



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 10, 2020 – 12:30 PM BST

PDB ID : 5JLV
Title : Receptor binding domain of Botulinum neurotoxin A in complex with human glycosylated SV2C
Authors : Yao, G.; Zhang, S.; Mahrhold, S.; Lam, K.; Stern, D.; Bagramyan, K.; Perry, K.; Kalkum, M.; Rummel, A.; Dong, M.; Jin, R.
Deposited on : 2016-04-27
Resolution : 2.00 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 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 : 2.13.1
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) : 2.13.1

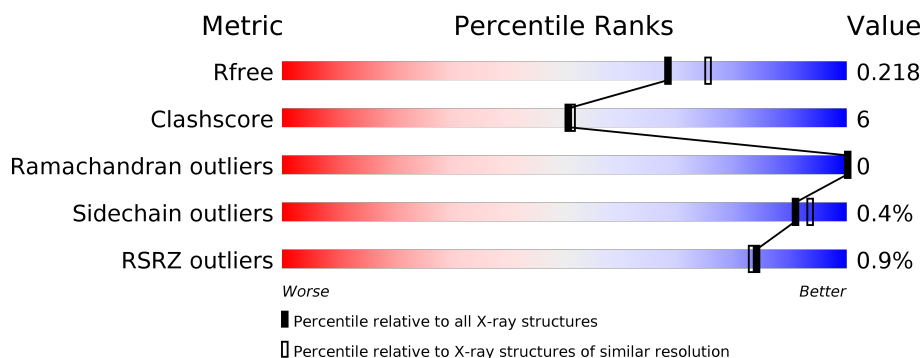
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 on 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	428	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 14%, green 86%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 86% 14% • </div> </div>
1	B	428	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 9%, green 90%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 90% 9% • </div> </div>
2	C	117	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 68%, yellow 15%, grey 17%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 68% 15% 17% </div> </div>
2	D	117	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 69%, yellow 9%, grey 22%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 69% 9% 22% </div> </div>
3	E	4	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 50%, yellow 50%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 50% 50% </div> </div>
4	F	3	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 67%, yellow 33%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 67% 33% </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	PO4	A	1303	-	X	X	-
6	PO4	B	1303	-	X	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 9428 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 Botulinum neurotoxin type A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	425	Total	C	N	O	S	3	5	0
			3538	2252	613	659	14			
1	B	425	Total	C	N	O	S	4	3	0
			3527	2246	610	657	14			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	869	GLY	-	expression tag	UNP P10845
A	870	PRO	-	expression tag	UNP P10845
A	871	MET	-	expression tag	UNP P10845
A	1158	ALA	THR	engineered mutation	UNP P10845
B	869	GLY	-	expression tag	UNP P10845
B	870	PRO	-	expression tag	UNP P10845
B	871	MET	-	expression tag	UNP P10845
B	1158	ALA	THR	engineered mutation	UNP P10845

- Molecule 2 is a protein called Synaptic vesicle glycoprotein 2C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	97	Total	C	N	O	S	0	0	0
			812	521	129	157	5			
2	D	91	Total	C	N	O	S	0	0	0
			764	493	118	148	5			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	451	SER	-	expression tag	UNP Q496J9
C	452	HIS	-	expression tag	UNP Q496J9
C	453	HIS	-	expression tag	UNP Q496J9
C	454	HIS	-	expression tag	UNP Q496J9

