



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 01:44 PM BST

PDB ID : 4XE4
Title : Coagulation Factor XII protease domain crystal structure
Authors : Pathak, M.; Wilmann, P.; Awford, J.; Li, C.; Fisher, P.M.; Dreveny, I.; Dekker, L.V.; Emsley, J.
Deposited on : 2014-12-22
Resolution : 2.40 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 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

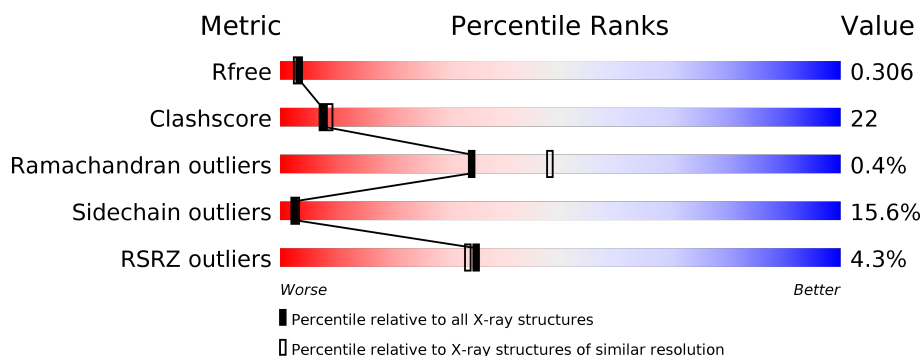
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	255	
2	B	3	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1894 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 Coagulation factor XII.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	231	1768	1113	309	334	12	7	1	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	467	SER	CYS	engineered mutation	UNP P00748
A	597	HIS	-	expression tag	UNP P00748
A	598	HIS	-	expression tag	UNP P00748
A	599	THR	-	expression tag	UNP P00748
A	600	GLY	-	expression tag	UNP P00748
A	601	THR	-	expression tag	UNP P00748
A	602	ARG	-	expression tag	UNP P00748
A	603	HIS	-	expression tag	UNP P00748
A	604	HIS	-	expression tag	UNP P00748
A	605	HIS	-	expression tag	UNP P00748
A	606	HIS	-	expression tag	UNP P00748
A	607	HIS	-	expression tag	UNP P00748
A	608	HIS	-	expression tag	UNP P00748

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(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
			Total	C	N	O			
2	B	3	42	24	3	15	0	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

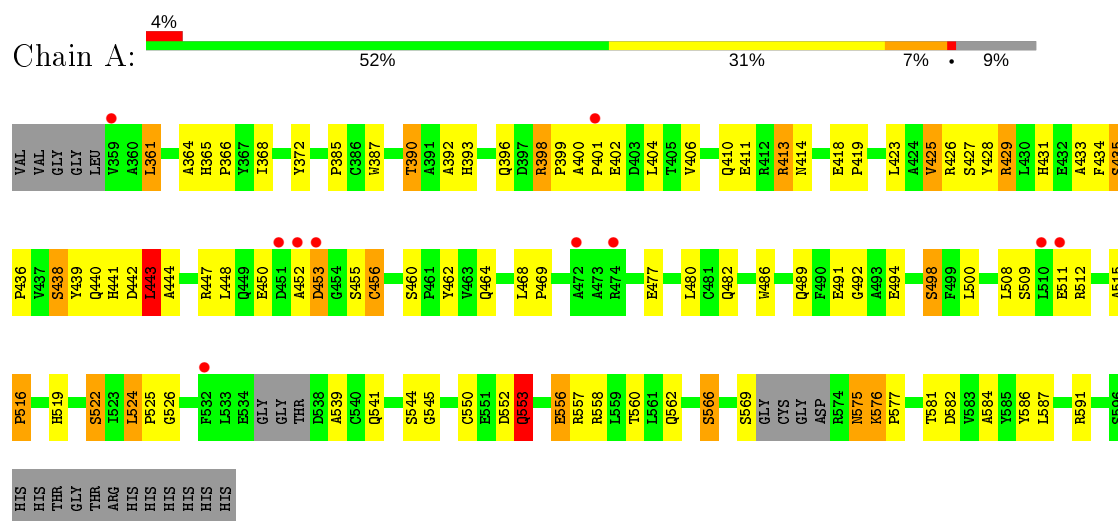
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	78	Total	O	0	0
			78	78		

