



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 06:21 AM UTC

PDB ID : 3QWQ / pdb_00003qwq
Title : Crystal structure of the extracellular domain of the epidermal growth factor receptor in complex with an adnectin
Authors : Sheriff, S.
Deposited on : 2011-02-28
Resolution : 2.75 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

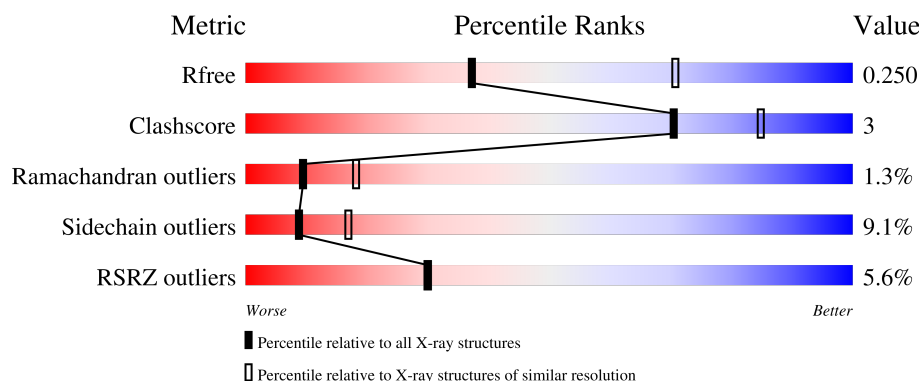
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	648	5% 77% 14% • 5%
2	B	114	5% 73% 12% • 12%
3	C	6	33% 67%
3	F	6	100%
4	D	3	100%

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Mol	Chain	Length	Quality of chain
5	E	2	 100%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5610 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 Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	613	Total	C	N	O	S	0	0	0
			4612	2844	815	893	60			

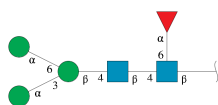
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	619	HIS	-	expression tag	UNP P00533
A	620	HIS	-	expression tag	UNP P00533
A	621	HIS	-	expression tag	UNP P00533
A	622	HIS	-	expression tag	UNP P00533
A	623	HIS	-	expression tag	UNP P00533
A	624	HIS	-	expression tag	UNP P00533

- Molecule 2 is a protein called ADNECTIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	0	0	0
			747	475	124	148			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[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
3	C	6	Total	C	N	O	0	0	0
			71	40	2	29			
3	F	6	Total	C	N	O	0	0	0
			71	40	2	29			

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



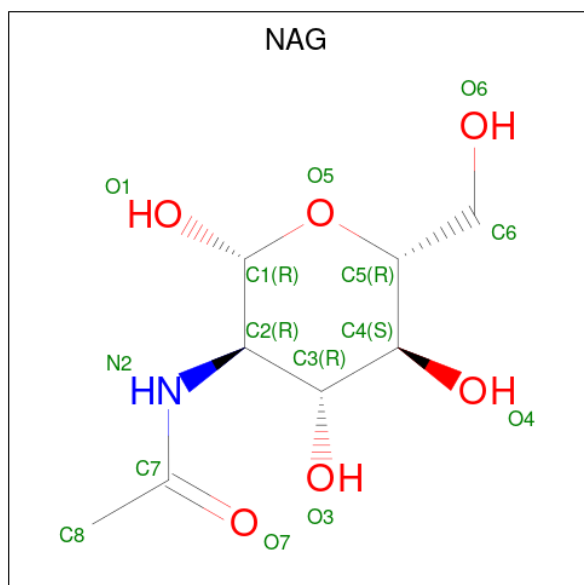
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	12	Total	O	0	0
			12	12		
7	B	2	Total	O	0	0
			2	2		

- Molecule 3: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-[α -L-fucopyranose-(1-6)]2-acetamido-2-deoxy- β -D-glucopyranose

Chain F:  100%

MAG1
MAG2
BMA3
MAN4
MAN5
FUC6

- Molecule 4: β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain D:  100%

MAG1
MAG2
BMA3

- Molecule 5: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain E:  100%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.00Å 72.10Å 262.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.47 – 2.75 49.47 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.47-2.75) 99.5 (49.47-2.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.04 (at 2.77Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.13.0, BUSTER 2.13.0	Depositor
R, R_{free}	0.202 , 0.246 0.210 , 0.250	Depositor DCC
R_{free} test set	1058 reflections (3.09%)	wwPDB-VP
Wilson B-factor (Å ²)	58.6	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 65.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5610	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, FUC, MAN, 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	0.90	3/4702 (0.1%)	1.39	35/6383 (0.5%)
2	B	0.89	0/770	1.32	4/1062 (0.4%)
All	All	0.90	3/5472 (0.1%)	1.38	39/7445 (0.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	87	MET	SD-CE	9.86	2.04	1.79
1	A	152	MET	SD-CE	-5.70	1.65	1.79
1	A	113	MET	SD-CE	-5.48	1.65	1.79

