



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:03 PM UTC

PDB ID : 7K7J / pdb_00007k7j
Title : EphB6 receptor ectodomain
Authors : Goldgur, Y.; Himanen, J.
Deposited on : 2020-09-22
Resolution : 3.00 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

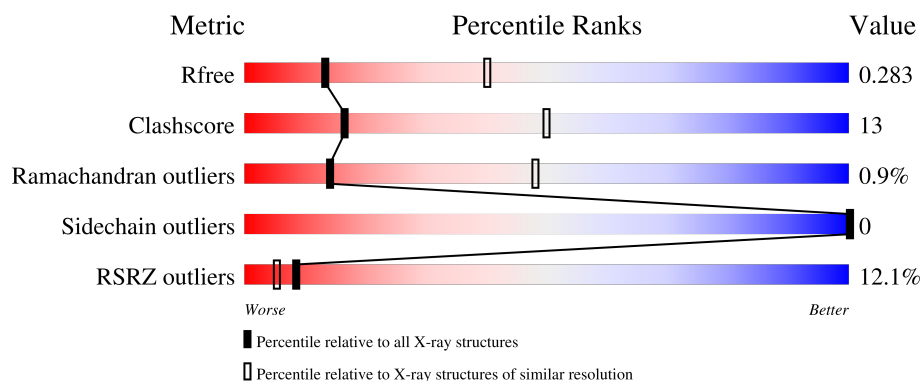
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	547	10% 56% 20% • 22%
2	B	2	50% 50%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 3255 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 Ephrin type-B receptor 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	S	0	0	0
			3227	2001	593	611	22			

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

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	105.31Å 105.31Å 188.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.05 – 3.00 48.05 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.05-3.00) 99.8 (48.05-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.16_3549: ???)	Depositor
R, R_{free}	0.235 , 0.288 0.233 , 0.283	Depositor DCC
R_{free} test set	1046 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	84.1	Xtriage
Anisotropy	0.243	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 70.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3255	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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

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	0.49	0/3312	0.82	4/4511 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	15

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	123	VAL	N-CA-C	-6.30	105.56	111.48
1	A	439	SER	N-CA-CB	6.03	121.09	110.37
1	A	10	GLU	CA-C-N	5.79	132.59	121.54
1	A	10	GLU	C-N-CA	5.79	132.59	121.54

There are no chirality outliers.

All (15) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	107	ASP	Peptide
1	A	11	THR	Peptide
1	A	114	SER	Peptide
1	A	215	GLN	Peptide
1	A	216	ALA	Peptide
1	A	250	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	A	262	MET	Peptide
1	A	293	ALA	Peptide
1	A	295	ALA	Peptide
1	A	33	GLN	Peptide
1	A	436	SER	Peptide
1	A	438	LEU	Peptide
1	A	439	SER	Peptide
1	A	440	PRO	Peptide
1	A	65	ARG	Peptide

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	3227	0	3069	82	0
2	B	28	0	25	2	0
All	All	3255	0	3094	82	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 13.