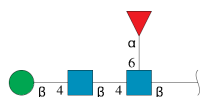
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Chain	Residue	Modelled	Actual	Comment	Reference
C	455	HIS	-	expression tag	UNP Q496J9
C	456	HIS	-	expression tag	UNP Q496J9
C	457	HIS	-	expression tag	UNP Q496J9
C	458	HIS	-	expression tag	UNP Q496J9
C	459	HIS	-	expression tag	UNP Q496J9
C	460	HIS	-	expression tag	UNP Q496J9
C	461	SER	-	expression tag	UNP Q496J9
C	462	GLY	-	expression tag	UNP Q496J9
C	463	GLY	-	expression tag	UNP Q496J9
C	464	GLY	-	expression tag	UNP Q496J9
C	465	SER	-	expression tag	UNP Q496J9
C	466	GLY	-	expression tag	UNP Q496J9
C	467	GLY	-	expression tag	UNP Q496J9
C	468	GLY	-	expression tag	UNP Q496J9
C	469	ILE	-	expression tag	UNP Q496J9
C	470	GLU	-	expression tag	UNP Q496J9
C	471	GLY	-	expression tag	UNP Q496J9
C	472	ARG	-	expression tag	UNP Q496J9
D	451	SER	-	expression tag	UNP Q496J9
D	452	HIS	-	expression tag	UNP Q496J9
D	453	HIS	-	expression tag	UNP Q496J9
D	454	HIS	-	expression tag	UNP Q496J9
D	455	HIS	-	expression tag	UNP Q496J9
D	456	HIS	-	expression tag	UNP Q496J9
D	457	HIS	-	expression tag	UNP Q496J9
D	458	HIS	-	expression tag	UNP Q496J9
D	459	HIS	-	expression tag	UNP Q496J9
D	460	HIS	-	expression tag	UNP Q496J9
D	461	SER	-	expression tag	UNP Q496J9
D	462	GLY	-	expression tag	UNP Q496J9
D	463	GLY	-	expression tag	UNP Q496J9
D	464	GLY	-	expression tag	UNP Q496J9
D	465	SER	-	expression tag	UNP Q496J9
D	466	GLY	-	expression tag	UNP Q496J9
D	467	GLY	-	expression tag	UNP Q496J9
D	468	GLY	-	expression tag	UNP Q496J9
D	469	ILE	-	expression tag	UNP Q496J9
D	470	GLU	-	expression tag	UNP Q496J9
D	471	GLY	-	expression tag	UNP Q496J9
D	472	ARG	-	expression tag	UNP Q496J9

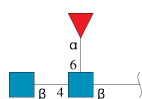
- Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopy

ranose.



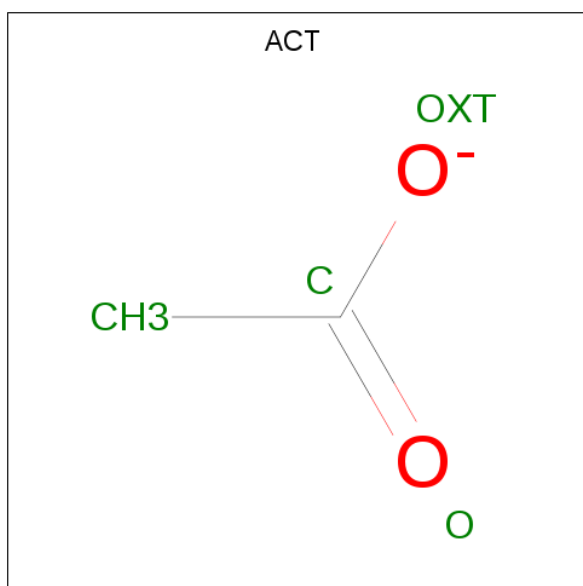
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	4	Total	C	N	O	0	0	0
			49	28	2	19			

- Molecule 4 is an oligosaccharide called 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
4	F	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	P	0	0
			5	4	1		
6	A	1	Total	O	P	0	0
			5	4	1		
6	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	276	Total	O	0	0
			276	276		
8	B	270	Total	O	0	0
			270	270		
8	C	81	Total	O	0	0
			81	81		
8	D	32	Total	O	0	0
			32	32		

Chain E:  50% 50%


MAG1
MAG2
BGL3
FUC4

- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  67% 33%


MAG1
MAG2
FUC3

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.00Å 111.85Å 126.25Å 90.00° 101.29° 90.00°	Depositor
Resolution (Å)	123.81 – 2.00 123.81 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.5 (123.81-2.00) 93.8 (123.81-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.175 , 0.216 0.179 , 0.218	Depositor DCC
R_{free} test set	4976 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.415	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 57.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9428	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage'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.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, PO4, BMA, 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	0.48	0/3614	0.54	0/4885
1	B	0.43	0/3600	0.53	0/4866
2	C	0.47	0/833	0.51	0/1119
2	D	0.39	0/785	0.44	0/1056
All	All	0.45	0/8832	0.53	0/11926

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	3538	0	3498	49	0
1	B	3527	0	3486	30	0
2	C	812	0	749	11	0
2	D	764	0	700	10	0
3	E	49	0	43	2	0
4	F	38	0	34	4	0
5	A	4	0	3	1	0
5	B	8	0	6	1	0
6	A	10	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	5	0	0	4	0
7	C	14	0	13	0	0
8	A	276	0	0	16	0
8	B	270	0	0	9	0
8	C	81	0	0	3	0
8	D	32	0	0	0	0
All	All	9428	0	8532	99	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 6.

