



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:43 PM UTC

PDB ID : 6ATT / pdb_00006att
Title : 39S Fab bound to HER2 ecd
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Deposited on : 2017-08-29
Resolution : 3.77 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

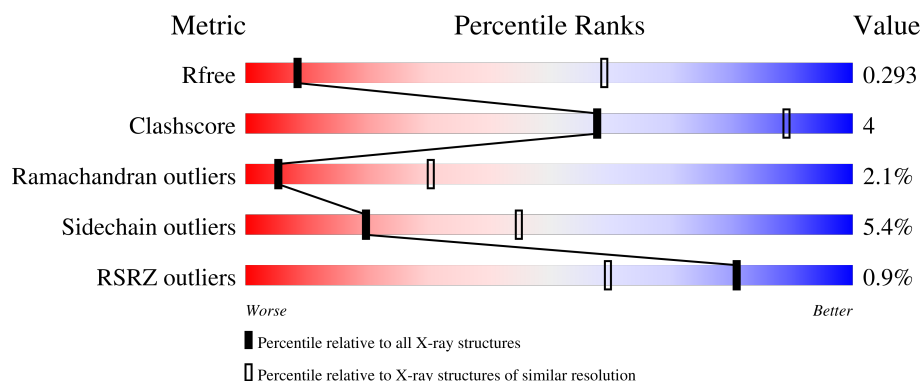
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.77 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	1085 (3.94-3.62)
Clashscore	190562	1029 (3.92-3.64)
Ramachandran outliers	187476	1064 (3.94-3.62)
Sidechain outliers	187428	1059 (3.94-3.62)
RSRZ outliers	180081	1084 (3.94-3.62)

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	630	80% 9% 10%
2	H	224	84% 13%
3	L	220	90% 10%
4	B	2	100%
4	C	2	100%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7819 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 Receptor tyrosine-protein kinase erbB-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	567	Total	C	N	O	S	0	0	0
			4370	2717	784	820	49			

- Molecule 2 is a protein called Antibody 39S Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	224	Total	C	N	O	S	0	0	0
			1679	1060	278	334	7			

- Molecule 3 is a protein called Antibody 39S Fab light chain.

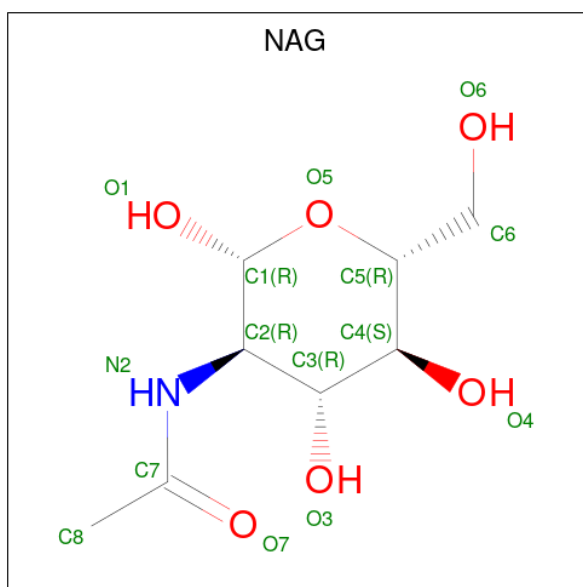
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	220	Total	C	N	O	S	0	0	0
			1700	1072	284	338	6			

- Molecule 4 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
4	B	2	Total	C	N	O		0	0	0
			28	16	2	10				
4	C	2	Total	C	N	O		0	0	0
			28	16	2	10				

