



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:39 AM UTC

PDB ID : 6U6T / pdb_00006u6t
Title : Neuronal growth regulator 1 (NEGR1)
Authors : Machius, M.; Venkannagari, H.; Misra, A.; Rudenko, G.
Deposited on : 2019-08-30
Resolution : 3.01 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 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

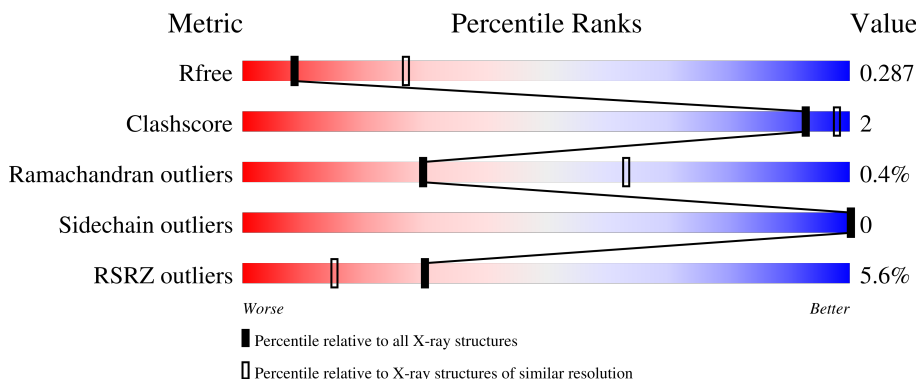
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 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	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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	291	5% 90% 8%
1	B	291	5% 89% 7%
2	C	3	100%
2	E	3	100%
3	D	2	100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8457 atoms, of which 4172 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 Neuronal growth regulator 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	268	Total	C	H	N	O	S	0	1	0
			4113	1301	2033	363	406	10			
1	B	271	Total	C	H	N	O	S	0	0	0
			4121	1306	2034	363	408	10			

There are 30 discrepancies between the modelled and reference sequences:

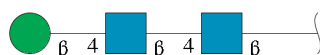
Chain	Residue	Modelled	Actual	Comment	Reference
A	33	GLU	-	expression tag	UNP Q7Z3B1
A	34	LEU	-	expression tag	UNP Q7Z3B1
A	35	GLY	-	expression tag	UNP Q7Z3B1
A	36	THR	-	expression tag	UNP Q7Z3B1
A	37	GLY	-	expression tag	UNP Q7Z3B1
A	314	ALA	-	expression tag	UNP Q7Z3B1
A	315	SER	-	expression tag	UNP Q7Z3B1
A	316	THR	-	expression tag	UNP Q7Z3B1
A	317	SER	-	expression tag	UNP Q7Z3B1
A	318	HIS	-	expression tag	UNP Q7Z3B1
A	319	HIS	-	expression tag	UNP Q7Z3B1
A	320	HIS	-	expression tag	UNP Q7Z3B1
A	321	HIS	-	expression tag	UNP Q7Z3B1
A	322	HIS	-	expression tag	UNP Q7Z3B1
A	323	HIS	-	expression tag	UNP Q7Z3B1
B	33	GLU	-	expression tag	UNP Q7Z3B1
B	34	LEU	-	expression tag	UNP Q7Z3B1
B	35	GLY	-	expression tag	UNP Q7Z3B1
B	36	THR	-	expression tag	UNP Q7Z3B1
B	37	GLY	-	expression tag	UNP Q7Z3B1
B	314	ALA	-	expression tag	UNP Q7Z3B1
B	315	SER	-	expression tag	UNP Q7Z3B1
B	316	THR	-	expression tag	UNP Q7Z3B1
B	317	SER	-	expression tag	UNP Q7Z3B1
B	318	HIS	-	expression tag	UNP Q7Z3B1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	319	HIS	-	expression tag	UNP Q7Z3B1
B	320	HIS	-	expression tag	UNP Q7Z3B1
B	321	HIS	-	expression tag	UNP Q7Z3B1
B	322	HIS	-	expression tag	UNP Q7Z3B1
B	323	HIS	-	expression tag	UNP Q7Z3B1

- Molecule 2 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
2	C	3	Total	C	H	N	O	0	0	0
			73	22	34	2	15			
2	E	3	Total	C	H	N	O	0	0	0
			73	22	34	2	15			

- Molecule 3 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.

