



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:06 PM UTC

PDB ID : 7CYV / pdb_00007cyv
Title : Crystal structure of FD20, a neutralizing single-chain variable fragment (scFv) in complex with SARS-CoV-2 Spike receptor-binding domain (RBD)
Authors : Li, Y.; Li, T.; Lai, Y.; Cai, H.; Yao, H.; Li, D.
Deposited on : 2020-09-04
Resolution : 3.13 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

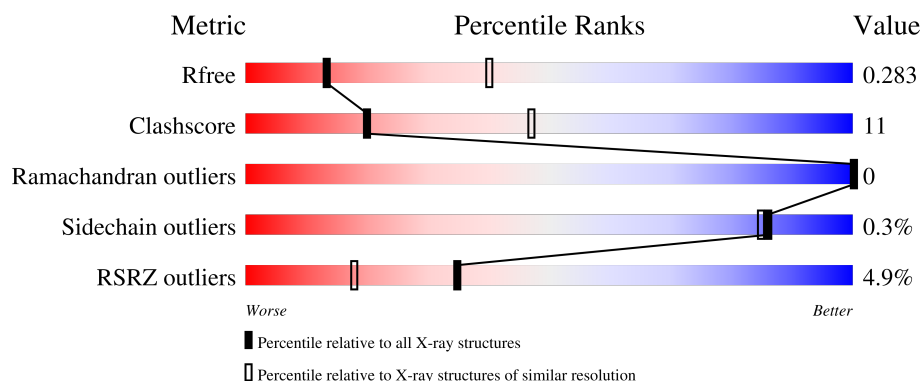
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.13 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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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	H	281	3% 63% 21% 16%
2	B	213	6% 70% 20% 9%
3	A	5	20% 40% 40%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3225 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 The heavy chain variable region of the scFv FD20, The light chain variable region of the scFv FD20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	236	Total	C	N	O	S	0	0	0
			1700	1085	275	332	8			

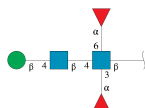
- Molecule 2 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	193	Total	C	N	O	P	S	0	0
			1466	937	248	272	1	8		

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	327	ALA	-	expression tag	UNP P0DTC2
B	328	GLY	-	expression tag	UNP P0DTC2
B	329	SER	-	expression tag	UNP P0DTC2
B	532	GLY	-	expression tag	UNP P0DTC2
B	533	THR	-	expression tag	UNP P0DTC2
B	534	LEU	-	expression tag	UNP P0DTC2
B	535	GLU	-	expression tag	UNP P0DTC2
B	536	VAL	-	expression tag	UNP P0DTC2
B	537	LEU	-	expression tag	UNP P0DTC2
B	538	PHE	-	expression tag	UNP P0DTC2
B	539	GLN	-	expression tag	UNP P0DTC2

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	A	5	Total	C	N	O	0	0	0
			59	34	2	23			

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	206.96Å 57.92Å 47.21Å 90.00° 100.43° 90.00°	Depositor
Resolution (Å)	46.43 – 3.13 46.43 – 3.13	Depositor EDS
% Data completeness (in resolution range)	98.9 (46.43-3.13) 99.2 (46.43-3.13)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.252 , 0.276 0.254 , 0.283	Depositor DCC
R_{free} test set	463 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	68.8	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.030 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3225	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.08% 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, BMA, FUC, TPO

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	H	0.17	0/1740	0.41	0/2383
2	B	0.18	0/1495	0.45	1/2040 (0.0%)
All	All	0.18	0/3235	0.43	1/4423 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	476	GLY	N-CA-C	6.13	117.75	111.56

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	H	1700	0	1560	44	0
2	B	1466	0	1315	25	0
3	A	59	0	52	1	0
All	All	3225	0	2927	70	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 11.