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

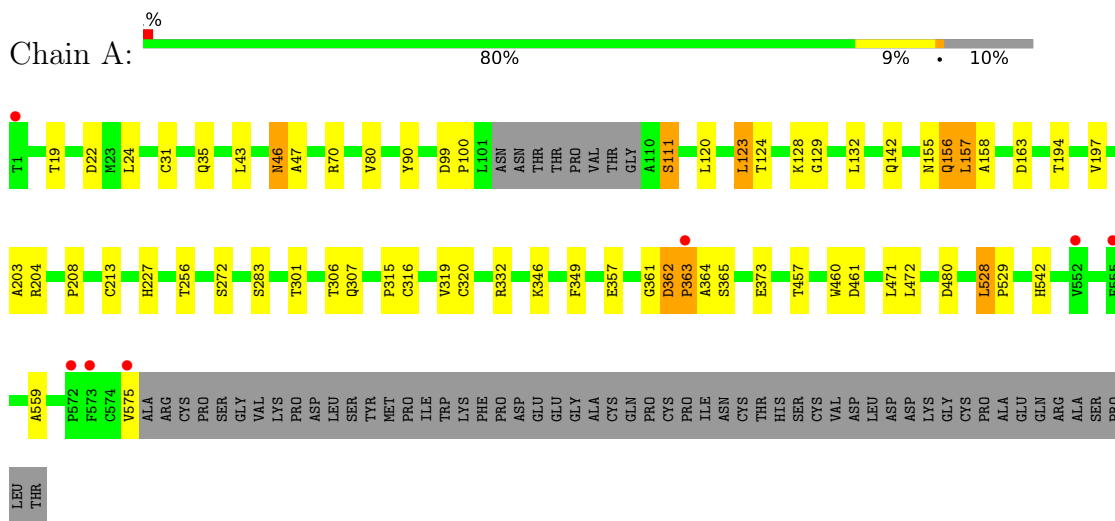


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	A	1	14	8	1	5	0	0

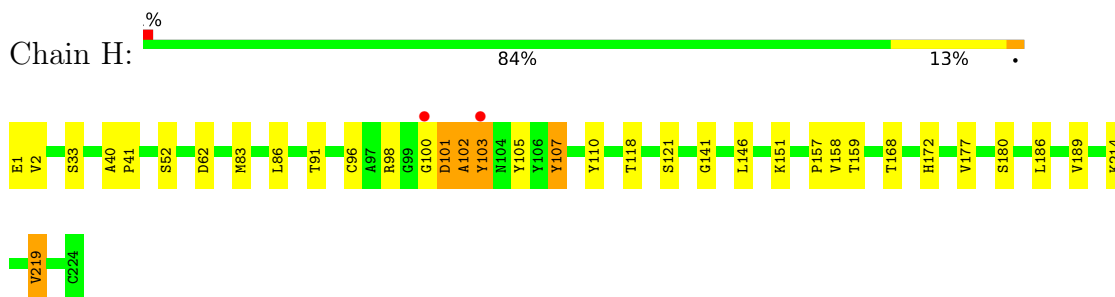
3 Residue-property plots

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.

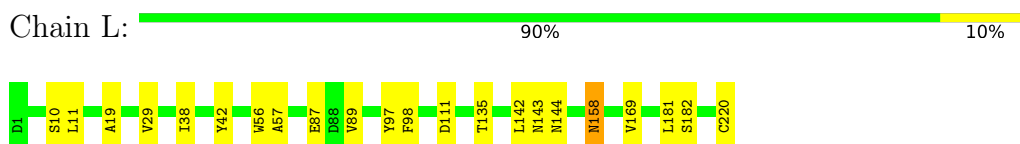
- Molecule 1: Receptor tyrosine-protein kinase erbB-2



- Molecule 2: Antibody 39S Fab heavy chain



- Molecule 3: Antibody 39S Fab light chain



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



MAG1
MAG2

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

Chain C:

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	67.33Å 108.17Å 348.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 3.77 40.00 – 3.77	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-3.77) 99.9 (40.00-3.77)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 3.78Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.251 , 0.289 0.265 , 0.293	Depositor DCC
R_{free} test set	1313 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	132.4	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 18.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7819	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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: 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	0.72	0/4472	0.95	2/6085 (0.0%)
2	H	0.70	0/1721	0.91	3/2342 (0.1%)
3	L	0.64	0/1741	0.82	0/2365
All	All	0.70	0/7934	0.91	5/10792 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	103	TYR	CB-CA-C	-6.51	103.09	111.50
2	H	107	TYR	CB-CA-C	-5.49	101.68	110.79
1	A	111	SER	CA-C-N	-5.25	115.17	120.31
1	A	111	SER	C-N-CA	-5.25	115.17	120.31
2	H	103	TYR	N-CA-C	5.11	119.17	111.81

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	4370	0	4186	44	0
2	H	1679	0	1626	31	0
3	L	1700	0	1652	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	28	0	25	3	0
4	C	28	0	25	0	0
5	A	14	0	13	0	0
All	All	7819	0	7527	69	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 4.