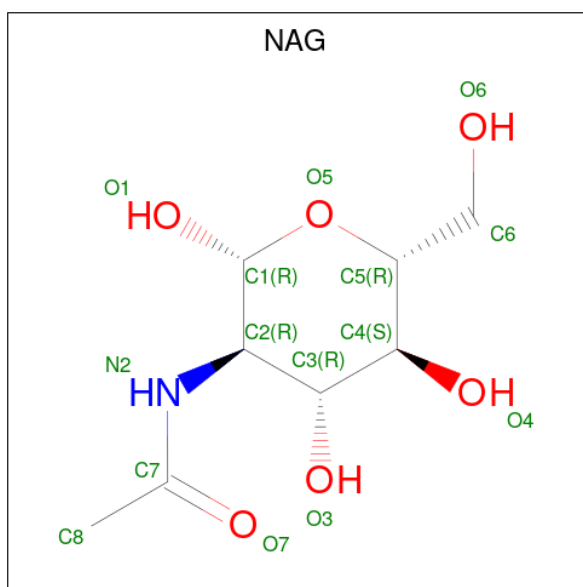


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	2	Total	C	H	N	O	0	0	0
			47	14	23	1	9			

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Ca	0	0
			2	2		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

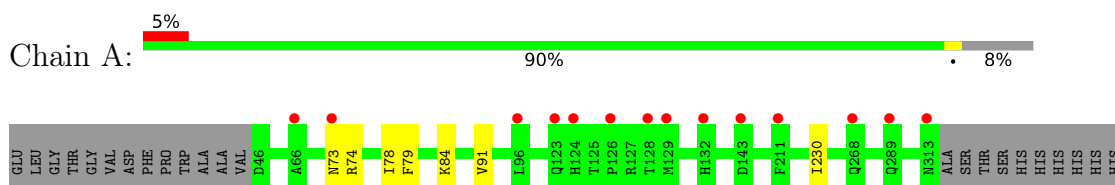


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
5	B	1	28	8	14	1	5	0	0

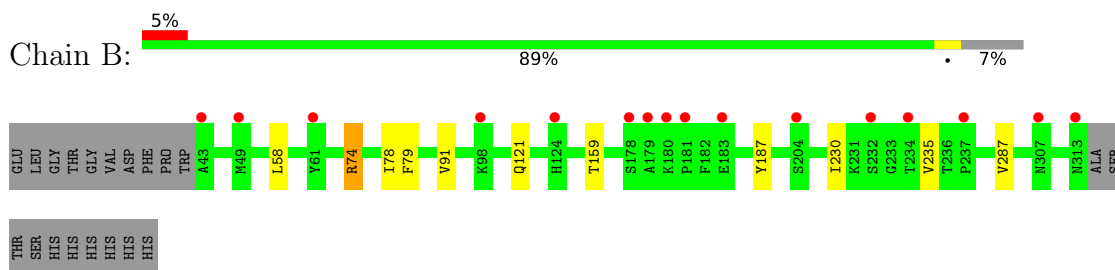
3 Residue-property plots [i](#)

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.

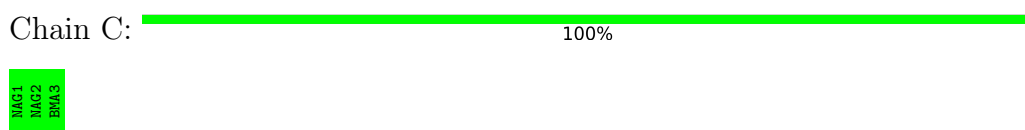
- Molecule 1: Neuronal growth regulator 1



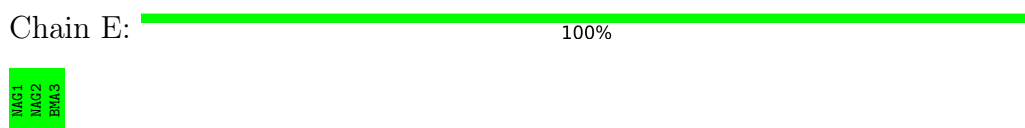
- Molecule 1: Neuronal growth regulator 1