All (70) 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)
3:A:2:NAG:H61	3:A:4:FUC:H63	1.62	0.81
1:H:97:ALA:HB3	1:H:114:MET:HE2	1.63	0.78
2:B:391:CYS:HA	2:B:525:CYS:HB3	1.69	0.73
1:H:38:ARG:NH1	1:H:90:ASP:OD1	2.22	0.71
1:H:39:GLN:OE1	1:H:1037:GLN:NE2	2.22	0.71
1:H:47:TRP:CD2	1:H:1097:VAL:HG22	2.26	0.71
1:H:1090:ALA:HA	1:H:1097:VAL:HA	1.75	0.67
2:B:365:TYR:HD2	2:B:388:ASN:HA	1.60	0.67
1:H:35:SER:HB3	1:H:114:MET:HE3	1.78	0.65
1:H:48:MET:HG2	1:H:64:PHE:HE2	1.60	0.64
1:H:51:ILE:HG13	1:H:58:ALA:HB2	1.82	0.62
2:B:417:LYS:HD2	2:B:455:LEU:HD12	1.82	0.62
1:H:1060:ARG:NH1	1:H:1081:ASP:OD2	2.33	0.61
1:H:39:GLN:NE2	1:H:43:GLN:O	2.31	0.61
1:H:1074:ILE:HB	1:H:1077:VAL:HG12	1.84	0.60
1:H:1079:ALA:HA	1:H:1107:VAL:HG11	1.84	0.59
2:B:484:GLU:OE2	2:B:484:GLU:N	2.33	0.59
2:B:433:VAL:HG22	2:B:512:VAL:HG13	1.86	0.57
1:H:47:TRP:HZ2	1:H:50:ARG:HG2	1.69	0.57
1:H:67:ARG:NH2	1:H:90:ASP:OD2	2.38	0.56
1:H:35:SER:OG	1:H:47:TRP:NE1	2.37	0.56
1:H:6:GLN:HG2	1:H:22:CYS:HB3	1.89	0.54
1:H:37:VAL:HG23	1:H:114:MET:HE1	1.88	0.54
1:H:59:ASN:ND2	1:H:1095:THR:O	2.40	0.54
1:H:100:GLU:O	1:H:112:TYR:HB2	2.06	0.54
2:B:365:TYR:CD2	2:B:388:ASN:HA	2.41	0.54
1:H:1036:GLN:HB2	1:H:1085:TYR:CE1	2.43	0.53
1:H:14:PRO:HG2	1:H:127:SER:HB2	1.90	0.53
1:H:50:ARG:HD2	1:H:111:TYR:CZ	2.44	0.52
2:B:386:LYS:O	2:B:390:LEU:HD23	2.10	0.52
2:B:457:ARG:HD3	2:B:459:SER:O	2.10	0.52
1:H:1006:GLN:NE2	1:H:1085:TYR:O	2.42	0.52
2:B:497:PHE:CE2	2:B:507:PRO:HB3	2.45	0.52
1:H:6:GLN:HA	1:H:22:CYS:HA	1.93	0.51
1:H:106:ASP:O	1:H:108:TYR:N	2.43	0.51
1:H:50:ARG:HD2	1:H:111:TYR:CE1	2.46	0.51
2:B:520:ALA:HB1	2:B:521:PRO:HD2	1.92	0.50
2:B:485:GLY:N	2:B:488:CYS:O	2.42	0.50
1:H:47:TRP:CG	1:H:1097:VAL:HG22	2.46	0.50
1:H:112:TYR:HA	1:H:1033:TYR:OH	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:391:CYS:CA	2:B:525:CYS:HB3	2.34	0.50
1:H:23:LYS:HG3	1:H:78:THR:HG22	1.94	0.49
1:H:1027:LEU:HD23	1:H:1070:VAL:HG23	1.93	0.49
2:B:474:GLN:HA	2:B:480:CYS:SG	2.52	0.49
1:H:36:TRP:HB3	1:H:48:MET:HE2	1.94	0.49
2:B:381:GLY:HA3	2:B:430:THR:HA	1.94	0.48
1:H:71:THR:HG23	1:H:80:TYR:HB2	1.95	0.48
1:H:10:GLU:HB2	1:H:123:VAL:HG22	1.96	0.47
2:B:453:TYR:CE1	2:B:493:GLN:HB3	2.49	0.47
2:B:457:ARG:HH11	2:B:467:ASP:HB2	1.79	0.47
2:B:350:VAL:HA	2:B:400:PHE:HB2	1.97	0.46
1:H:1027:LEU:N	1:H:1028:PRO:HD2	2.31	0.46
1:H:36:TRP:CE2	1:H:81:MET:HB2	2.52	0.45
1:H:101:VAL:HG23	1:H:111:TYR:HA	1.98	0.45
1:H:1006:GLN:HG2	1:H:1103:THR:OG1	2.17	0.45
2:B:377:PHE:CD2	2:B:434:ILE:HG12	2.52	0.45
1:H:98:LYS:HB3	1:H:98:LYS:HE3	1.76	0.44
2:B:440:ASN:OD1	2:B:441:LEU:HG	2.18	0.44
2:B:353:TRP:NE1	2:B:466:ARG:HB2	2.33	0.43
2:B:410:ILE:O	2:B:425:LEU:HD13	2.18	0.43
1:H:1004:LEU:HG	1:H:1098:VAL:HG13	2.00	0.43
1:H:32:TYR:CD1	1:H:98:LYS:HD3	2.55	0.42
1:H:36:TRP:CH2	1:H:96:CYS:HB2	2.55	0.42
1:H:91:THR:HG23	1:H:124:THR:HA	2.02	0.41
2:B:384:PRO:HA	2:B:387:LEU:HB2	2.01	0.41
2:B:382:VAL:HG21	2:B:390:LEU:HG	2.01	0.41
1:H:114:MET:O	1:H:1035:TYR:HE1	2.04	0.41
2:B:445:VAL:HG22	2:B:499:PRO:HG2	2.03	0.41
2:B:497:PHE:CZ	2:B:507:PRO:HB3	2.54	0.41
1:H:35:SER:HA	1:H:50:ARG:HA	2.03	0.40