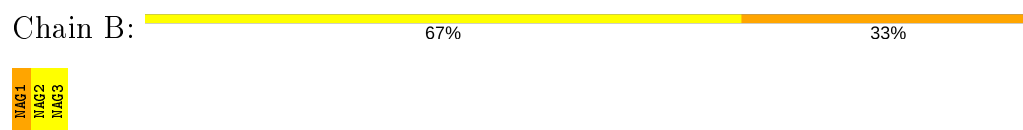
3 Residue-property plots [i](#)

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: Coagulation factor XII



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	137.06 Å 137.06 Å 37.01 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.57 – 2.40 39.57 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.2 (39.57-2.40) 98.2 (39.57-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.39 Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.218 , 0.295 0.232 , 0.306	Depositor DCC
R_{free} test set	723 reflections (4.66%)	wwPDB-VP
Wilson B-factor (Å ²)	61.7	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.052 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	1894	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.29% 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: GOL, NAG

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	0.81	0/1817	1.06	9/2478 (0.4%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	539	ALA	CB-CA-C	-8.76	96.97	110.10
1	A	515	ALA	CB-CA-C	-8.11	97.94	110.10
1	A	443	LEU	CA-CB-CG	7.24	131.96	115.30
1	A	429	ARG	CB-CA-C	6.66	123.72	110.40
1	A	576	LYS	CB-CA-C	6.05	122.51	110.40
1	A	456	CYS	N-CA-C	6.02	127.25	111.00
1	A	556	GLU	CB-CA-C	5.90	122.19	110.40
1	A	553	GLN	CB-CA-C	-5.85	98.69	110.40
1	A	516	PRO	N-CA-CB	-5.08	97.01	102.60

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	1768	0	1679	78	0
2	B	42	0	37	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	6	0	8	0	0
4	A	78	0	0	13	0
All	All	1894	0	1724	78	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 22.