All (82) 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:325:ARG:NH1	1:A:333:ALA:O	2.08	0.87
1:A:115:TRP:O	1:A:117:LEU:N	2.11	0.80
1:A:115:TRP:CD1	1:A:115:TRP:H	2.02	0.77
1:A:104:GLU:HG3	1:A:175:GLN:HB3	1.66	0.76
1:A:402:PHE:CD2	1:A:406:GLN:HB2	2.24	0.72
1:A:65:ARG:HB2	1:A:177:GLY:HA2	1.70	0.72
1:A:71:ALA:HB3	1:A:170:PHE:HD2	1.53	0.71
1:A:18:THR:HG21	1:A:24:TRP:H	1.56	0.70
1:A:96:PHE:HB3	1:A:189:LEU:HD11	1.77	0.66
1:A:376:CYS:SG	1:A:377:LYS:N	2.69	0.65
1:A:65:ARG:HD3	1:A:175:GLN:O	1.96	0.65
1:A:19:TYR:HB3	1:A:20:PRO:HD3	1.79	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:THR:HG22	1:A:125:THR:HG23	1.82	0.61
1:A:317:CYS:O	1:A:325:ARG:NH2	2.34	0.60
1:A:304:PRO:HB2	1:A:334:PRO:HA	1.85	0.59
1:A:347:PHE:HD1	1:A:354:LEU:HD11	1.70	0.57
1:A:410:THR:OG1	1:A:411:GLU:OE1	2.21	0.57
1:A:290:LEU:HD22	1:A:302:PRO:HA	1.87	0.56
1:A:208:PHE:HZ	1:A:276:PRO:HB3	1.71	0.56
1:A:71:ALA:HB3	1:A:170:PHE:CD2	2.38	0.55
1:A:208:PHE:CZ	1:A:276:PRO:HB3	2.43	0.54
1:A:405:ARG:O	1:A:406:GLN:HB3	2.08	0.54
1:A:347:PHE:CD1	1:A:354:LEU:HD11	2.43	0.53
1:A:107:ASP:HB2	1:A:179:TYR:OH	2.09	0.52
1:A:101:ARG:NH2	1:A:107:ASP:OD2	2.43	0.52
1:A:269:ARG:NH1	1:A:295:ALA:HB1	2.24	0.52
1:A:115:TRP:CD1	1:A:115:TRP:N	2.74	0.51
1:A:355:MET:HE2	1:A:414:VAL:C	2.36	0.51
1:A:5:LEU:HB3	1:A:196:LEU:HB2	1.92	0.51
1:A:65:ARG:HD2	1:A:173:LEU:HB3	1.93	0.51
1:A:251:PRO:HB3	1:A:267:GLY:HA2	1.93	0.50
1:A:370:LEU:HD21	1:A:372:PHE:CZ	2.45	0.50
1:A:449:ASN:ND2	2:B:1:NAG:O7	2.45	0.50
1:A:238:GLU:HG2	1:A:267:GLY:O	2.12	0.50
1:A:290:LEU:HD22	1:A:302:PRO:CA	2.42	0.50
1:A:349:VAL:C	1:A:350:GLN:HG2	2.37	0.50
1:A:231:CYS:HB3	1:A:235:ALA:HB3	1.94	0.49
1:A:115:TRP:C	1:A:117:LEU:H	2.12	0.49
1:A:123:VAL:HG21	1:A:170:PHE:HE1	1.77	0.49
1:A:292:LYS:NZ	1:A:296:GLY:O	2.28	0.49
1:A:253:LEU:HD13	1:A:264:ALA:HB2	1.94	0.49
1:A:403:ASP:HB2	1:A:415:LEU:HB3	1.95	0.48
1:A:253:LEU:HG	1:A:261:TRP:CE3	2.49	0.48
1:A:370:LEU:HD12	1:A:432:VAL:O	2.14	0.48
1:A:362:ARG:HE	1:A:362:ARG:HB3	1.42	0.47