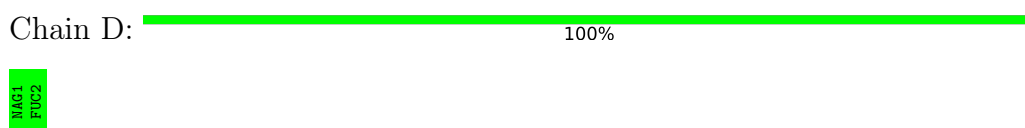
- Molecule 2: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



- Molecule 3: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	73.57Å 97.22Å 252.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.81 – 3.01 31.81 – 3.01	Depositor EDS
% Data completeness (in resolution range)	77.3 (31.81-3.01) 77.5 (31.81-3.01)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.67 (at 3.00Å)	Xtriage
Refinement program	PHENIX 1.16-3549	Depositor
R, R_{free}	0.234 , 0.286 0.233 , 0.287	Depositor DCC
R_{free} test set	500 reflections (2.70%)	wwPDB-VP
Wilson B-factor (Å ²)	49.1	Xtriage
Anisotropy	0.277	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	8457	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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: BMA, CA, NAG, FUC

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	0.10	0/2125	0.26	0/2893
1	B	0.10	0/2131	0.26	0/2902
All	All	0.10	0/4256	0.26	0/5795

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2080	2033	2032	5	1
1	B	2087	2034	2045	8	1
2	C	39	34	34	0	0
2	E	39	34	34	0	0
3	D	24	23	22	0	0
4	A	2	0	0	0	0
5	B	14	14	13	0	0
All	All	4285	4172	4180	13	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (13) 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:B:78:ILE:HG21	1:B:91:VAL:HG11	1.85	0.57
1:A:79:PHE:HE2	1:A:91:VAL:HG12	1.71	0.56
1:A:78:ILE:HG21	1:A:91:VAL:HG11	1.88	0.56
1:B:230:ILE:O	1:B:230:ILE:HG23	2.11	0.50
1:A:73:ASN:HB2	1:A:78:ILE:HD11	1.97	0.46
1:B:74:ARG:HH11	1:B:74:ARG:HG2	1.80	0.46
1:B:159:THR:HG23	1:B:187:TYR:CE1	2.51	0.45
1:A:78:ILE:CG2	1:A:91:VAL:HG11	2.48	0.43
1:B:78:ILE:CG2	1:B:91:VAL:HG11	2.48	0.43
1:B:58:LEU:HD12	1:B:58:LEU:N	2.34	0.43
1:B:235:VAL:HG11	1:B:287:VAL:HG21	1.99	0.43
1:A:230:ILE:HG23	1:A:230:ILE:O	2.18	0.43
1:B:79:PHE:HE2	1:B:91:VAL:HG12	1.84	0.42

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:LYS:O	1:B:121:GLN:HE21[6_445]	1.57	0.03

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	267/291 (92%)	253 (95%)	13 (5%)	1 (0%)	30	63
1	B	269/291 (92%)	256 (95%)	12 (4%)	1 (0%)	30	63
All	All	536/582 (92%)	509 (95%)	25 (5%)	2 (0%)	30	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	74	ARG
1	B	74	ARG

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	234/251 (93%)	234 (100%)	0	100	100
1	B	234/251 (93%)	234 (100%)	0	100	100
All	All	468/502 (93%)	468 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

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](#)

8 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	C	1	2,1	14,14,15	0.20	0	17,19,21	0.50	0
2	NAG	C	2	2	14,14,15	0.20	0	17,19,21	0.42	0
2	BMA	C	3	2	11,11,12	0.59	0	15,15,17	0.68	0
3	NAG	D	1	1,3	14,14,15	0.22	0	17,19,21	0.46	0
3	FUC	D	2	3	10,10,11	0.74	0	14,14,16	0.76	0
2	NAG	E	1	2,1	14,14,15	0.19	0	17,19,21	0.54	0
2	NAG	E	2	2	14,14,15	0.22	0	17,19,21	0.41	0
2	BMA	E	3	2	11,11,12	0.60	0	15,15,17	0.69	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
2	NAG	C	1	2,1	-	5/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
3	NAG	D	1	1,3	-	1/6/23/26	0/1/1/1
3	FUC	D	2	3	-	-	0/1/1/1
2	NAG	E	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	2	NAG	O5-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
2	C	1	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6