The worst 5 of 99 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:1294:ARG:HG2	8:A:1430:HOH:O	1.72	0.89
1:A:996:ARG:HH22	6:A:1303:PO4:P	1.97	0.87
1:B:996:ARG:NH2	6:B:1303:PO4:O2	2.07	0.87
1:B:996:ARG:HH22	6:B:1303:PO4:P	2.07	0.76
2:C:476:ASP:HA	8:C:733:HOH:O	1.88	0.74

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	426/428 (100%)	408 (96%)	18 (4%)	0	100	100
1	B	424/428 (99%)	405 (96%)	19 (4%)	0	100	100
2	C	95/117 (81%)	94 (99%)	1 (1%)	0	100	100
2	D	89/117 (76%)	88 (99%)	1 (1%)	0	100	100
All	All	1034/1090 (95%)	995 (96%)	39 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	397/393 (101%)	396 (100%)	1 (0%)	92	95
1	B	395/393 (100%)	393 (100%)	2 (0%)	88	92
2	C	92/106 (87%)	92 (100%)	0	100	100
2	D	87/106 (82%)	86 (99%)	1 (1%)	73	78
All	All	971/998 (97%)	967 (100%)	4 (0%)	91	93

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1065	ARG
1	B	937	MET
1	B	1065	ARG
2	D	563	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	1106	GLN
1	B	1172	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 ⓘ

7 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	E	1	3,2	14,14,15	0.30	0	17,19,21	0.61	0
3	NAG	E	2	3	14,14,15	0.29	0	17,19,21	0.62	0
3	BMA	E	3	3	11,11,12	0.27	0	15,15,17	0.64	0
3	FUC	E	4	3	10,10,11	0.29	0	14,14,16	0.62	0
4	NAG	F	1	2,4	14,14,15	0.29	0	17,19,21	0.62	0
4	NAG	F	2	4	14,14,15	0.30	0	17,19,21	0.61	0
4	FUC	F	3	4	10,10,11	0.30	0	14,14,16	0.63	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
3	NAG	E	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
3	FUC	E	4	3	-	-	0/1/1/1
4	NAG	F	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	1/6/23/26	0/1/1/1
4	FUC	F	3	4	-	-	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