All (78) 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:423:LEU:HD11	1:A:456:CYS:O	1.44	1.16
1:A:460:SER:O	1:A:464:GLN:NE2	1.91	1.02
1:A:423:LEU:CD1	1:A:456:CYS:O	2.20	0.90
1:A:399:PRO:O	1:A:428:TYR:OH	1.93	0.84
1:A:439:TYR:HA	4:A:878:HOH:O	1.78	0.83
1:A:431:HIS:CD2	1:A:441:HIS:HB3	2.24	0.72
1:A:489:GLN:OE1	1:A:498:SER:N	2.23	0.72
1:A:398:ARG:N	1:A:399:PRO:HD3	2.05	0.72
1:A:441:HIS:CE1	4:A:875:HOH:O	2.43	0.70
1:A:434:PHE:O	1:A:436:PRO:HD3	1.92	0.70
1:A:522:SER:O	4:A:876:HOH:O	2.11	0.68
1:A:516:PRO:HD3	4:A:816:HOH:O	1.95	0.67
1:A:440:GLN:HE21	1:A:524:LEU:HB3	1.64	0.63
1:A:519:HIS:CD2	1:A:577:PRO:HG2	2.35	0.61
1:A:361:LEU:HD12	4:A:803:HOH:O	2.00	0.61
1:A:439:TYR:C	4:A:878:HOH:O	2.39	0.60
1:A:584:ALA:O	1:A:587:LEU:HB2	2.01	0.60
1:A:550:CYS:SG	1:A:562:GLN:HG3	2.42	0.60
1:A:418:GLU:HB2	1:A:419:PRO:HD3	1.83	0.59
1:A:404:LEU:HB2	1:A:425:VAL:CG1	2.32	0.59
1:A:440:GLN:HG3	1:A:524:LEU:HD23	1.84	0.59
1:A:550:CYS:HB2	1:A:562:GLN:HG3	1.85	0.57
1:A:400:ALA:HB1	1:A:402:GLU:OE1	2.06	0.56
1:A:414:ASN:HA	2:B:1:NAG:O7	2.07	0.55
1:A:440:GLN:N	4:A:878:HOH:O	2.40	0.55
1:A:433:ALA:O	1:A:441:HIS:HD2	1.90	0.55
1:A:526:GLY:C	4:A:870:HOH:O	2.44	0.55
1:A:429:ARG:O	1:A:444:ALA:HB1	2.08	0.53
1:A:398:ARG:N	1:A:399:PRO:CD	2.71	0.53
1:A:440:GLN:OE1	1:A:441:HIS:NE2	2.41	0.53
1:A:519:HIS:HD2	1:A:577:PRO:HG2	1.76	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:525:PRO:O	4:A:875:HOH:O	2.19	0.50
1:A:393:HIS:HA	1:A:396:GLN:HG2	1.93	0.50
1:A:587:LEU:O	1:A:591:ARG:HG3	2.12	0.50
1:A:491:GLU:CD	1:A:492:GLY:N	2.66	0.49
1:A:575:ASN:C	1:A:575:ASN:HD22	2.17	0.49
1:A:477:GLU:OE2	1:A:509:SER:OG	2.17	0.48
1:A:469:PRO:HA	1:A:560:THR:OG1	2.14	0.47
1:A:393:HIS:HE1	1:A:544:SER:HB3	1.78	0.47
1:A:435:SER:HG	1:A:438:SER:H	1.63	0.47
1:A:428:TYR:O	1:A:429:ARG:HB2	2.15	0.47
1:A:406:VAL:HG11	1:A:448:LEU:CD2	2.44	0.47
1:A:450:GLU:OE1	1:A:450:GLU:N	2.44	0.47
1:A:477:GLU:OE2	1:A:509:SER:N	2.41	0.47
1:A:491:GLU:OE2	1:A:492:GLY:N	2.48	0.46
1:A:447:ARG:NH1	1:A:448:LEU:O	2.49	0.46
1:A:550:CYS:CB	1:A:562:GLN:HG3	2.45	0.46
1:A:434:PHE:HE1	4:A:878:HOH:O	1.93	0.46
1:A:365:HIS:O	1:A:368:ILE:HG22	2.16	0.46
1:A:434:PHE:CE1	4:A:878:HOH:O	2.56	0.45
1:A:426:ARG:O	1:A:427:SER:HB2	2.16	0.45
1:A:443:LEU:HD21	1:A:586:TYR:CB	2.47	0.44
1:A:516:PRO:CD	4:A:816:HOH:O	2.57	0.44
1:A:404:LEU:CB	1:A:425:VAL:CG1	2.96	0.44
1:A:576:LYS:HA	1:A:576:LYS:HD2	1.67	0.44
1:A:413:ARG:HG3	1:A:486:TRP:HB3	2.00	0.44
1:A:569:SER:O	1:A:569:SER:OG	2.36	0.44
1:A:418:GLU:N	1:A:419:PRO:HD2	2.34	0.43
1:A:582:ASP:C	1:A:582:ASP:OD1	2.56	0.43
1:A:385:PRO:O	1:A:456:CYS:HB2	2.19	0.43
1:A:404:LEU:HB2	1:A:425:VAL:HG13	2.00	0.43
1:A:372:TYR:OH	1:A:410:GLN:NE2	2.51	0.42
1:A:468:LEU:O	1:A:560:THR:OG1	2.38	0.42
1:A:392:ALA:N	1:A:442:ASP:OD1	2.51	0.42
1:A:440:GLN:HG3	1:A:524:LEU:CB	2.49	0.42
1:A:452:ALA:O	1:A:453:ASP:CB	2.68	0.42
1:A:468:LEU:HB3	1:A:469:PRO:CD	2.50	0.42
1:A:364:ALA:O	1:A:365:HIS:C	2.58	0.42
1:A:390:THR:HG22	1:A:545:GLY:HA3	2.02	0.41
1:A:435:SER:OG	1:A:438:SER:HB3	2.20	0.41
1:A:418:GLU:N	1:A:419:PRO:CD	2.83	0.41
1:A:434:PHE:CD1	4:A:878:HOH:O	2.73	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:552:ASP:O	1:A:553:GLN:HB2	2.21	0.41
1:A:442:ASP:OD2	1:A:566:SER:OG	2.34	0.41
1:A:442:ASP:O	1:A:581:THR:OG1	2.26	0.40
1:A:387:TRP:CH2	1:A:447:ARG:HD3	2.57	0.40
1:A:433:ALA:O	1:A:441:HIS:CD2	2.73	0.40
1:A:462:TYR:N	1:A:462:TYR:CD1	2.89	0.40