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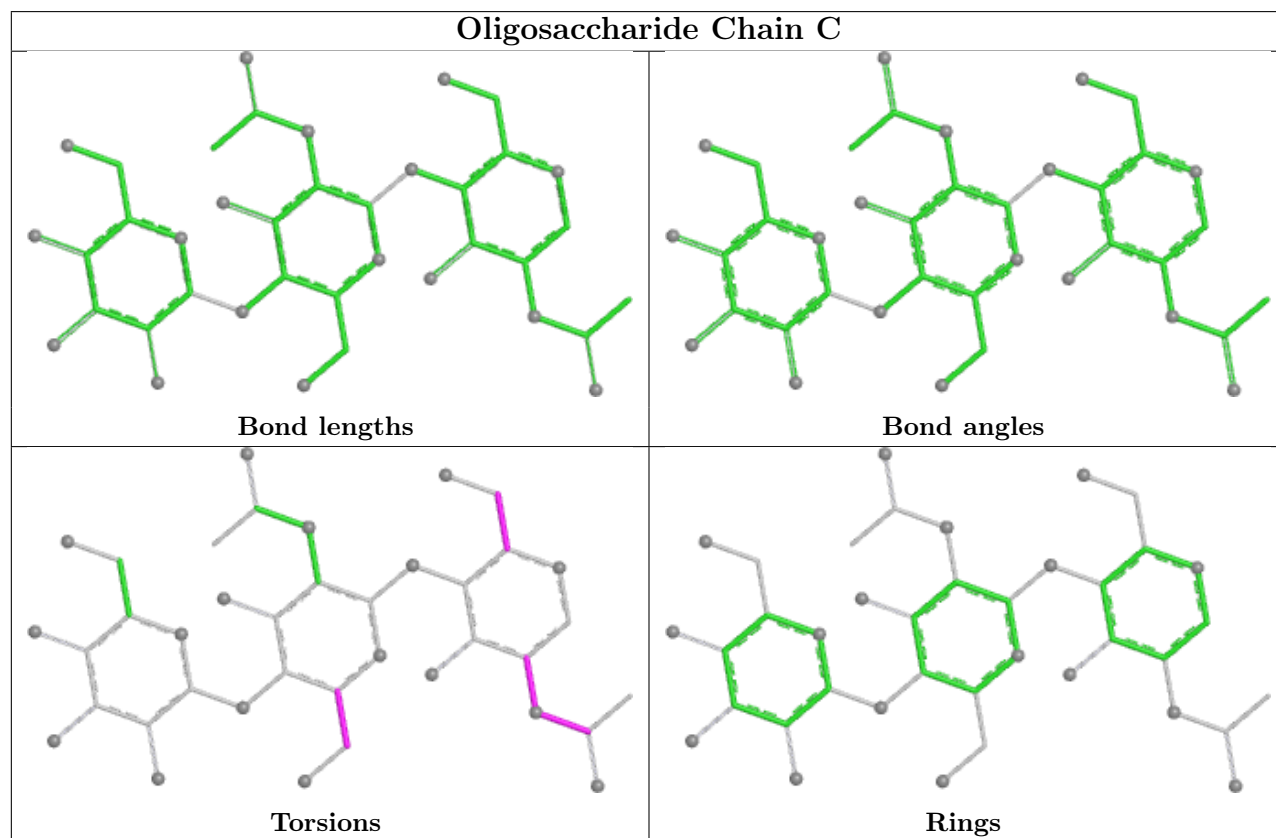
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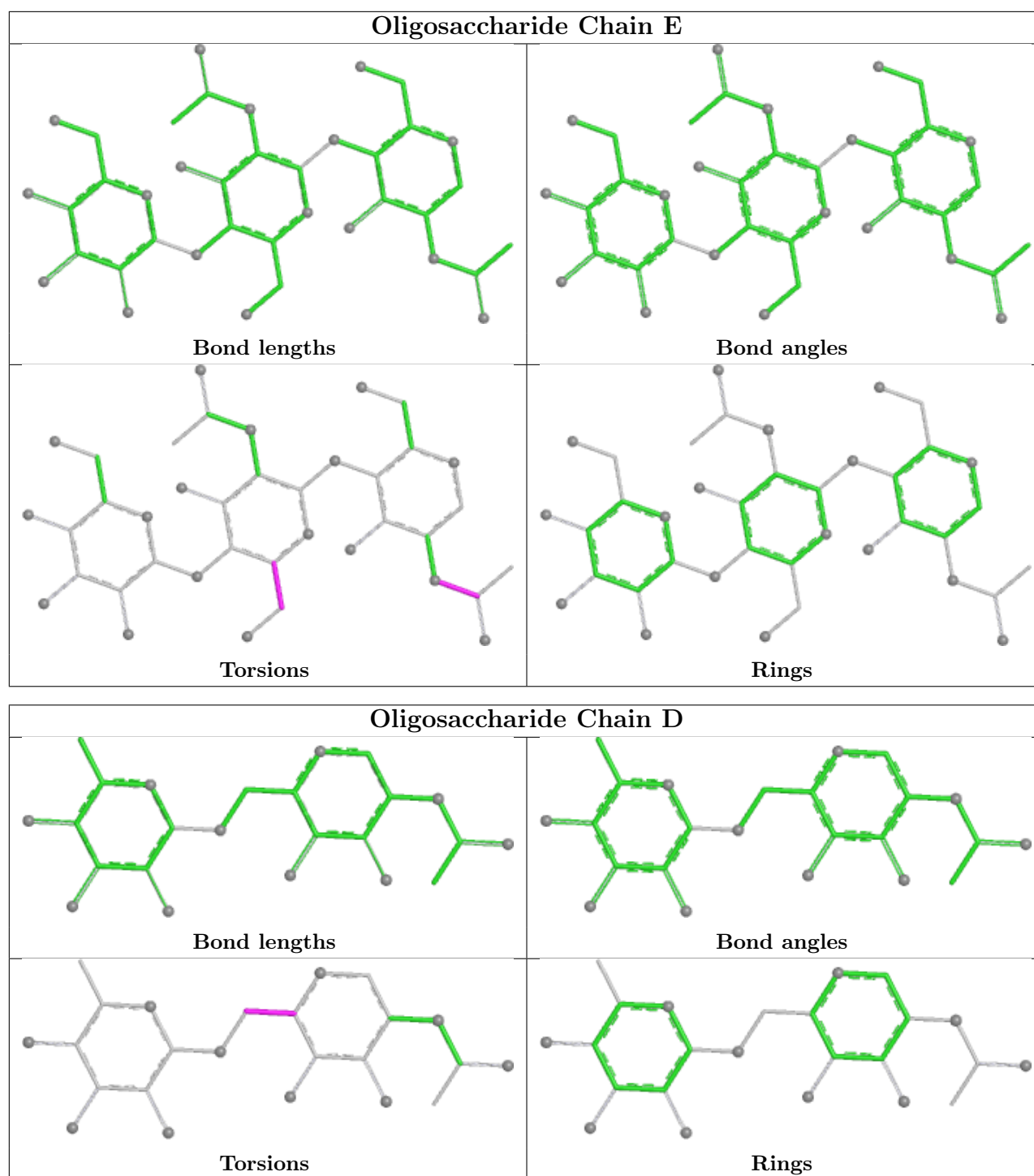
Mol	Chain	Res	Type	Atoms
3	D	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C1-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