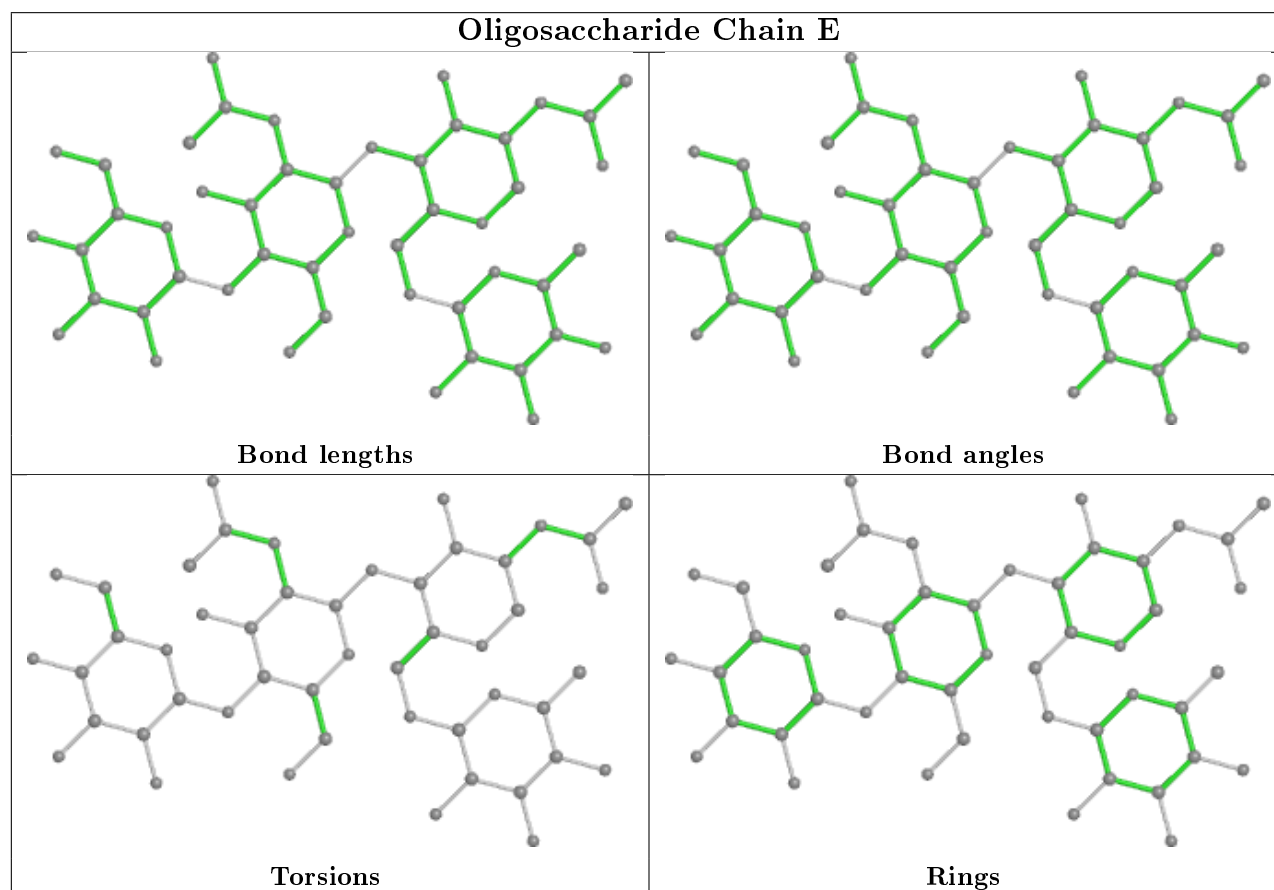
Mol	Chain	Res	Type	Atoms
4	F	2	NAG	O5-C5-C6-O6

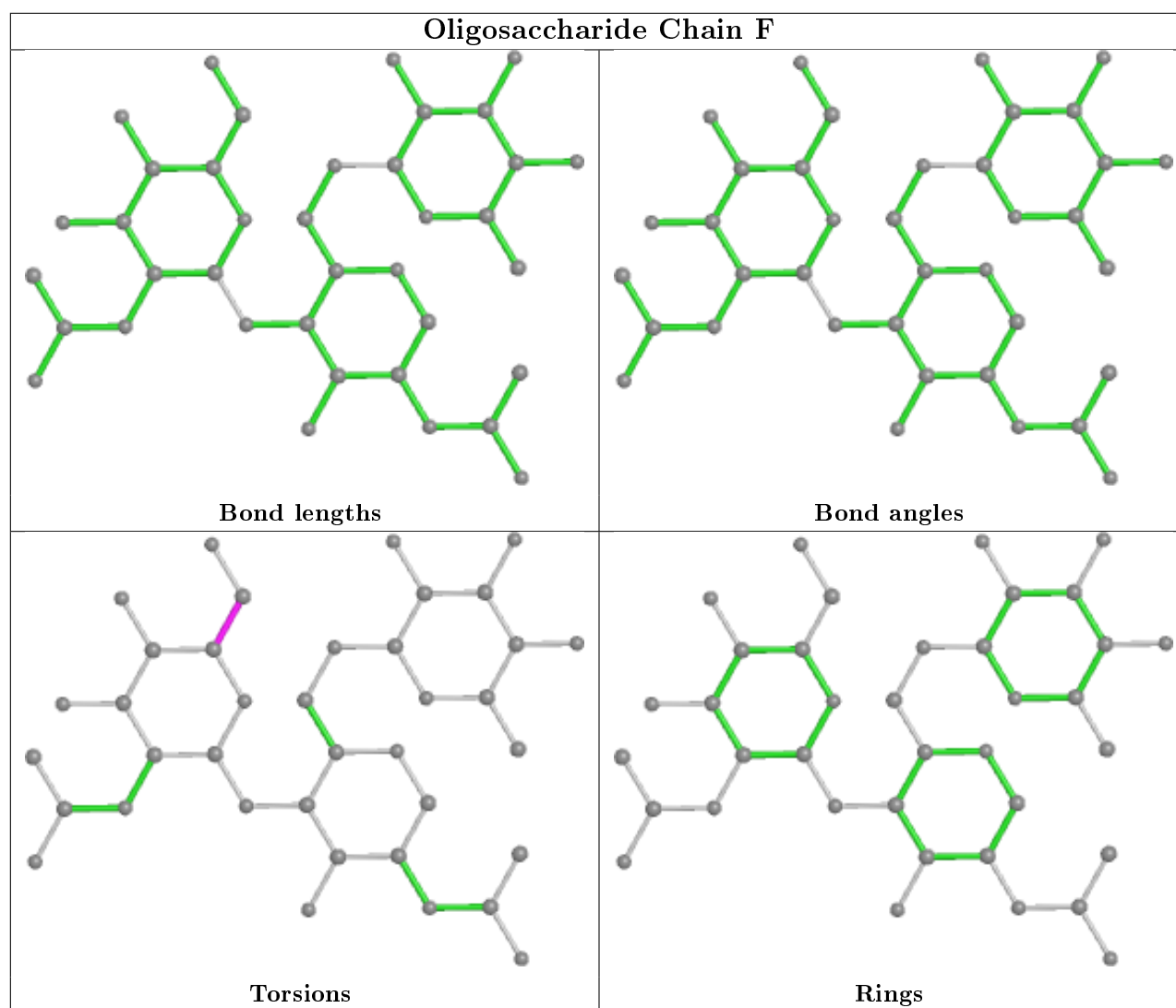
There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	4	FUC	1	0
3	E	1	NAG	1	0
4	F	1	NAG	4	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](#)