All (69) 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:129:GLY:H	1:A:155:ASN:ND2	1.54	1.03
1:A:197:VAL:HG13	2:H:103:TYR:CE1	1.97	0.99
1:A:362:ASP:H	1:A:363:PRO:CD	1.75	0.97
1:A:155:ASN:O	1:A:157:LEU:N	1.98	0.96
2:H:101:ASP:O	2:H:102:ALA:O	1.89	0.91
1:A:362:ASP:N	1:A:363:PRO:CD	2.34	0.88
2:H:33:SER:HB2	2:H:102:ALA:O	1.72	0.86
1:A:363:PRO:O	1:A:365:SER:N	2.12	0.82
4:B:1:NAG:H61	4:B:2:NAG:C7	2.09	0.82
1:A:363:PRO:C	1:A:365:SER:H	1.88	0.81
1:A:46:ASN:OD1	4:B:1:NAG:O7	1.99	0.81
1:A:197:VAL:CG1	2:H:103:TYR:CE1	2.65	0.80
3:L:11:LEU:HD21	3:L:19:ALA:HB1	1.64	0.79
1:A:129:GLY:H	1:A:155:ASN:HD21	1.30	0.76
1:A:129:GLY:N	1:A:155:ASN:ND2	2.31	0.75
1:A:197:VAL:CG1	2:H:103:TYR:HE1	1.98	0.75
1:A:362:ASP:H	1:A:363:PRO:HD3	1.50	0.71
1:A:197:VAL:HG13	2:H:103:TYR:CZ	2.26	0.71
2:H:107:TYR:O	3:L:42:TYR:OH	2.10	0.70
1:A:197:VAL:HG13	2:H:103:TYR:HE1	1.54	0.69
1:A:129:GLY:H	1:A:155:ASN:HD22	1.40	0.68
1:A:542:HIS:CD2	1:A:559:ALA:HB2	2.29	0.67
1:A:362:ASP:N	1:A:363:PRO:HD2	2.10	0.64
1:A:208:PRO:CB	2:H:103:TYR:HD1	2.11	0.63
2:H:103:TYR:O	2:H:103:TYR:CD1	2.52	0.63
2:H:101:ASP:O	2:H:102:ALA:C	2.37	0.62
1:A:129:GLY:N	1:A:155:ASN:HD22	1.95	0.62
1:A:363:PRO:C	1:A:365:SER:N	2.51	0.62
1:A:208:PRO:HB2	2:H:103:TYR:HD1	1.66	0.60
1:A:120:LEU:HD13	1:A:123:LEU:HD22	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:52:SER:HB2	2:H:103:TYR:HB3	1.86	0.57
1:A:362:ASP:OD1	1:A:362:ASP:O	2.23	0.56
1:A:197:VAL:HG12	2:H:103:TYR:HE1	1.71	0.55
1:A:208:PRO:HB3	2:H:103:TYR:CD1	2.43	0.54
1:A:208:PRO:CB	2:H:103:TYR:CD1	2.91	0.54
1:A:301:THR:HB	1:A:307:GLN:HE21	1.73	0.53
2:H:83:MET:HB3	2:H:86:LEU:HD21	1.91	0.51
1:A:362:ASP:O	1:A:363:PRO:O	2.30	0.50
2:H:103:TYR:HD1	2:H:103:TYR:O	1.93	0.50
1:A:319:VAL:HG11	1:A:349:PHE:HE1	1.76	0.50
1:A:203:ALA:O	1:A:204:ARG:HG2	2.11	0.50
2:H:105:TYR:CZ	3:L:56:TRP:HH2	2.30	0.50
1:A:19:THR:HG23	1:A:472:LEU:HB3	1.95	0.48
1:A:24:LEU:CD1	1:A:43:LEU:HD21	2.43	0.48
1:A:19:THR:CG2	1:A:472:LEU:HD13	2.44	0.47
2:H:2:VAL:HG21	2:H:110:TYR:CE2	2.49	0.47
2:H:146:LEU:HD11	2:H:219:VAL:CG1	2.44	0.47
3:L:29:VAL:HG13	3:L:98:PHE:HB2	1.97	0.47
1:A:128:LYS:HA	1:A:155:ASN:HD22	1.79	0.47
1:A:197:VAL:HG13	2:H:103:TYR:OH	2.16	0.46
2:H:91:THR:HG23	2:H:118:THR:HA	1.97	0.46
1:A:361:GLY:HA2	1:A:363:PRO:HD3	1.98	0.46
1:A:208:PRO:HB2	2:H:103:TYR:O	2.15	0.46
2:H:100:GLY:C	2:H:101:ASP:O	2.58	0.46
1:A:460:TRP:CE2	1:A:471:LEU:HD22	2.50	0.46
2:H:100:GLY:O	2:H:101:ASP:O	2.35	0.45
1:A:357:GLU:O	1:A:361:GLY:N	2.49	0.45
1:A:90:TYR:CD1	1:A:132:LEU:HB2	2.54	0.43
2:H:33:SER:CB	2:H:102:ALA:O	2.56	0.43
2:H:52:SER:CB	2:H:103:TYR:HB3	2.48	0.43
2:H:40:ALA:HB1	2:H:41:PRO:HD2	2.01	0.42
1:A:362:ASP:O	1:A:365:SER:HA	2.19	0.42
2:H:172:HIS:CD2	3:L:143:ASN:HD21	2.37	0.41
3:L:169:VAL:HG22	3:L:181:LEU:HD13	2.02	0.41
1:A:319:VAL:HG11	1:A:349:PHE:CE1	2.55	0.41
1:A:528:LEU:CB	1:A:529:PRO:CD	2.99	0.41
2:H:158:VAL:CG2	2:H:186:LEU:HD22	2.51	0.41
3:L:38:ILE:HG23	3:L:97:TYR:CE2	2.55	0.41
4:B:1:NAG:H61	4:B:2:NAG:C8	2.49	0.41