The worst 5 of 39 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	520	GLY	N-CA-C	8.80	124.04	112.25
1	A	159	HIS	CA-C-N	8.12	137.05	121.54
1	A	159	HIS	C-N-CA	8.12	137.05	121.54
2	B	23	ASP	CA-CB-CG	7.14	119.74	112.60
1	A	138	ILE	N-CA-C	7.13	118.64	109.30

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	4612	0	4321	34	0
2	B	747	0	684	3	0
3	C	71	0	61	0	0
3	F	71	0	61	0	0
4	D	39	0	34	0	0
5	E	28	0	25	0	0
6	A	28	0	26	0	0
7	A	12	0	0	0	0
7	B	2	0	0	0	0
All	All	5610	0	5212	37	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.

The worst 5 of 37 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:87:MET:CE	1:A:87:MET:SD	2.04	1.46
1:A:521:GLU:HB2	1:A:522:PRO:HD3	1.49	0.94
1:A:327:ILE:HD11	1:A:345:LEU:HG	1.54	0.89
1:A:146:SER:HA	1:A:149:LEU:HD23	1.56	0.86
1:A:144:VAL:HG11	1:A:152:MET:HE3	1.75	0.69

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	611/648 (94%)	559 (92%)	46 (8%)	6 (1%)	12	24
2	B	98/114 (86%)	91 (93%)	4 (4%)	3 (3%)	3	6
All	All	709/762 (93%)	650 (92%)	50 (7%)	9 (1%)	9	18

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	306	GLU
1	A	336	LYS
2	B	4	VAL
1	A	159	HIS
2	B	40	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	508/559 (91%)	462 (91%)	46 (9%)	9	17
2	B	77/99 (78%)	70 (91%)	7 (9%)	9	17
All	All	585/658 (89%)	532 (91%)	53 (9%)	9	17

5 of 53 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	400	GLU
1	A	449	ASN
2	B	49	THR
1	A	406	THR
1	A	439	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	384	GLN
1	A	580	ASN
1	A	394	HIS
2	B	42	ASN
1	A	528	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 ⓘ

17 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	C	1	1,3	14,14,15	0.27	0	17,19,21	0.84	0
3	NAG	C	2	3	14,14,15	0.38	0	17,19,21	0.77	0
3	BMA	C	3	3	11,11,12	0.32	0	15,15,17	1.95	1 (6%)
3	MAN	C	4	3	11,11,12	0.43	0	15,15,17	0.85	1 (6%)
3	MAN	C	5	3	11,11,12	0.33	0	15,15,17	1.09	1 (6%)
3	FUC	C	6	3	10,10,11	0.47	0	14,14,16	0.87	1 (7%)
4	NAG	D	1	1,4	14,14,15	0.29	0	17,19,21	0.89	1 (5%)
4	NAG	D	2	4	14,14,15	0.24	0	17,19,21	2.47	3 (17%)
4	BMA	D	3	4	11,11,12	0.47	0	15,15,17	1.10	1 (6%)
5	NAG	E	1	1,5	14,14,15	0.34	0	17,19,21	1.82	2 (11%)
5	NAG	E	2	5	14,14,15	0.34	0	17,19,21	1.57	2 (11%)
3	NAG	F	1	1,3	14,14,15	0.28	0	17,19,21	1.21	1 (5%)
3	NAG	F	2	3	14,14,15	0.26	0	17,19,21	0.77	1 (5%)
3	BMA	F	3	3	11,11,12	0.45	0	15,15,17	2.08	2 (13%)
3	MAN	F	4	3	11,11,12	0.50	0	15,15,17	1.52	1 (6%)
3	MAN	F	5	3	11,11,12	0.42	0	15,15,17	1.37	1 (6%)
3	FUC	F	6	3	10,10,11	0.43	0	14,14,16	0.73	1 (7%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

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

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

There are no bond length outliers.

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	2	NAG	C1-C2-N2	7.91	122.90	110.43
3	C	3	BMA	C1-O5-C5	7.01	121.58	112.19
3	F	3	BMA	C1-O5-C5	6.95	121.51	112.19
5	E	1	NAG	C1-O5-C5	5.64	119.74	112.19
3	F	4	MAN	C1-O5-C5	5.30	119.28	112.19

There are no chirality outliers.

