.2	
1	С	57	GLY	2.7	
1	С	3	ILE	2.3	
1	С	32	THR	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)

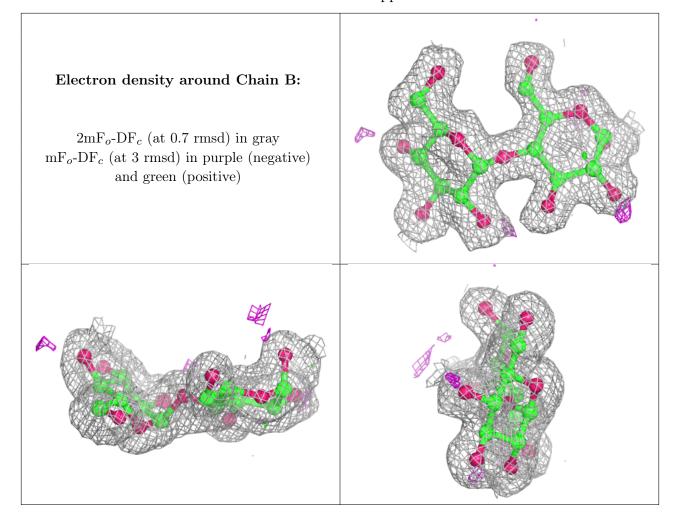
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
2	GLC	В	1[A]	12/12	0.92	0.11	4,8,14,16	0
2	GLC	D	1[A]	12/12	0.94	0.09	5,12,15,18	0
2	GLC	D	2[A]	11/12	0.95	0.08	6,8,11,11	0
2	GLC	В	2[A]	11/12	0.96	0.10	5,8,11,12	0

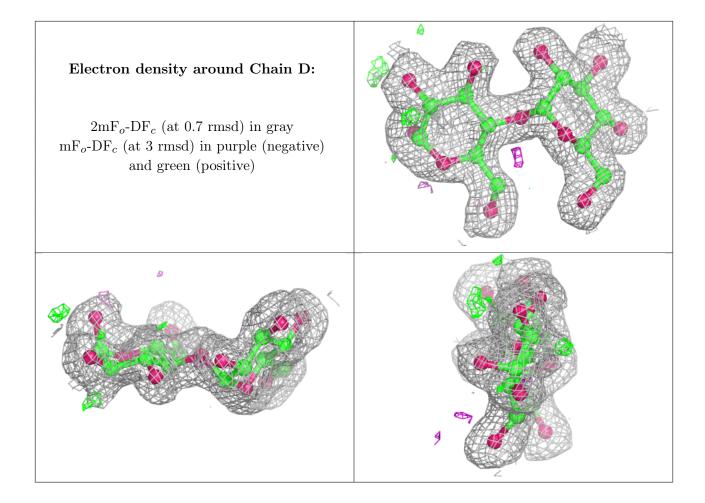
The following is a graphical depiction of the model fit to experimental electron density for oligosac-



charide. Each fit is shown from different orientation to approximate a three-dimensional view.







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^{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.

Mo	l Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	NI	A	373	1/1	0.99	0.08	16,16,16,16	0
3	NI	С	373	1/1	1.00	0.09	12,12,12,12	0

6.5 Other polymers (i)

There are no such residues in this entry.