7 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	PO4	A	1302	-	4,4,4	0.58	0	6,6,6	0.69	0
5	ACT	B	1302	-	1,3,3	1.68	0	0,3,3	0.00	-
6	PO4	B	1303	-	4,4,4	2.96	4 (100%)	6,6,6	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	ACT	B	1301	-	1,3,3	1.34	0	0,3,3	0.00	-
7	NAG	C	605	2	14,14,15	0.30	0	17,19,21	0.62	0
6	PO4	A	1303	-	4,4,4	2.96	4 (100%)	6,6,6	0.43	0
5	ACT	A	1301	-	1,3,3	2.64	1 (100%)	0,3,3	0.00	-

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
7	NAG	C	605	2	-	2/6/23/26	0/1/1/1

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	1303	PO4	P-O1	4.51	1.61	1.50
6	A	1303	PO4	P-O1	4.51	1.61	1.50
5	A	1301	ACT	CH3-C	2.64	1.52	1.48
6	B	1303	PO4	P-O3	2.30	1.61	1.54
6	A	1303	PO4	P-O3	2.30	1.61	1.54

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	C	605	NAG	O5-C5-C6-O6
7	C	605	NAG	C4-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1303	PO4	4	0
5	B	1301	ACT	1	0
6	A	1303	PO4	5	0
5	A	1301	ACT	1	0

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	425/428 (99%)	-0.21	3 (0%) 87 87	19, 38, 68, 105	1 (0%)
1	B	425/428 (99%)	-0.20	6 (1%) 75 74	20, 42, 77, 104	1 (0%)
2	C	97/117 (82%)	-0.30	0 100 100	27, 42, 67, 83	0
2	D	91/117 (77%)	0.06	0 100 100	32, 58, 81, 115	0
All	All	1038/1090 (95%)	-0.19	9 (0%) 84 83	19, 41, 75, 115	2 (0%)

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1273	ARG	4.6
1	B	1296	LEU	4.2
1	A	1229	GLN	3.2
1	A	1170	LYS	2.5
1	B	1256	ASN	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, 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.