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





5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
5	NAG	B	404	1	14,14,15	0.22	0	17,19,21	0.48	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
5	NAG	B	404	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	268/291 (92%)	0.28	14 (5%)	33 16	24, 52, 105, 147	1 (0%)
1	B	271/291 (93%)	0.43	16 (5%)	28 14	27, 59, 105, 169	0
All	All	539/582 (92%)	0.36	30 (5%)	30 15	24, 56, 106, 169	1 (0%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	132[A]	HIS	5.5
1	B	43	ALA	4.9
1	A	129	MET	4.6
1	A	66	ALA	4.1
1	B	307	ASN	3.7
1	A	313	ASN	3.7
1	B	181	PRO	3.6
1	A	73	ASN	3.4
1	B	61	TYR	3.2
1	B	180	LYS	3.2
1	A	128	THR	3.2
1	B	183	GLU	3.0
1	B	179	ALA	2.8
1	B	178	SER	2.8
1	B	234	THR	2.7
1	B	124	HIS	2.5
1	A	143	ASP	2.5
1	B	313	ASN	2.5
1	A	96	LEU	2.5
1	A	123	GLN	2.3
1	B	98	LYS	2.3
1	A	289	GLN	2.2
1	A	126	PRO	2.1
1	B	237	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	124	HIS	2.1
1	B	232	SER	2.1
1	A	211	PHE	2.1
1	A	268	GLN	2.0
1	B	204	SER	2.0
1	B	49	MET	2.0