5 of 15 torsion outliers are listed below:

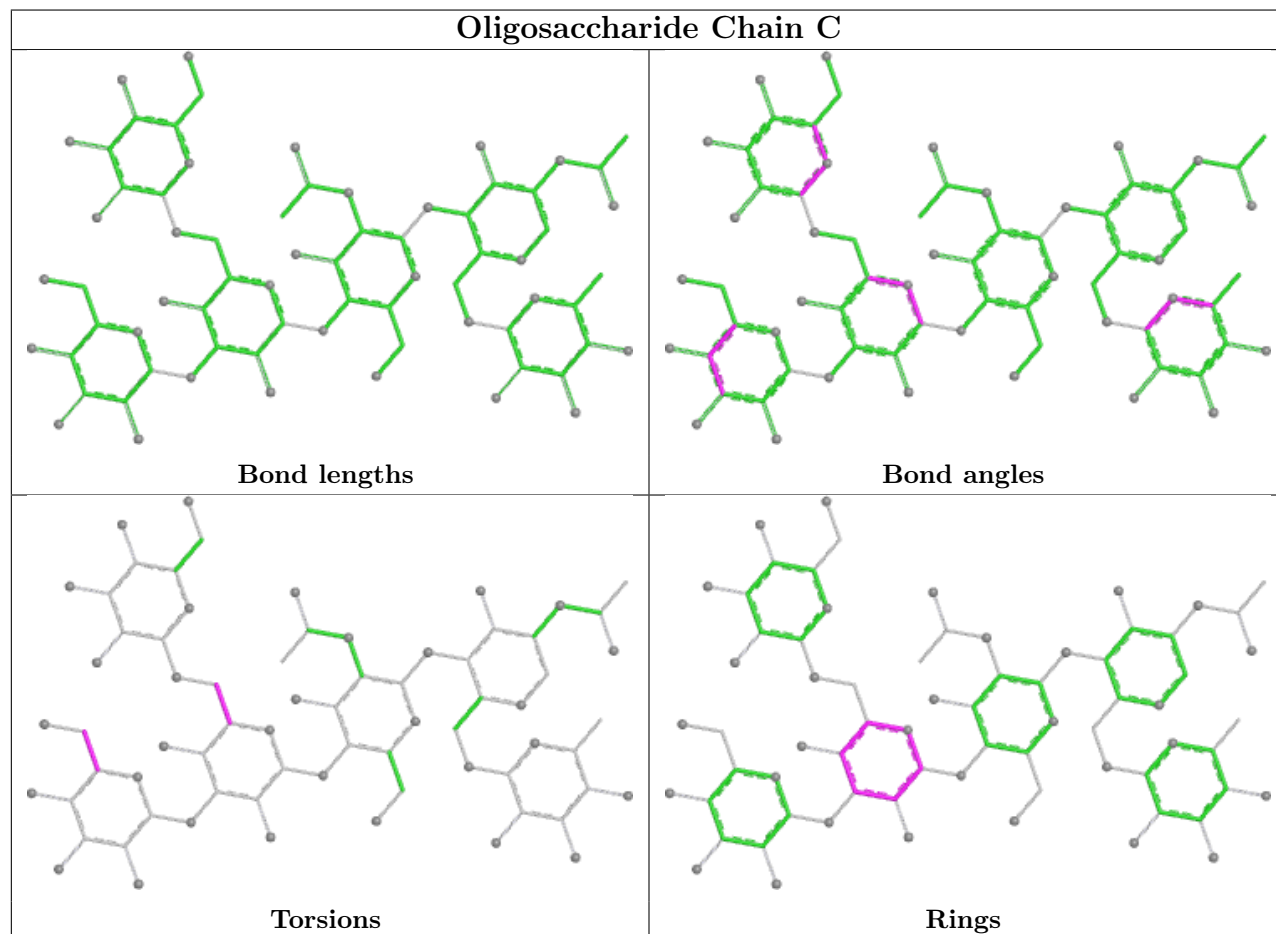
Mol	Chain	Res	Type	Atoms
4	D	2	NAG	C1-C2-N2-C7
3	F	3	BMA	O5-C5-C6-O6
3	C	3	BMA	O5-C5-C6-O6
3	C	3	BMA	C4-C5-C6-O6
3	F	4	MAN	C4-C5-C6-O6