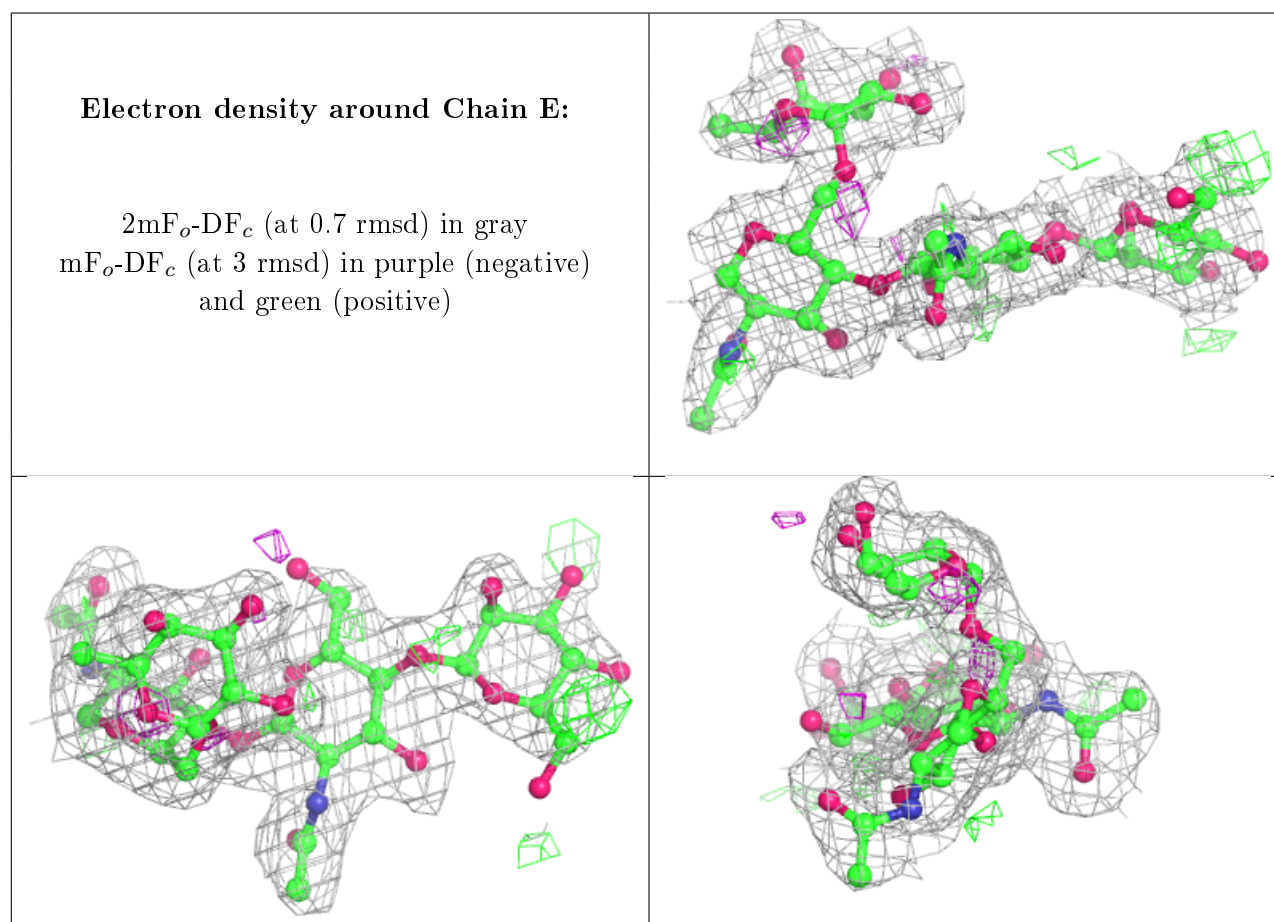
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	BMA	E	3	11/12	0.92	0.19	56,72,81,88	1
4	NAG	F	1	14/15	0.93	0.12	36,41,46,49	0