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](#)

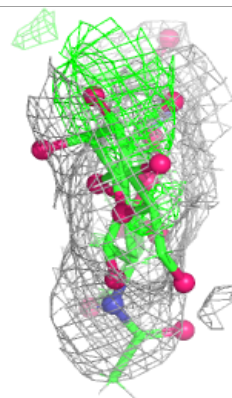
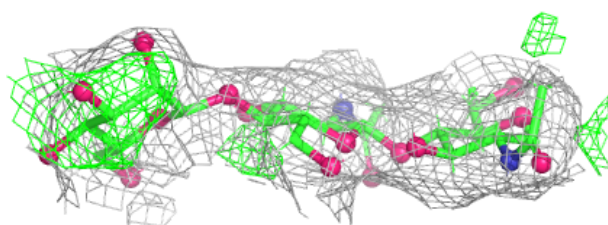
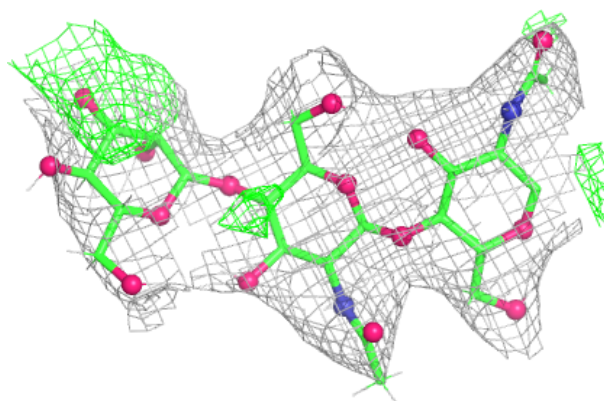
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BMA	C	3	11/12	0.29	0.27	137,161,186,209	0
2	BMA	E	3	11/12	0.72	0.13	47,81,132,142	0
2	NAG	C	2	14/15	0.78	0.21	63,88,115,117	0
3	FUC	D	2	10/11	0.82	0.12	51,70,90,104	0
2	NAG	E	2	14/15	0.85	0.11	52,63,76,92	0
3	NAG	D	1	14/15	0.86	0.11	50,69,93,93	0
2	NAG	E	1	14/15	0.89	0.14	53,64,80,81	0
2	NAG	C	1	14/15	0.91	0.13	54,62,72,74	0

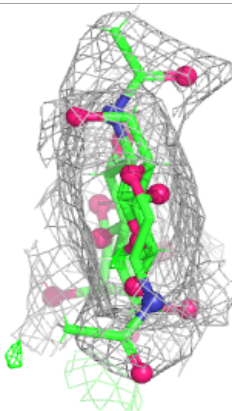
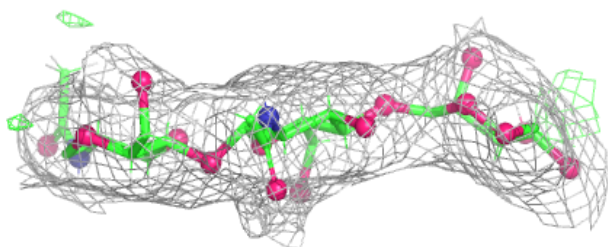
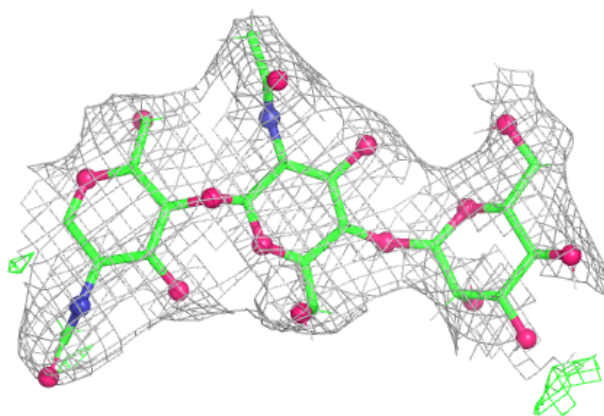
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