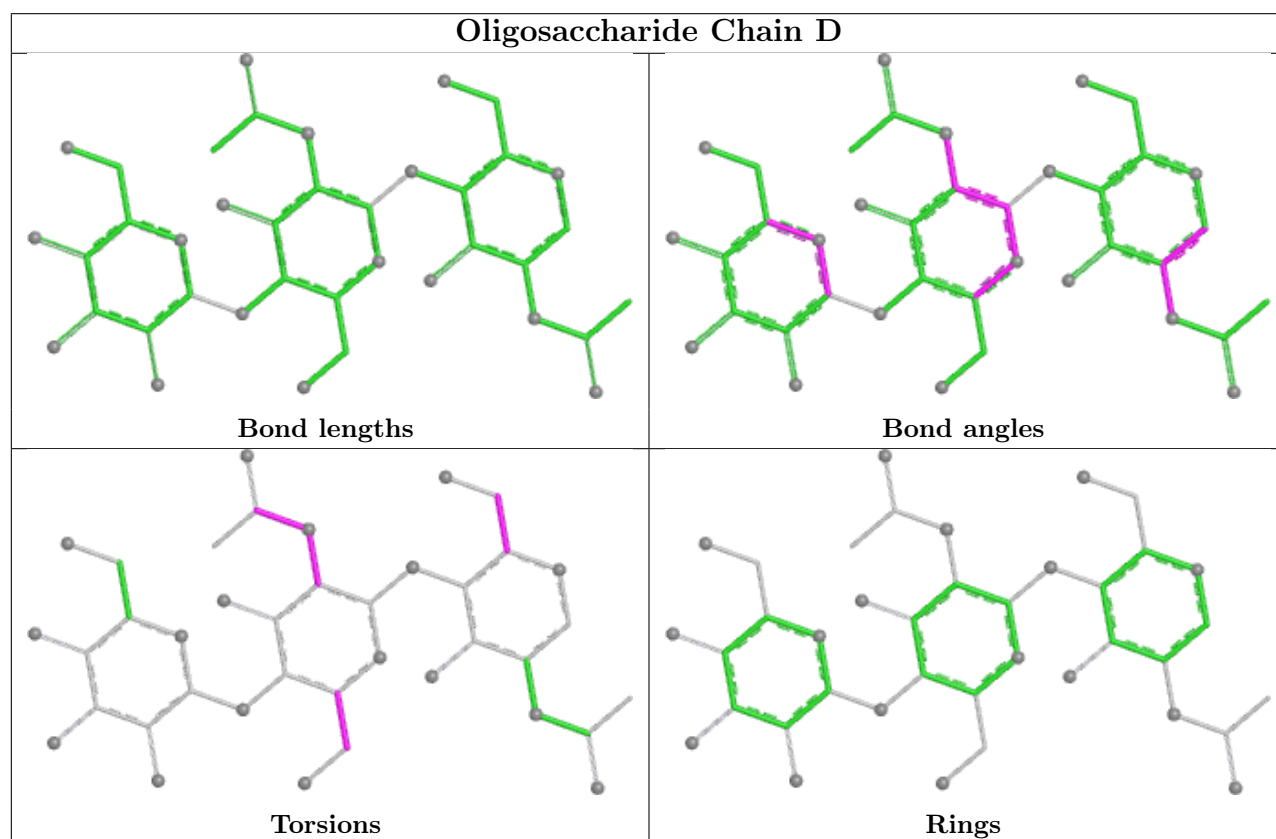
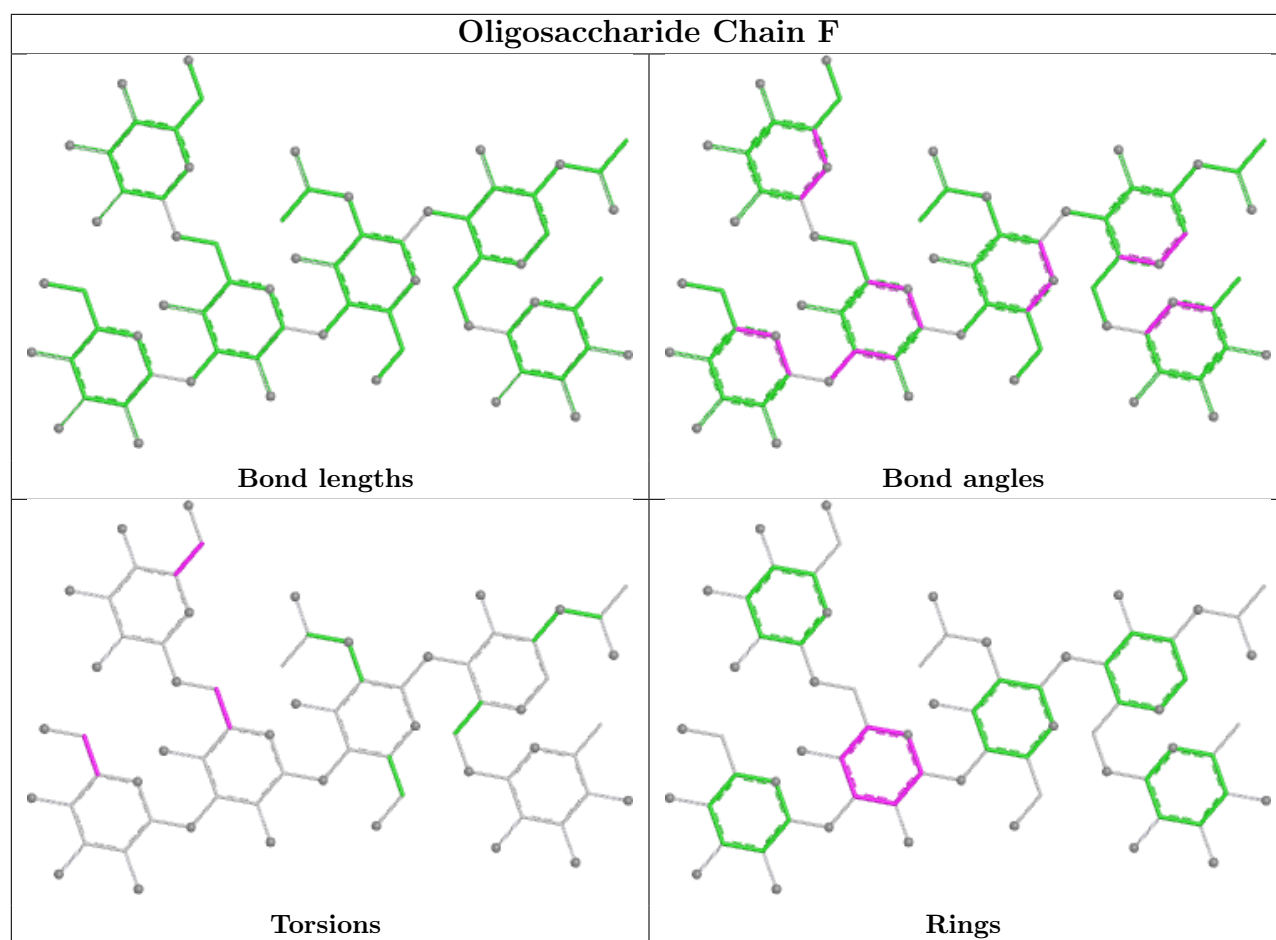
All (2) ring outliers are listed below:

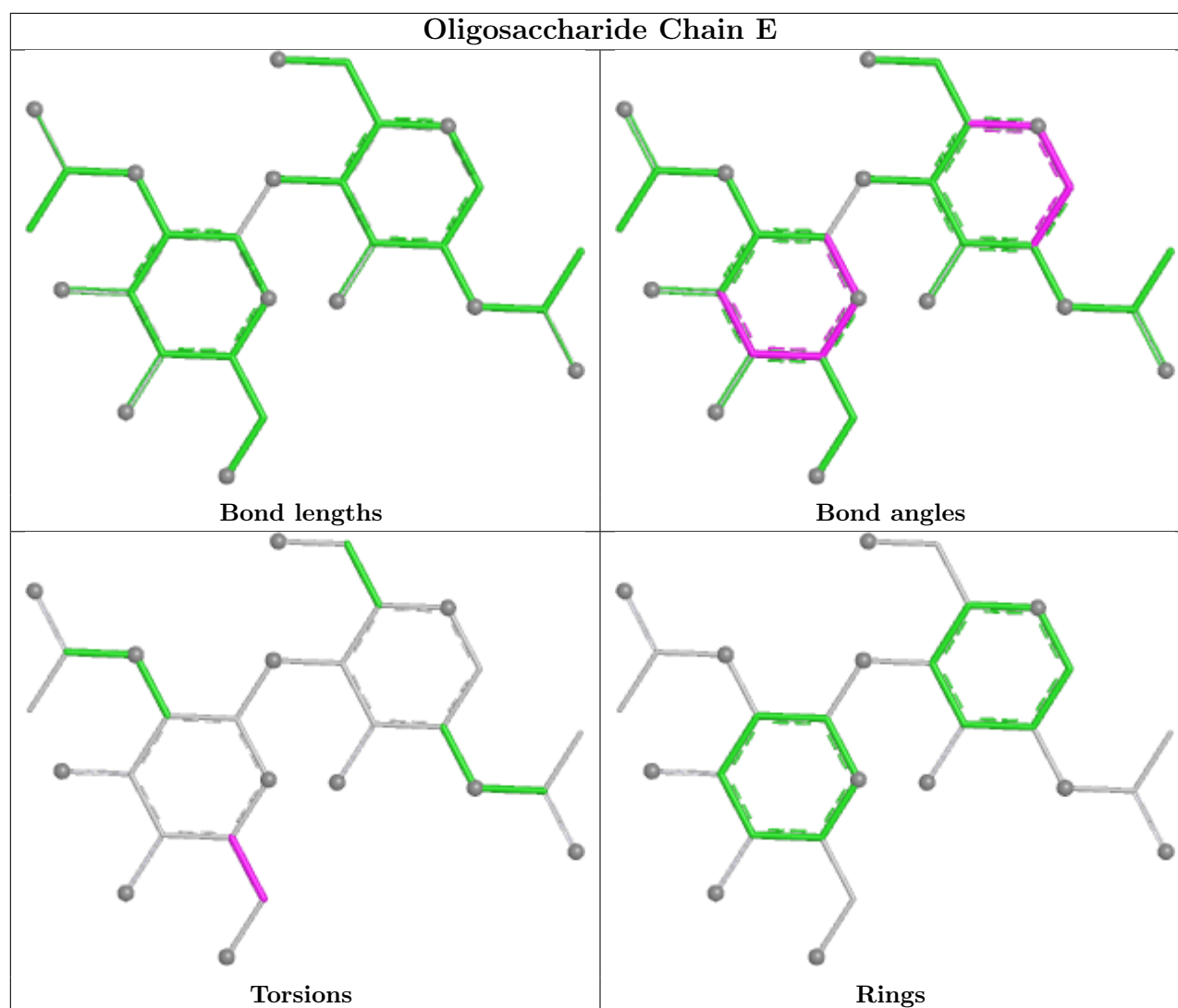
Mol	Chain	Res	Type	Atoms
3	C	3	BMA	C1-C2-C3-C4-C5-O5
3	F	3	BMA	C1-C2-C3-C4-C5-O5

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.