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	A	563/630 (89%)	503 (89%)	47 (8%)	13 (2%)	5	29
2	H	222/224 (99%)	205 (92%)	13 (6%)	4 (2%)	6	32
3	L	218/220 (99%)	205 (94%)	9 (4%)	4 (2%)	6	32
All	All	1003/1074 (93%)	913 (91%)	69 (7%)	21 (2%)	5	30

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	PRO
1	A	111	SER
1	A	156	GLN
1	A	362	ASP
1	A	363	PRO
1	A	364	ALA
1	A	528	LEU
2	H	102	ALA
1	A	47	ALA
1	A	123	LEU
1	A	158	ALA
1	A	320	CYS
2	H	141	GLY
2	H	101	ASP
3	L	144	ASN
3	L	57	ALA
3	L	158	ASN
1	A	315	PRO
3	L	89	VAL
1	A	99	ASP
2	H	157	PRO

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	488/543 (90%)	462 (95%)	26 (5%)	20	45
2	H	188/188 (100%)	175 (93%)	13 (7%)	14	40
3	L	195/195 (100%)	187 (96%)	8 (4%)	27	51
All	All	871/926 (94%)	824 (95%)	47 (5%)	20	45

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	31	CYS
1	A	35	GLN
1	A	46	ASN
1	A	70	ARG
1	A	80	VAL
1	A	124	THR
1	A	142	GLN
1	A	156	GLN
1	A	157	LEU
1	A	163	ASP
1	A	194	THR
1	A	213	CYS
1	A	227	HIS
1	A	256	THR
1	A	272	SER
1	A	283	SER
1	A	306	THR
1	A	316	CYS
1	A	332	ARG
1	A	346	LYS
1	A	373	GLU
1	A	457	THR
1	A	461	ASP
1	A	480	ASP
1	A	575	VAL

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Mol	Chain	Res	Type
2	H	1	GLU
2	H	62	ASP
2	H	96	CYS
2	H	98	ARG
2	H	121	SER
2	H	151	LYS
2	H	159	THR
2	H	168	THR
2	H	177	VAL
2	H	180	SER
2	H	189	VAL
2	H	214	LYS
2	H	219	VAL
3	L	10	SER
3	L	87	GLU
3	L	111	ASP
3	L	135	THR
3	L	142	LEU
3	L	158	ASN
3	L	182	SER
3	L	220	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	142	GLN
1	A	152	HIS
1	A	154	ASN
1	A	155	ASN
1	A	171	HIS
1	A	217	GLN
1	A	235	HIS
1	A	245	HIS
1	A	307	GLN
1	A	415	HIS
1	A	490	HIS
1	A	542	HIS
1	A	561	GLN
2	H	39	GLN
2	H	77	ASN
2	H	82	GLN

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Mol	Chain	Res	Type
2	H	84	ASN
3	L	35	ASN
3	L	44	GLN
3	L	143	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 ⓘ

4 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	$\# Z > 2$
4	NAG	B	1	1,4	14,14,15	0.53	0	17,19,21	1.50	2 (11%)
4	NAG	B	2	4	14,14,15	0.95	1 (7%)	17,19,21	1.27	1 (5%)
4	NAG	C	1	1,4	14,14,15	0.55	0	17,19,21	1.31	3 (17%)
4	NAG	C	2	4	14,14,15	0.62	0	17,19,21	1.08	1 (5%)