1:A:39:THR:HB	1:A:191:LEU:O	2.14	0.47
1:A:378:GLU:HG2	1:A:425:TYR:CE2	2.49	0.46
1:A:339:PRO:HD3	1:A:433:ASN:ND2	2.31	0.46
1:A:438:LEU:HD23	1:A:438:LEU:HA	1.50	0.46
1:A:208:PHE:CZ	1:A:280:ASP:HA	2.50	0.46
1:A:18:THR:CG2	1:A:24:TRP:H	2.28	0.46
1:A:400:VAL:O	1:A:401:HIS:C	2.59	0.46
1:A:377:LYS:HB2	1:A:426:ILE:CG2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:TRP:CD1	1:A:16:TRP:N	2.84	0.45
1:A:161:LEU:HD12	1:A:161:LEU:HA	1.88	0.45
1:A:269:ARG:HH11	1:A:295:ALA:HB1	1.82	0.45
1:A:454:HIS:O	1:A:455:GLU:HG3	2.17	0.45
1:A:389:GLY:HA2	2:B:2:NAG:H81	1.97	0.45
1:A:201:CYS:HB3	1:A:261:TRP:CE2	2.52	0.44
1:A:403:ASP:CB	1:A:415:LEU:HB3	2.48	0.43
1:A:216:ALA:O	1:A:225:VAL:HG21	2.17	0.43
1:A:381:GLY:HA2	1:A:390:GLY:O	2.19	0.43
1:A:161:LEU:C	1:A:163:LEU:H	2.27	0.43
1:A:210:SER:O	1:A:229:GLY:HA2	2.18	0.43
1:A:116:HIS:O	1:A:116:HIS:CG	2.72	0.43
1:A:76:HIS:HA	1:A:164:ASN:O	2.18	0.42
1:A:292:LYS:HE2	1:A:294:SER:O	2.20	0.42
1:A:277:ALA:C	1:A:279:GLY:H	2.27	0.42
1:A:441:ASP:O	1:A:442:PRO:C	2.62	0.42
1:A:48:ALA:O	1:A:51:GLY:N	2.50	0.41
1:A:256:ASN:HB2	1:A:262:MET:CG	2.50	0.41
1:A:308:HIS:ND1	1:A:320:LEU:HD23	2.35	0.41
1:A:97:THR:HA	1:A:125:THR:HA	2.02	0.41
1:A:279:GLY:O	1:A:281:LYS:HB2	2.20	0.41
1:A:289:GLY:CA	1:A:309:ALA:HB3	2.51	0.41
1:A:123:VAL:HG21	1:A:170:PHE:CE1	2.55	0.41
1:A:274:TYR:O	1:A:292:LYS:HD3	2.21	0.41
1:A:427:LEU:HD23	1:A:427:LEU:HA	1.83	0.40
1:A:108:SER:HB3	1:A:112:VAL:HB	2.03	0.40
1:A:74:ARG:HD3	1:A:167:GLU:OE1	2.21	0.40
1:A:402:PHE:CG	1:A:406:GLN:HB2	2.55	0.40
1:A:440:PRO:HA	1:A:441:ASP:O	2.22	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	A	423/547 (77%)	349 (82%)	70 (16%)	4 (1%)	14	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	HIS
1	A	280	ASP
1	A	12	SER
1	A	404	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	339/434 (78%)	339 (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. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	ASN
1	A	234	HIS
1	A	254	HIS
1	A	454	HIS