5.6 Ligand geometry [i](#)

2 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	NAG	A	1420	1	14,14,15	0.35	0	17,19,21	2.54	2 (11%)
6	NAG	A	1389	1	14,14,15	0.37	0	17,19,21	2.36	5 (29%)

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	NAG	A	1420	1	-	1/6/23/26	0/1/1/1
6	NAG	A	1389	1	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1420	NAG	C1-O5-C5	7.29	121.96	112.19
6	A	1420	NAG	O5-C1-C2	6.67	121.61	111.29
6	A	1389	NAG	C1-C2-N2	6.41	120.54	110.43
6	A	1389	NAG	C1-O5-C5	5.64	119.74	112.19
6	A	1389	NAG	O5-C1-C2	2.93	115.83	111.29

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1389	NAG	C1-C2-N2-C7
6	A	1389	NAG	O5-C5-C6-O6
6	A	1420	NAG	O5-C5-C6-O6
6	A	1389	NAG	C4-C5-C6-O6

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, 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	613/648 (94%)	0.32	34 (5%) 30 30	32, 65, 95, 141	0
2	B	100/114 (87%)	0.07	6 (6%) 27 28	32, 45, 74, 107	0
All	All	713/762 (93%)	0.29	40 (5%) 30 30	32, 63, 94, 141	0

The worst 5 of 40 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	160	LEU	9.5
2	B	97	ASP	6.0
1	A	614	THR	4.9
1	A	159	HIS	4.2
1	A	602	TYR	4.0

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

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

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	E	2	14/15	0.32	0.20	121,126,128,129	0
4	BMA	D	3	11/12	0.55	0.18	102,106,108,108	0
3	MAN	C	4	11/12	0.71	0.18	102,106,108,109	0
3	MAN	C	5	11/12	0.73	0.18	104,106,108,108	0
3	MAN	F	4	11/12	0.73	0.18	105,106,109,109	0

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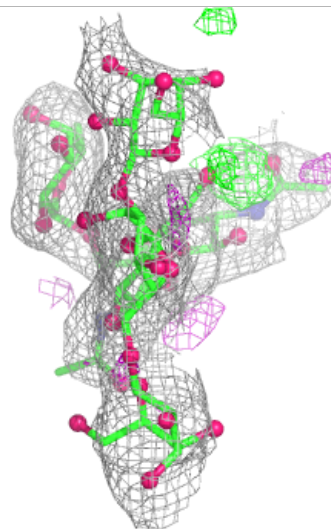
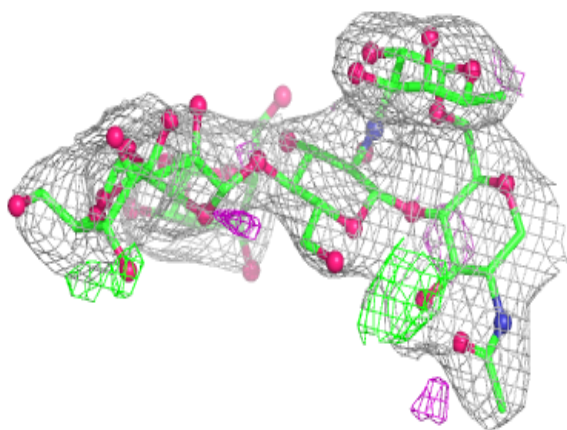
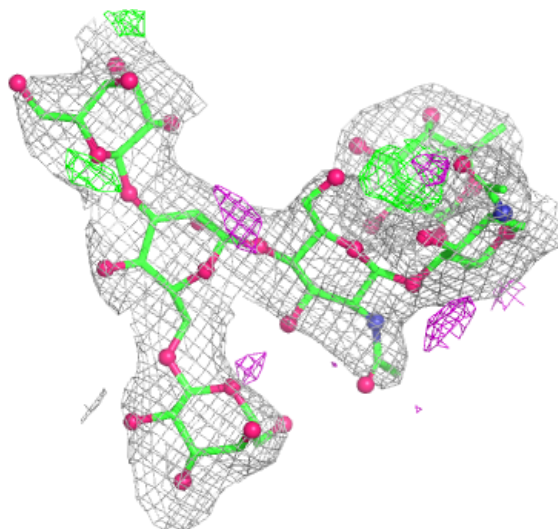
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MAN	F	5	11/12	0.74	0.17	112,116,119,119	0
5	NAG	E	1	14/15	0.77	0.19	124,129,130,130	0
4	NAG	D	2	14/15	0.83	0.17	82,89,96,96	0
3	FUC	F	6	10/11	0.84	0.14	100,102,102,103	0
3	BMA	C	3	11/12	0.85	0.14	97,99,102,106	0
3	NAG	C	2	14/15	0.86	0.14	81,89,94,97	0
3	BMA	F	3	11/12	0.88	0.11	93,96,102,107	0
3	NAG	C	1	14/15	0.93	0.09	31,53,70,70	0
3	NAG	F	2	14/15	0.93	0.11	76,80,83,87	0
3	NAG	F	1	14/15	0.94	0.10	74,84,93,97	0
3	FUC	C	6	10/11	0.94	0.09	66,71,75,76	0
4	NAG	D	1	14/15	0.95	0.09	65,72,76,78	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.