There are no symmetry-related clashes.

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	H	232/281 (83%)	224 (97%)	8 (3%)	0	100	100
2	B	190/213 (89%)	181 (95%)	9 (5%)	0	100	100
All	All	422/494 (85%)	405 (96%)	17 (4%)	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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	164/222 (74%)	164 (100%)	0	100	100
2	B	143/182 (79%)	142 (99%)	1 (1%)	76	79
All	All	307/404 (76%)	306 (100%)	1 (0%)	86	85

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	440	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	43	GLN
1	H	65	GLN
2	B	360	ASN
2	B	394	ASN
2	B	409	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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	TPO	B	345	2	8,10,11	1.22	0	10,14,16	1.24	1 (10%)

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	TPO	B	345	2	-	3/9/11/13	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	345	TPO	O3P-P-OG1	2.75	116.57	105.85

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	345	TPO	O-C-CA-CB
2	B	345	TPO	CB-OG1-P-O1P
2	B	345	TPO	CG2-CB-OG1-P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

5 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	A	1	3,2	14,14,15	0.25	0	17,19,21	1.81	4 (23%)
3	NAG	A	2	3	14,14,15	0.69	1 (7%)	17,19,21	0.87	0
3	BMA	A	3	3	11,11,12	0.70	1 (9%)	15,15,17	1.13	1 (6%)
3	FUC	A	4	3	10,10,11	1.35	1 (10%)	14,14,16	1.70	4 (28%)
3	FUC	A	5	3	10,10,11	0.69	0	14,14,16	0.90	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	A	1	3,2	-	1/6/23/26	0/1/1/1
3	NAG	A	2	3	-	0/6/23/26	0/1/1/1
3	BMA	A	3	3	-	2/2/19/22	0/1/1/1
3	FUC	A	4	3	-	-	0/1/1/1
3	FUC	A	5	3	-	-	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	4	FUC	C1-C2	3.39	1.60	1.52
3	A	2	NAG	C1-C2	2.28	1.55	1.52
3	A	3	BMA	C1-C2	2.20	1.57	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1	NAG	C1-O5-C5	3.94	117.47	112.19
3	A	1	NAG	O3-C3-C4	3.83	119.41	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	4	FUC	C1-O5-C5	3.00	120.04	112.97
3	A	4	FUC	C1-C2-C3	2.84	113.78	109.64
3	A	3	BMA	C1-O5-C5	2.78	115.91	112.19
3	A	4	FUC	O5-C1-C2	2.59	116.96	110.79
3	A	4	FUC	O5-C5-C4	2.52	114.08	109.55
3	A	1	NAG	C4-C3-C2	-2.33	107.60	111.02
3	A	1	NAG	C2-N2-C7	2.29	125.97	122.90

There are no chirality outliers.