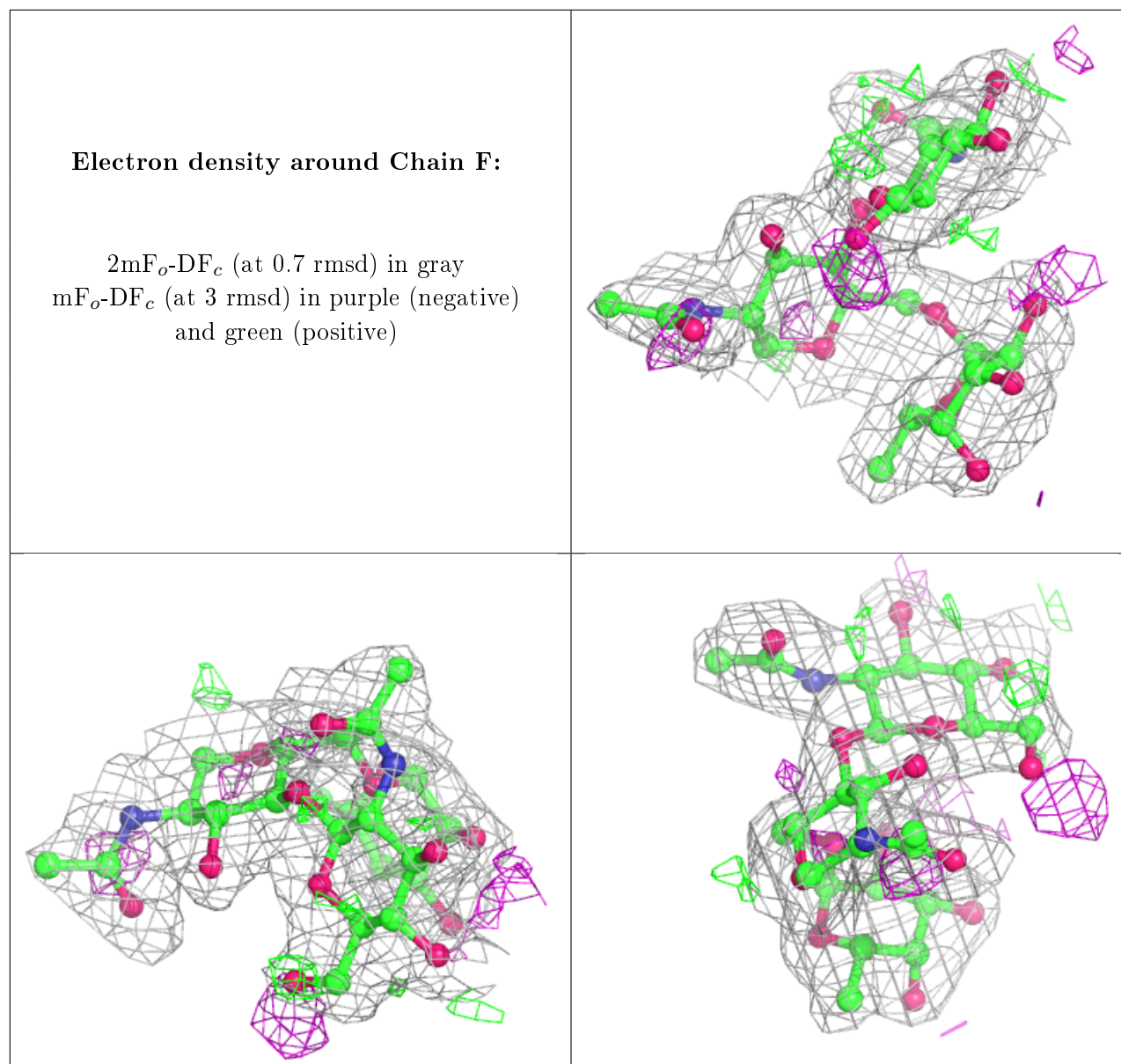
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	NAG	F	2	14/15	0.93	0.09	37,43,60,70	0
3	NAG	E	1	14/15	0.95	0.09	30,37,46,46	0
3	FUC	E	4	10/11	0.96	0.09	32,37,42,45	0
3	NAG	E	2	14/15	0.96	0.13	32,38,57,84	0
4	FUC	F	3	10/11	0.96	0.10	41,52,57,62	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.





6.4 Ligands ⓘ

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	ACT	B	1302	4/4	0.68	0.21	64,75,76,84	0
5	ACT	A	1301	4/4	0.82	0.21	53,59,66,68	0
7	NAG	C	605	14/15	0.84	0.22	63,82,90,93	0
6	PO4	B	1303	5/5	0.84	0.22	105,112,130,135	0
6	PO4	A	1303	5/5	0.89	0.24	81,109,143,149	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	ACT	B	1301	4/4	0.92	0.10	45,49,53,57	0
6	PO4	A	1302	5/5	0.94	0.11	43,51,56,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.