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	226/255 (89%)	187 (83%)	38 (17%)	1 (0%)	34	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	401	PRO

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	187/203 (92%)	158 (84%)	29 (16%)	2	3

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	361	LEU
1	A	366	PRO
1	A	390	THR
1	A	398	ARG
1	A	411	GLU
1	A	413	ARG
1	A	425	VAL
1	A	435	SER
1	A	438	SER
1	A	443	LEU
1	A	453	ASP
1	A	455	SER
1	A	480	LEU
1	A	482	GLN
1	A	494	GLU
1	A	498	SER
1	A	500	LEU
1	A	508	LEU
1	A	511	GLU
1	A	512	ARG
1	A	522	SER
1	A	524	LEU
1	A	541	GLN
1	A	553	GLN
1	A	556	GLU
1	A	557	ARG
1	A	558	ARG
1	A	566	SER
1	A	575	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	393	HIS
1	A	410	GLN
1	A	482	GLN
1	A	501	GLN
1	A	541	GLN
1	A	575	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 ⓘ

3 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	B	1	1,2	14,14,15	0.80	0	17,19,21	2.59	7 (41%)
2	NAG	B	2	2	14,14,15	0.93	1 (7%)	17,19,21	2.39	4 (23%)
2	NAG	B	3	2	14,14,15	0.78	0	17,19,21	2.46	8 (47%)

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	B	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	NAG	B	3	2	-	4/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	NAG	C1-C2	2.23	1.55	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	O4-C4-C3	6.22	124.73	110.35
2	B	1	NAG	O4-C4-C5	5.68	123.40	109.30
2	B	3	NAG	C1-O5-C5	5.06	119.05	112.19
2	B	1	NAG	C1-C2-N2	4.80	118.69	110.49
2	B	1	NAG	C2-N2-C7	4.44	129.22	122.90
2	B	2	NAG	C1-C2-N2	4.35	117.93	110.49
2	B	2	NAG	O5-C1-C2	-4.06	104.87	111.29
2	B	3	NAG	C1-C2-N2	3.82	117.00	110.49
2	B	3	NAG	C2-N2-C7	3.73	128.22	122.90
2	B	3	NAG	C8-C7-N2	3.50	122.02	116.10
2	B	3	NAG	C3-C4-C5	3.45	116.39	110.24
2	B	1	NAG	O3-C3-C4	3.41	118.24	110.35
2	B	3	NAG	O5-C1-C2	-3.30	106.08	111.29
2	B	1	NAG	C1-O5-C5	2.56	115.66	112.19
2	B	1	NAG	O5-C1-C2	-2.48	107.37	111.29
2	B	3	NAG	O7-C7-C8	-2.40	117.59	122.06
2	B	2	NAG	O7-C7-C8	-2.31	117.76	122.06
2	B	3	NAG	O5-C5-C4	2.23	116.26	110.83
2	B	1	NAG	C6-C5-C4	2.08	117.86	113.00

There are no chirality outliers.

All (5) torsion outliers are listed below:

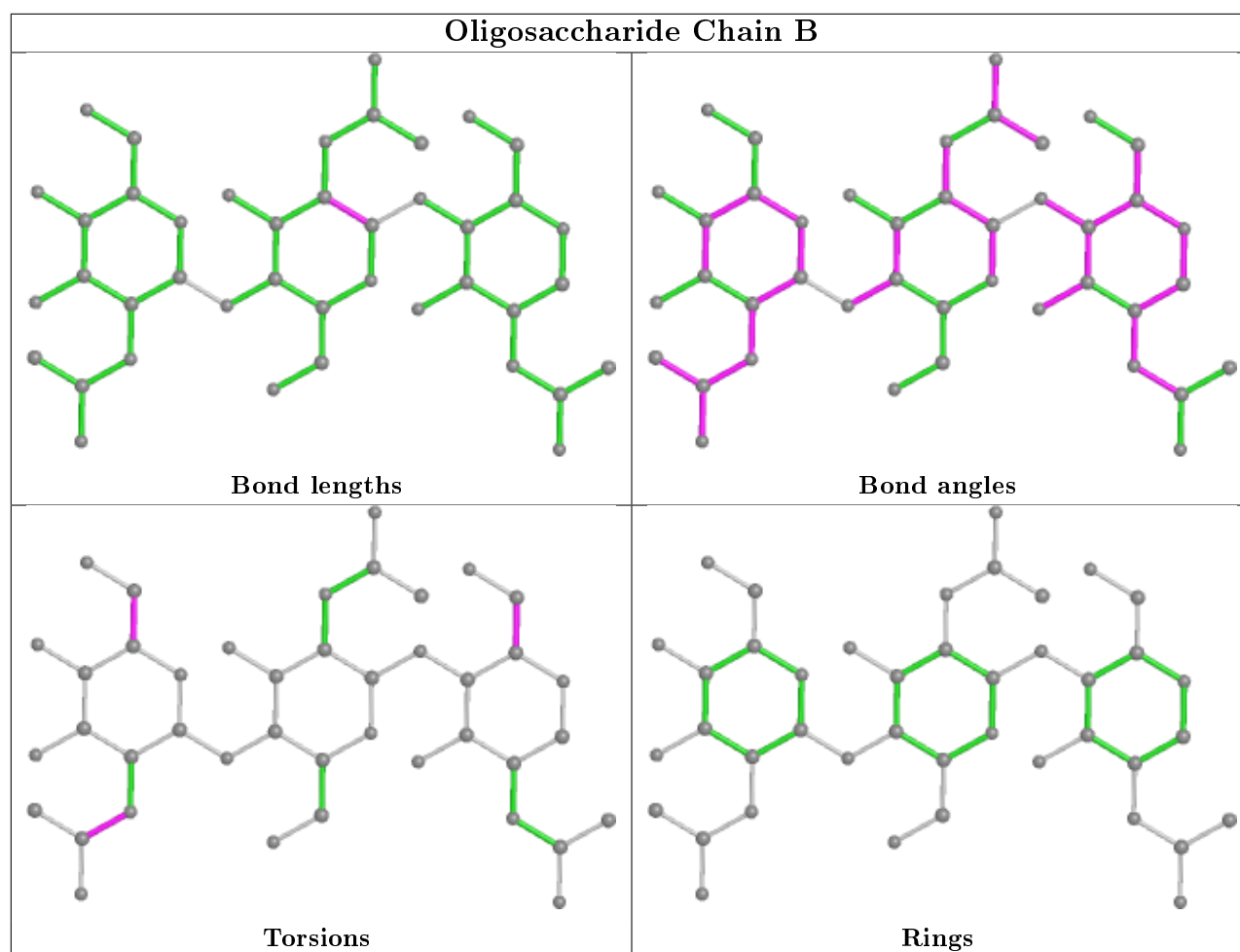
Mol	Chain	Res	Type	Atoms
2	B	3	NAG	C4-C5-C6-O6
2	B	3	NAG	C8-C7-N2-C2
2	B	3	NAG	O7-C7-N2-C2
2	B	3	NAG	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	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](#)

1 ligand 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	$\# Z > 2$
3	GOL	A	704	-	5,5,5	0.23	0	5,5,5	0.50	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	GOL	A	704	-	-	0/4/4/4	-

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, 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	231/255 (90%)	0.16	10 (4%) 35 33	38, 67, 125, 141	7 (3%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	359	VAL	4.0
1	A	452	ALA	3.5
1	A	472	ALA	3.5
1	A	532	PHE	3.1
1	A	451	ASP	2.3
1	A	453	ASP	2.3
1	A	474	ARG	2.3
1	A	510	LEU	2.3
1	A	511	GLU	2.1
1	A	401	PRO	2.1

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.