All (3) torsion outliers are listed below:

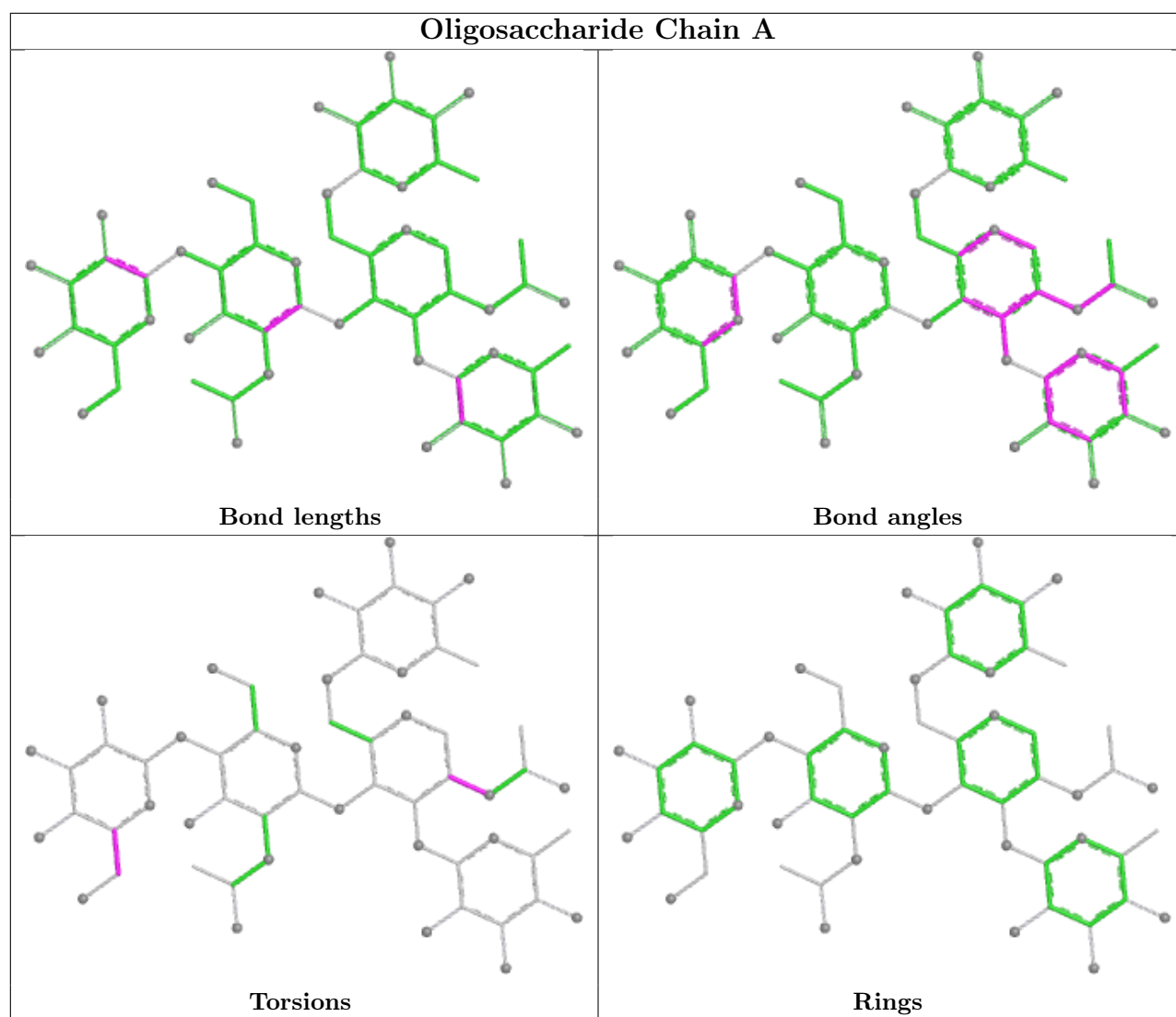
Mol	Chain	Res	Type	Atoms
3	A	3	BMA	O5-C5-C6-O6
3	A	3	BMA	C4-C5-C6-O6
3	A	1	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	4	FUC	1	0
3	A	2	NAG	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](#)

There are no ligands in this entry.

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	H	236/281 (83%)	0.62	9 (3%)	44	24	45, 61, 80, 96	1 (0%)
2	B	192/213 (90%)	0.74	12 (6%)	26	13	40, 56, 101, 121	0
All	All	428/494 (86%)	0.68	21 (4%)	35	18	40, 60, 94, 121	1 (0%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	489	TYR	4.2
1	H	18	VAL	3.2
1	H	26	GLY	3.1
1	H	122	THR	2.8
2	B	334	ASN	2.6
2	B	389	ASP	2.6
2	B	456	PHE	2.5
1	H	1106	THR	2.4
2	B	516	GLU	2.3
2	B	453	TYR	2.3
1	H	1089	SER	2.3
1	H	22	CYS	2.2
2	B	449	TYR	2.2
1	H	27	GLY	2.1
2	B	388	ASN	2.1
2	B	403	ARG	2.1
1	H	1005	THR	2.1
1	H	1082	GLU	2.1
2	B	368	LEU	2.0
2	B	360	ASN	2.0
2	B	495	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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.