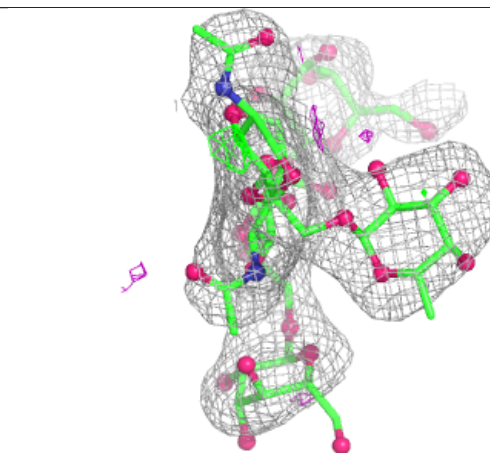
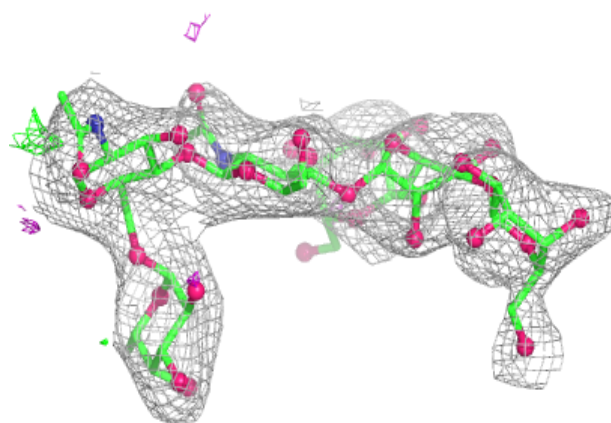
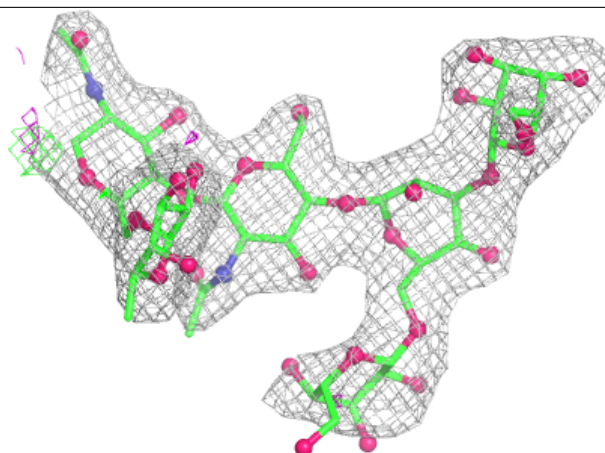
Electron density around Chain C:

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

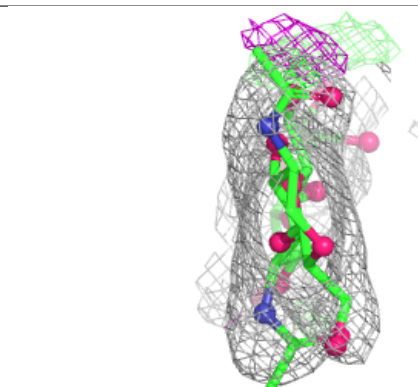
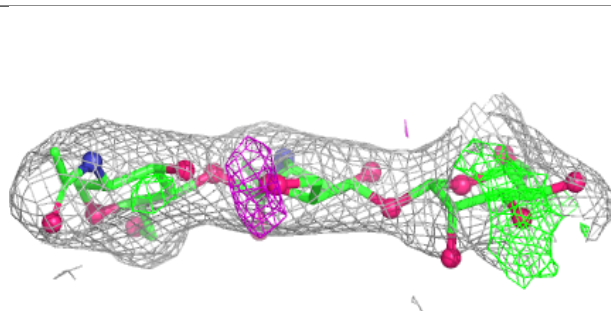
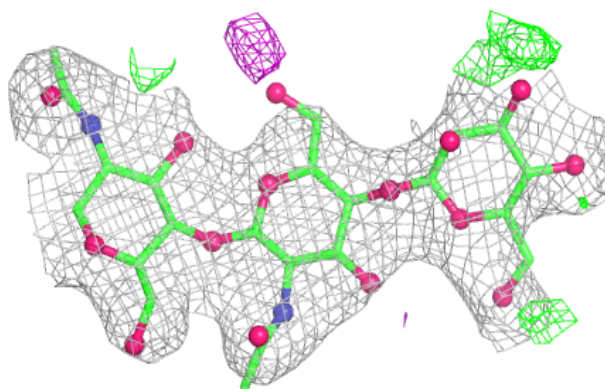