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	C	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	C	2	4	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	2	NAG	C1-C2	2.06	1.55	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1	NAG	C1-C2-N2	-4.21	103.81	110.43
4	B	1	NAG	C4-C3-C2	3.24	115.76	111.02
4	C	1	NAG	C4-C3-C2	2.88	115.23	111.02
4	B	2	NAG	C2-N2-C7	2.56	126.33	122.90
4	C	1	NAG	C1-O5-C5	2.30	115.27	112.19
4	C	1	NAG	O5-C1-C2	-2.20	107.89	111.29
4	C	2	NAG	O5-C1-C2	-2.13	108.00	111.29

There are no chirality outliers.

All (2) torsion outliers are listed below:

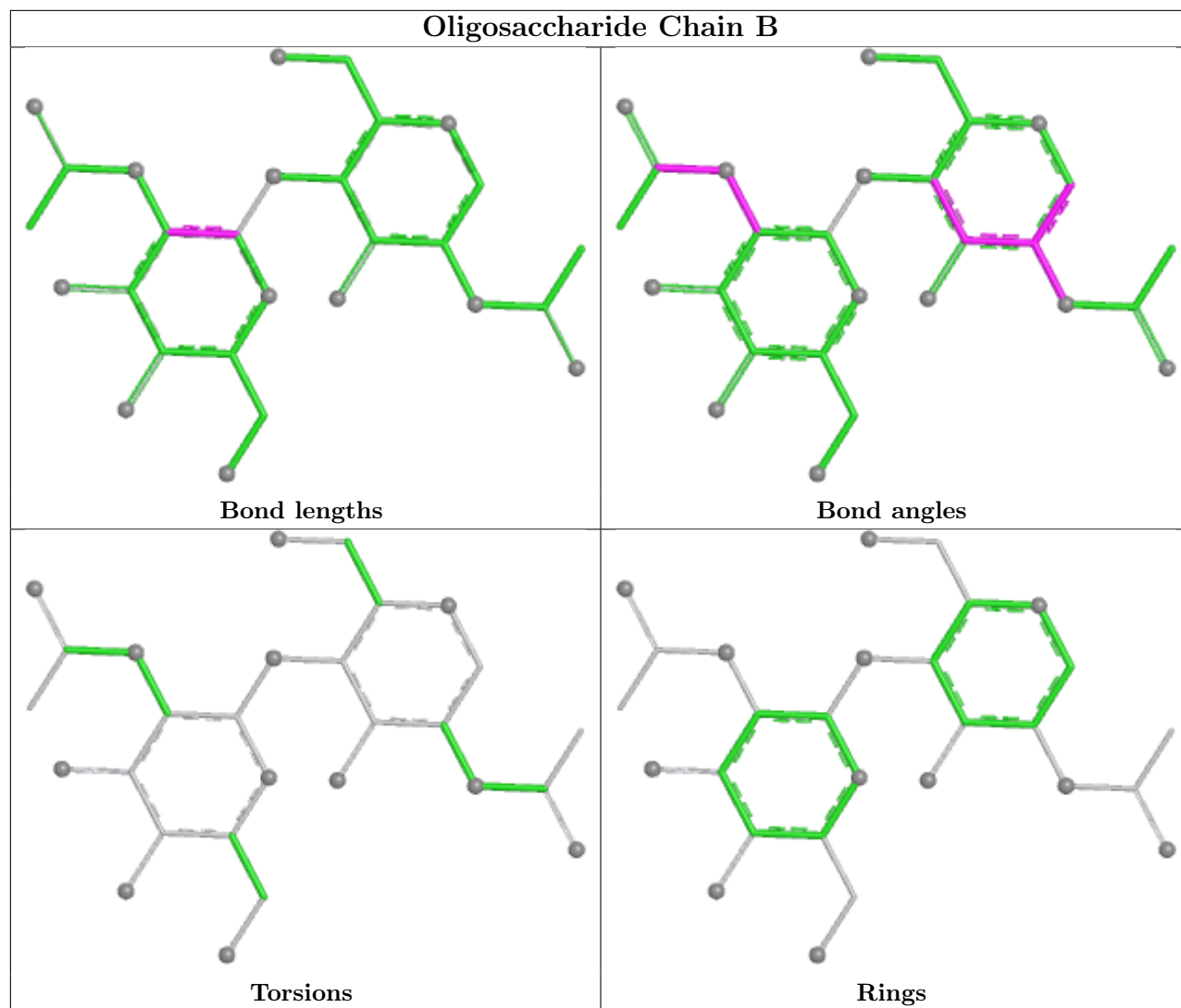
Mol	Chain	Res	Type	Atoms
4	C	2	NAG	C4-C5-C6-O6
4	C	2	NAG	O5-C5-C6-O6