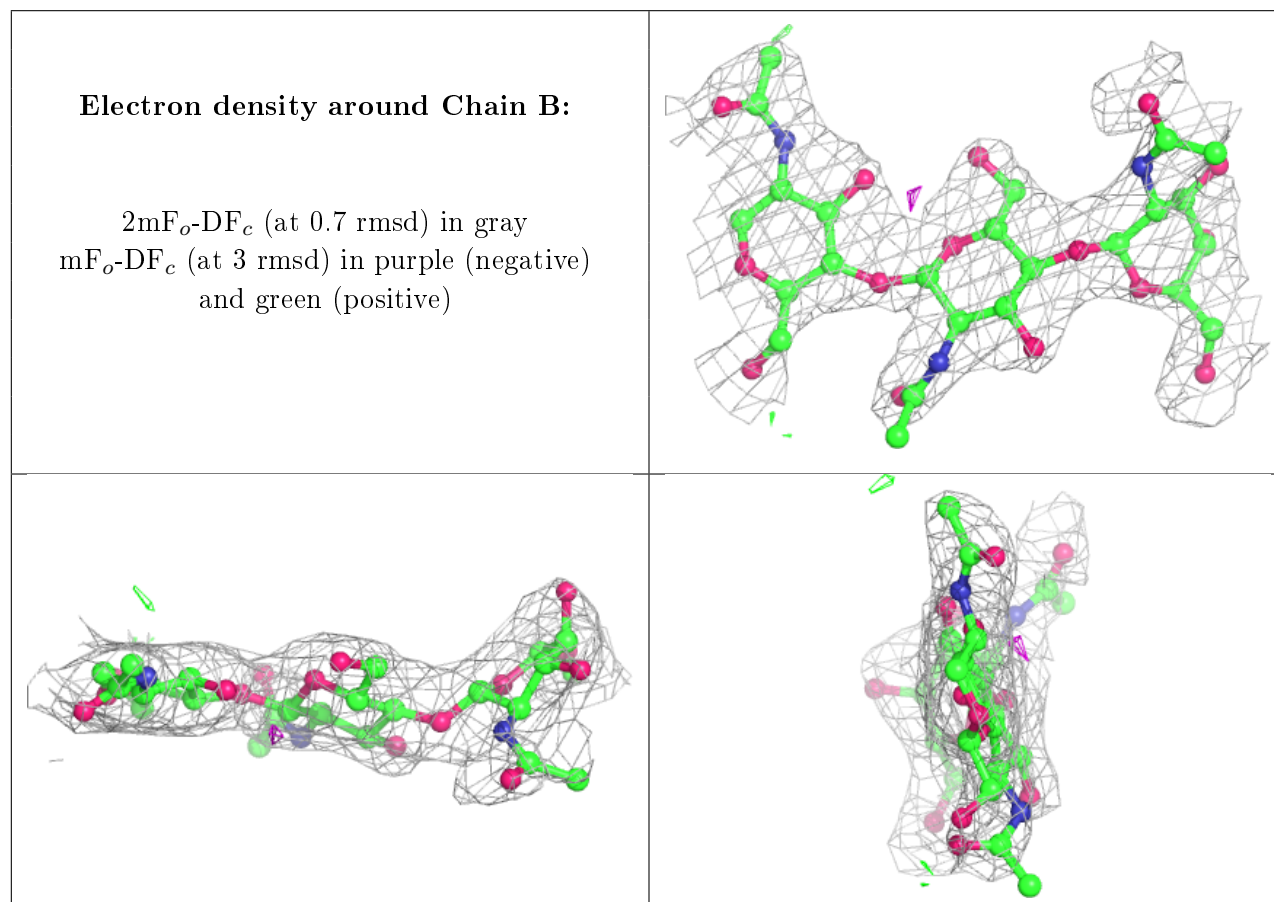
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	B	3	14/15	0.71	0.31	109,134,138,138	0
2	NAG	B	1	14/15	0.88	0.13	71,80,94,103	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	B	2	14/15	0.88	0.28	97,108,116,120	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](#)

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(\AA^2)	Q<0.9
3	GOL	A	704	6/6	0.92	0.17	66,80,82,96	0

6.5 Other polymers [i](#)

There are no such residues in this entry.