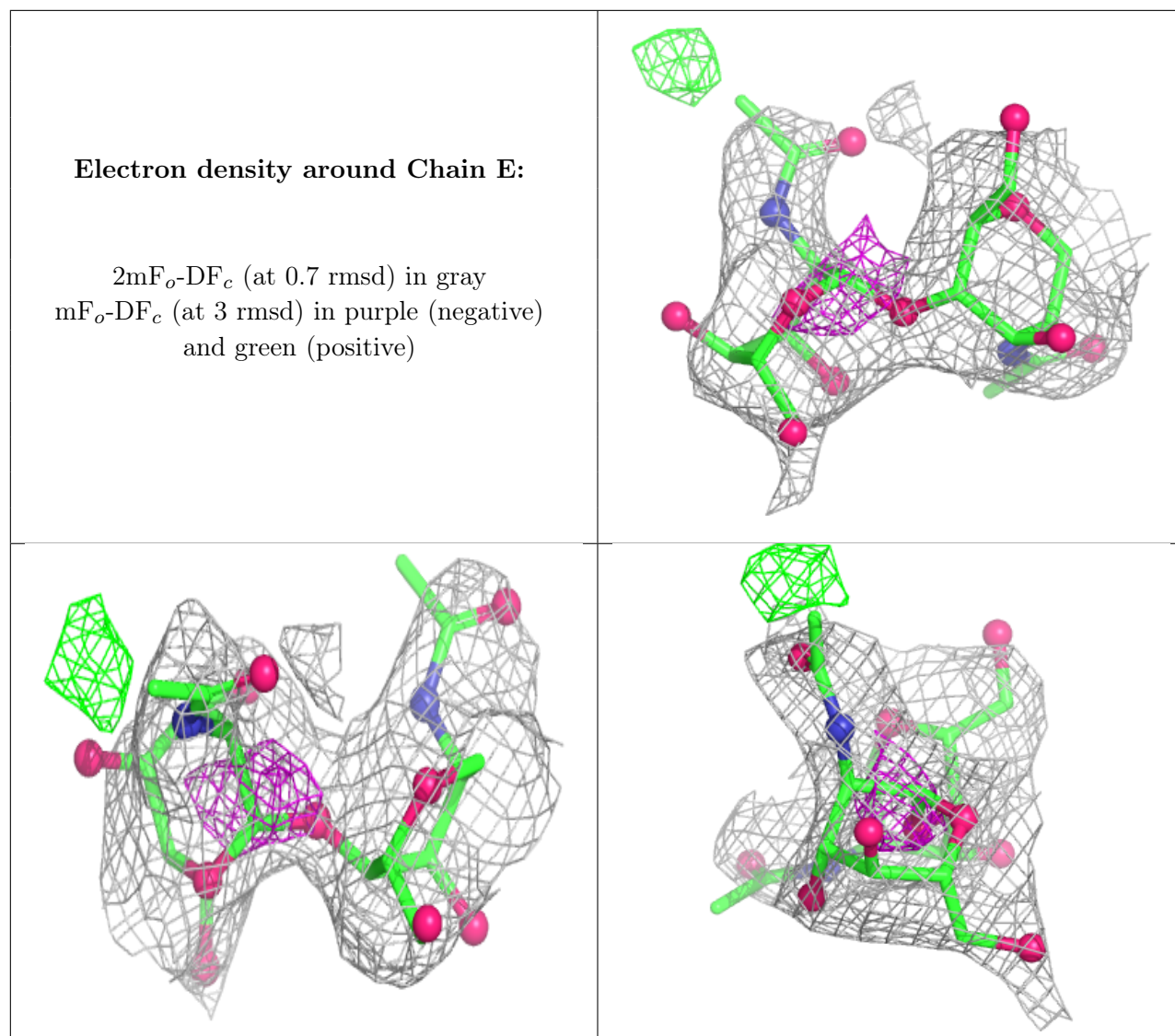
Electron density around Chain F:

$2mF_o-DF_c$ (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, 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
6	NAG	A	1389	14/15	0.41	0.18	138,140,142,144	0
6	NAG	A	1420	14/15	0.55	0.24	115,123,127,128	0

6.5 Other polymers [i](#)

There are no such residues in this entry.