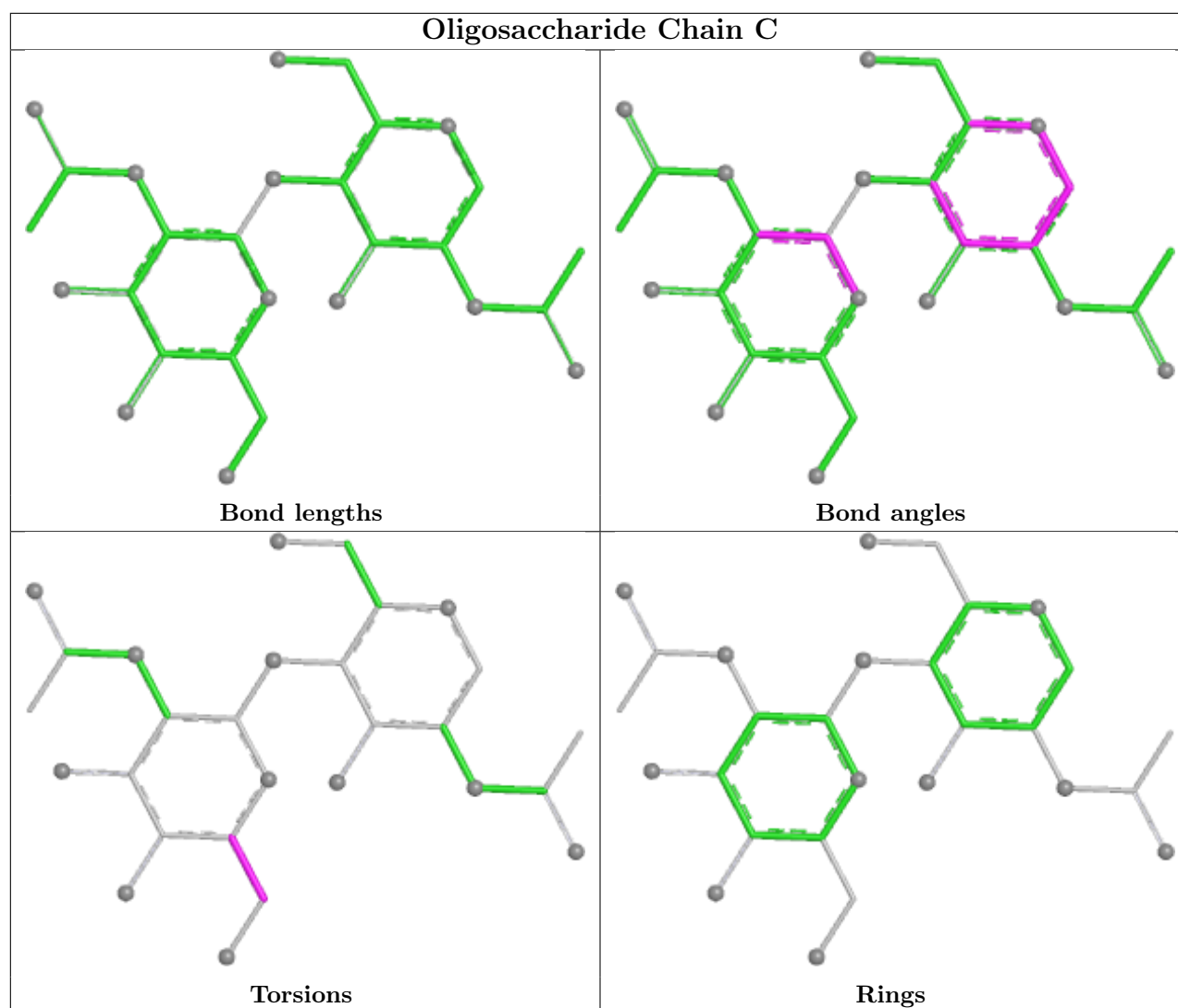
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2	NAG	2	0
4	B	1	NAG	3	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	$\# Z > 2$
5	NAG	A	703	1	14,14,15	0.81	0	17,19,21	1.72	3 (17%)