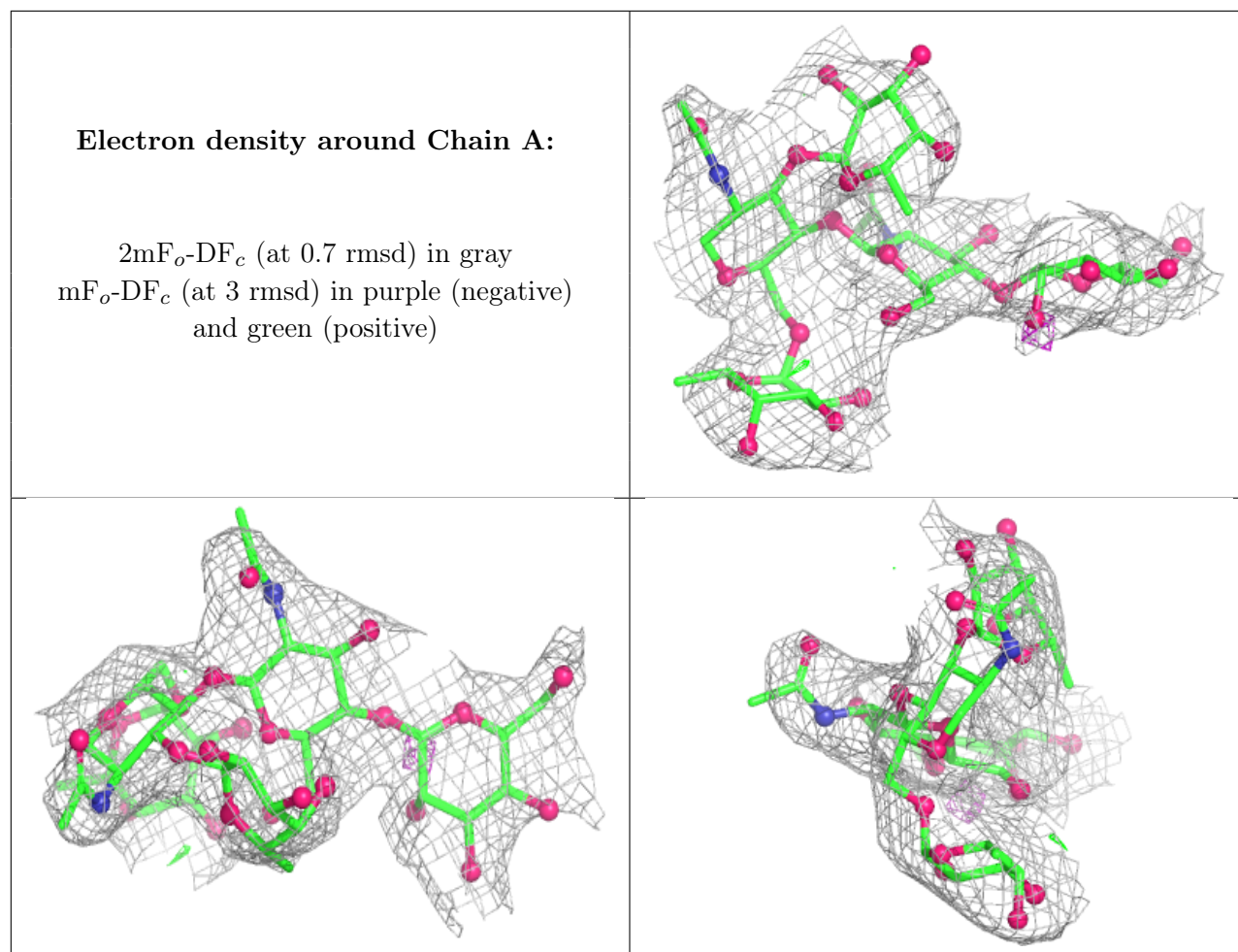
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	TPO	B	345	11/12	0.80	0.16	57,57,57,57	0

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	NAG	A	1	14/15	-	-	76,76,76,76	0
3	NAG	A	2	14/15	-	-	85,85,85,85	0
3	BMA	A	3	11/12	-	-	98,98,98,98	0
3	FUC	A	4	10/11	-	-	86,86,86,86	0
3	FUC	A	5	10/11	-	-	77,77,77,77	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 [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.