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

2 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	2,1	14,14,15	0.67	1 (7%)	17,19,21	0.56	0
2	NAG	B	2	2	14,14,15	0.30	0	17,19,21	0.54	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	B	1	2,1	-	3/6/23/26	0/1/1/1
2	NAG	B	2	2	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	NAG	C1-C2	2.38	1.55	1.52

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

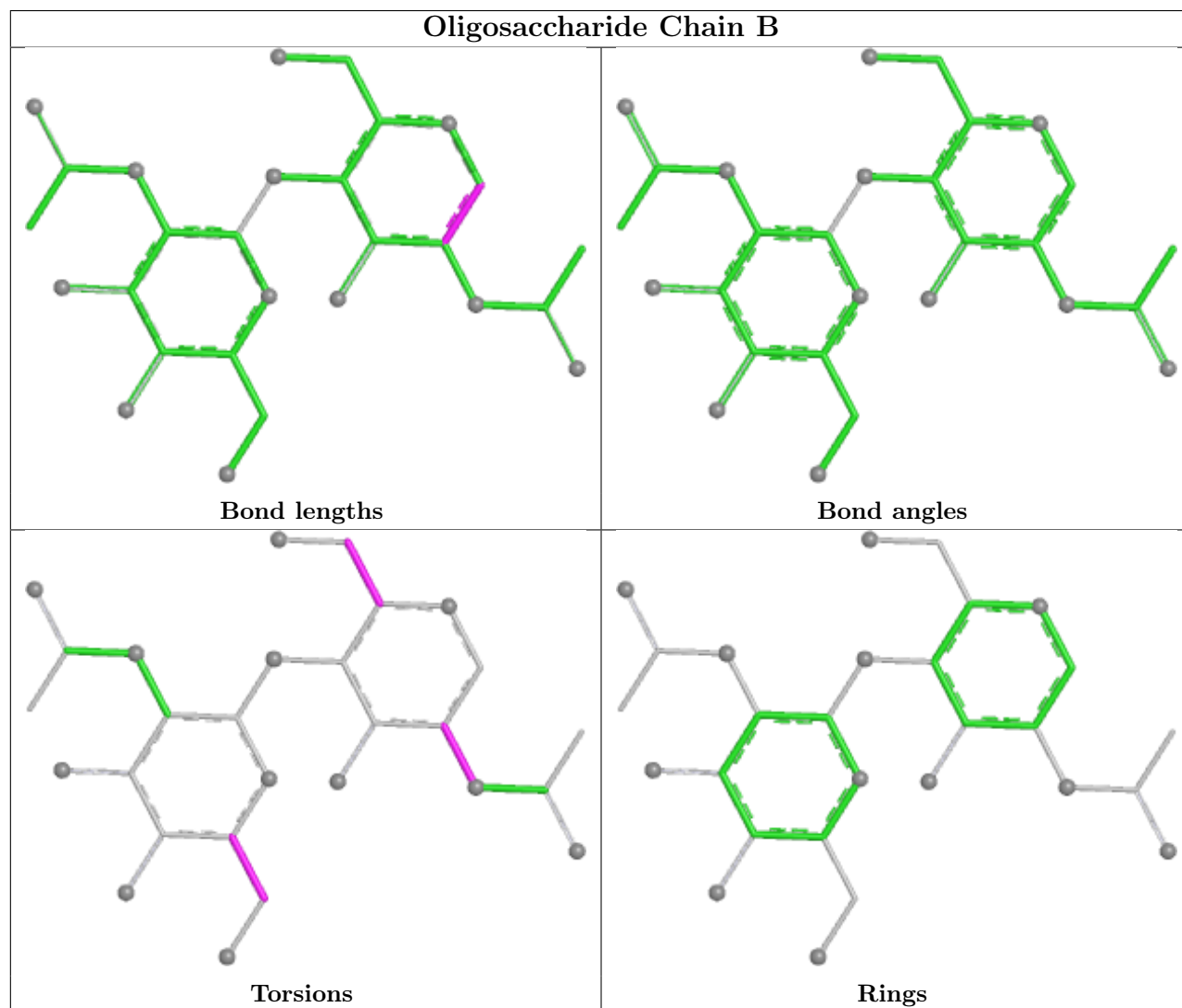
Mol	Chain	Res	Type	Atoms
2	B	1	NAG	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
2	B	2	NAG	O5-C5-C6-O6
2	B	1	NAG	C1-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	NAG	1	0
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](#)

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 ⓘ

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	429/547 (78%)	0.48	52 (12%) 8 5	44, 79, 176, 249	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	110	ASP	7.3
1	A	161	LEU	6.2
1	A	87	VAL	5.7
1	A	456	VAL	5.4
1	A	19	TYR	4.5
1	A	387	SER	4.4
1	A	250	PRO	4.4
1	A	154	GLY	4.3
1	A	13	GLU	4.2
1	A	12	SER	4.1
1	A	217	SER	4.0
1	A	401	HIS	3.9
1	A	111	SER	3.9
1	A	263	VAL	3.8
1	A	130	GLU	3.7
1	A	384	GLU	3.7
1	A	386	ALA	3.7
1	A	112	VAL	3.7
1	A	385	PRO	3.5
1	A	34	ARG	3.5
1	A	11	THR	3.5
1	A	460	VAL	3.3
1	A	49	PRO	3.2
1	A	239	GLU	3.1
1	A	88	SER	3.1
1	A	159	ALA	3.0
1	A	452	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	160	GLY	3.0
1	A	440	PRO	2.9
1	A	457	PRO	2.9
1	A	89	GLY	2.8
1	A	171	GLY	2.8
1	A	51	GLY	2.8
1	A	170	PHE	2.6
1	A	52	THR	2.6
1	A	352	SER	2.6
1	A	216	ALA	2.6
1	A	454	HIS	2.6
1	A	156	GLY	2.5
1	A	109	PRO	2.5
1	A	116	HIS	2.4
1	A	459	ALA	2.4
1	A	295	ALA	2.2
1	A	403	ASP	2.2
1	A	249	SER	2.1
1	A	281	LYS	2.1
1	A	397	ARG	2.1
1	A	53	GLY	2.1
1	A	294	SER	2.1
1	A	453	SER	2.0
1	A	157	GLN	2.0
1	A	35	ARG	2.0