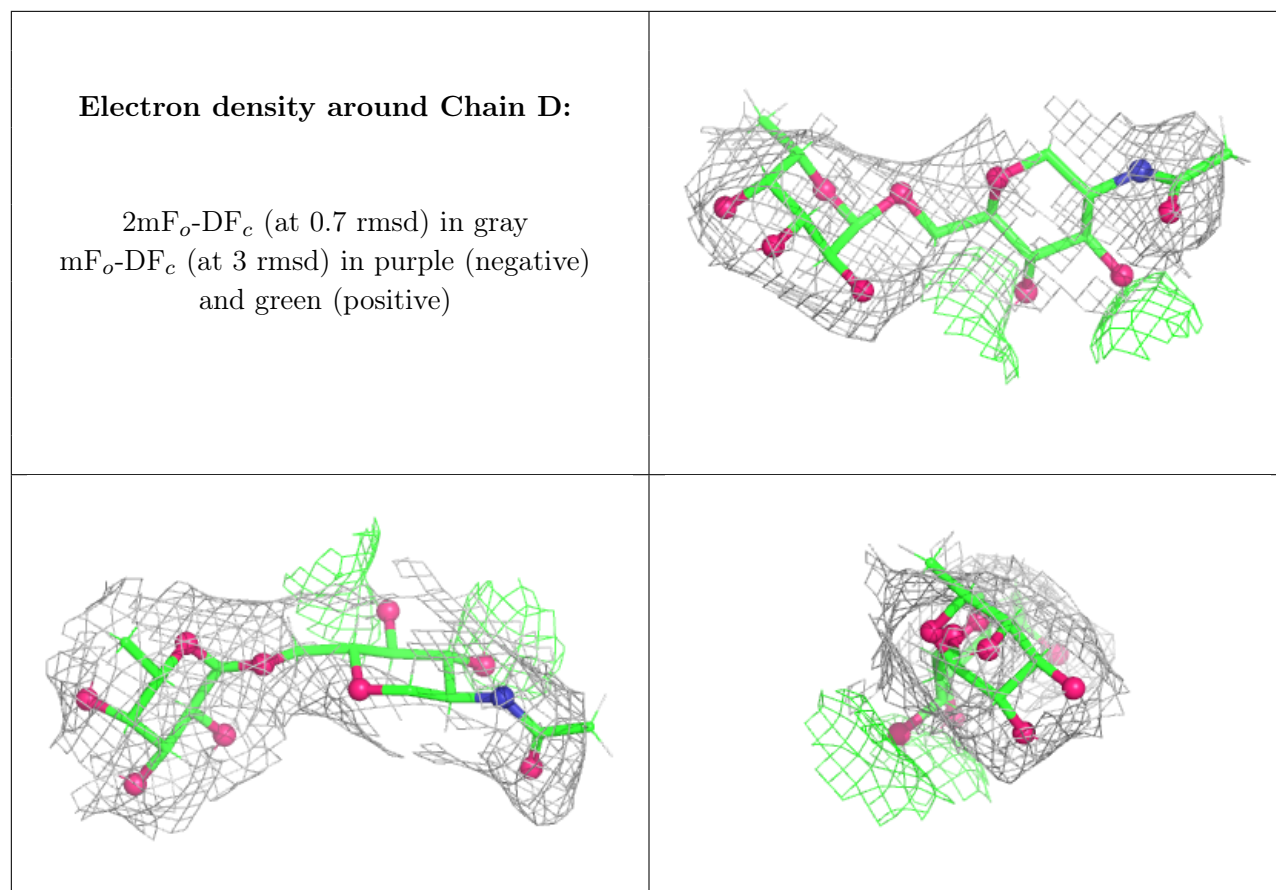
Electron density around Chain C:

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

**Electron density around Chain E:**

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





6.4 Ligands [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
4	CA	A	407	1/1	0.42	0.30	98,98,98,98	1
5	NAG	B	404	14/15	0.60	0.18	56,73,114,114	0
4	CA	A	406	1/1	0.86	0.18	58,58,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.