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	A	703	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	703	NAG	C1-O5-C5	5.84	120.02	112.19
5	A	703	NAG	C2-N2-C7	2.35	126.06	122.90
5	A	703	NAG	O5-C1-C2	2.06	114.48	111.29

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	567/630 (90%)	-0.22	7 (1%) 76 54	58, 73, 115, 148	0
2	H	224/224 (100%)	-0.24	2 (0%) 81 60	60, 76, 110, 128	0
3	L	220/220 (100%)	-0.42	0 100 100	64, 86, 105, 114	0
All	All	1011/1074 (94%)	-0.27	9 (0%) 81 60	58, 76, 111, 148	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	103	TYR	4.9
1	A	573	PHE	3.0
1	A	572	PRO	2.8
1	A	1	THR	2.6
1	A	363	PRO	2.4
2	H	100	GLY	2.3
1	A	575	VAL	2.3
1	A	555	PHE	2.2
1	A	552	VAL	2.1

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

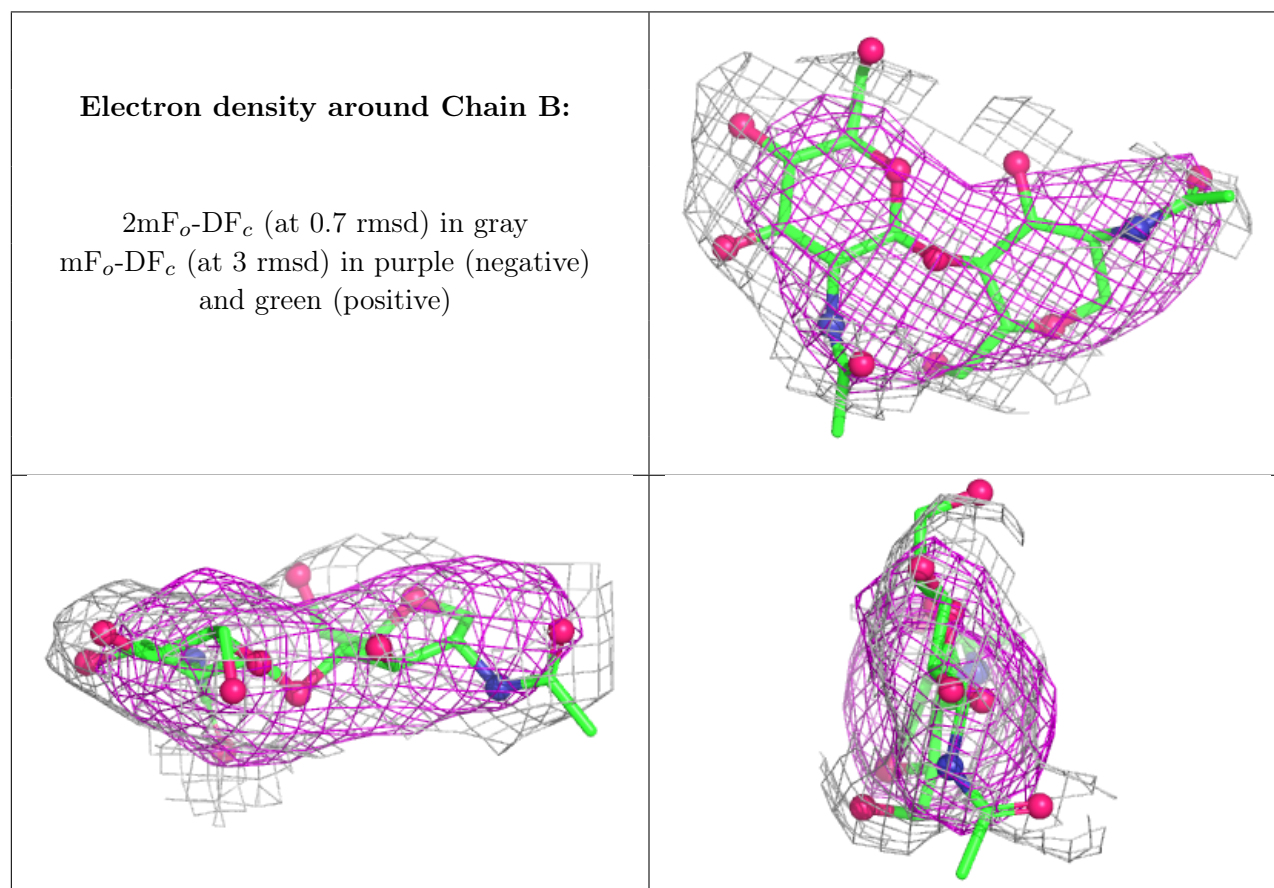
There are no non-standard protein/DNA/RNA residues in this entry.