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

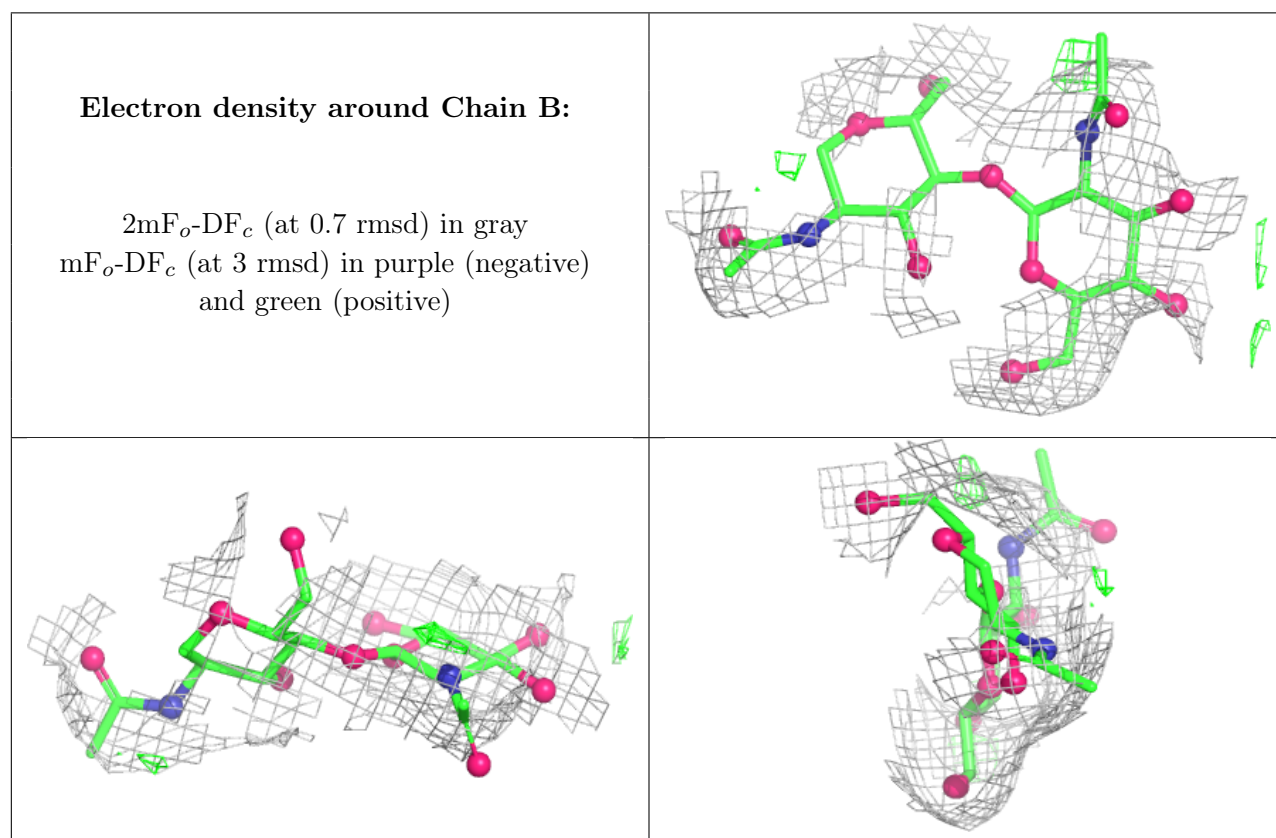
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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	NAG	B	2	14/15	0.68	0.15	165,190,194,197	0
2	NAG	B	1	14/15	0.86	0.12	124,147,183,183	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.