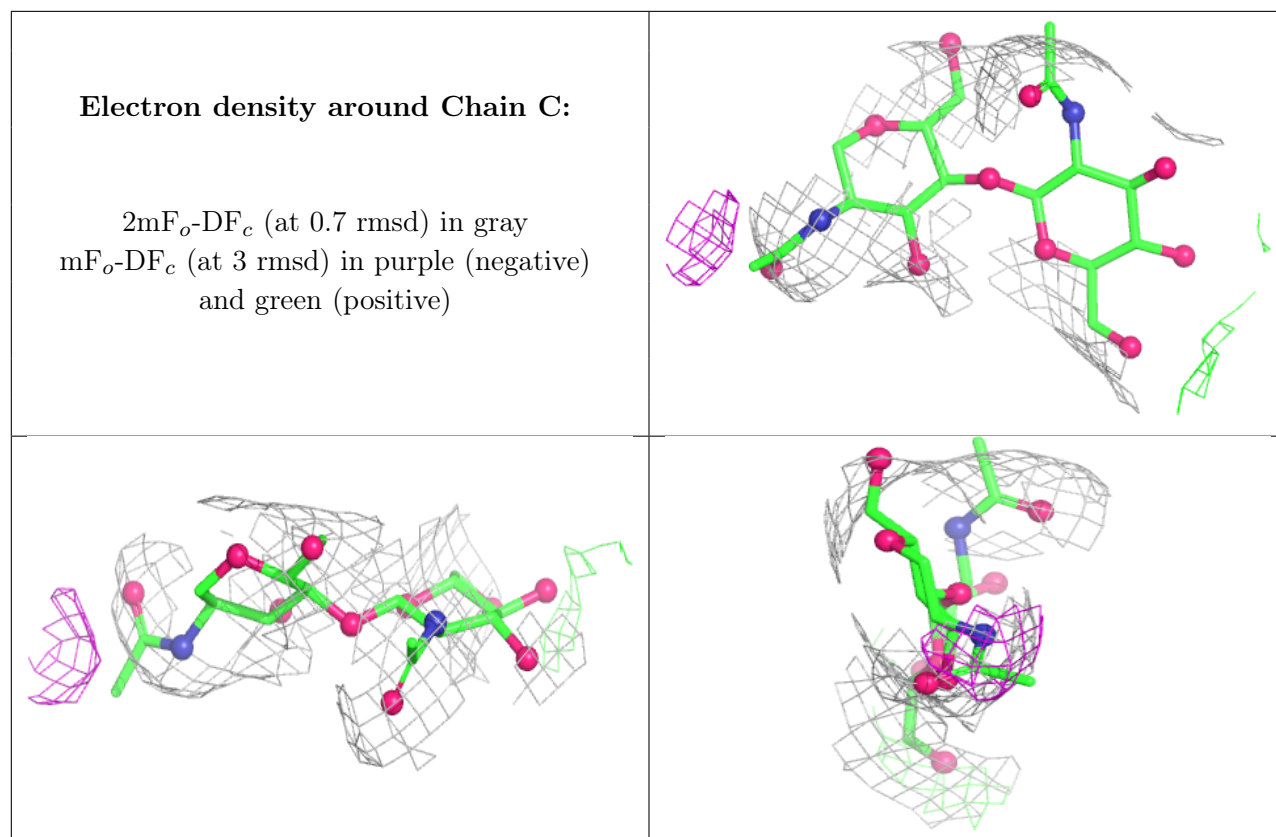
6.3 Carbohydrates [i](#)

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($\text{\AA}^2$)	Q<0.9
4	NAG	B	1	14/15	-	-	61,66,78,98	0
4	NAG	B	2	14/15	-	-	76,76,76,76	0
4	NAG	C	1	14/15	-	-	76,76,76,76	0
4	NAG	C	2	14/15	-	-	76,76,76,76	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(Å ²)	Q<0.9
5	NAG	A	703	14/15	0.80	0.11	76,76,76,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.