



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:46 PM UTC

PDB ID : 3QWR / pdb_00003qwr
Title : Crystal structure of IL-23 in complex with an adnectin
Authors : Wei, A.; Sheriff, S.
Deposited on : 2011-02-28
Resolution : 3.25 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

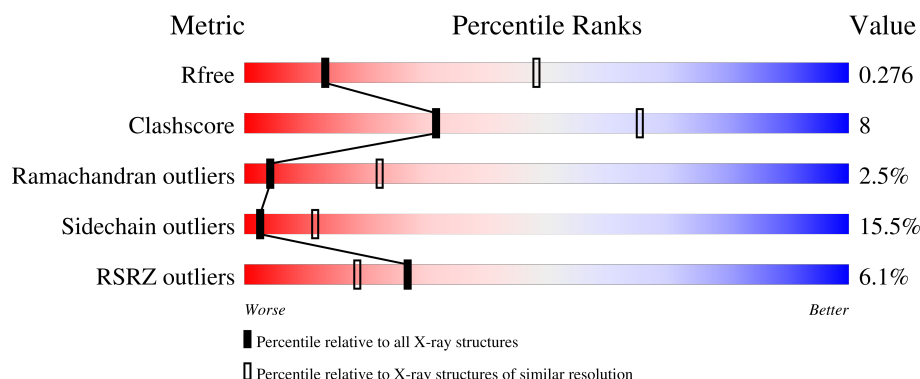
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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	1605 (3.30-3.22)
Clashscore	190562	1660 (3.30-3.22)
Ramachandran outliers	187476	1630 (3.30-3.22)
Sidechain outliers	187428	1629 (3.30-3.22)
RSRZ outliers	180081	1605 (3.30-3.22)

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	306	5% 63% 26% 5% . .
2	B	170	7% 68% 11% • 17%
3	D	109	5% 51% 30% •• 17%
4	C	4	25% 75%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4040 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 Interleukin-12 subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	295	Total	C	N	O	S	0	0	0
			2266	1432	370	453	11			

- Molecule 2 is a protein called Interleukin-23 subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	141	Total	C	N	O	S	0	0	0
			1008	650	179	173	6			

- Molecule 3 is a protein called ADNECTIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	91	Total	C	N	O	S	0	0	0
			715	462	110	142	1			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(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
4	C	4	Total	C	N	O		0	0	0
			50	28	2	20				

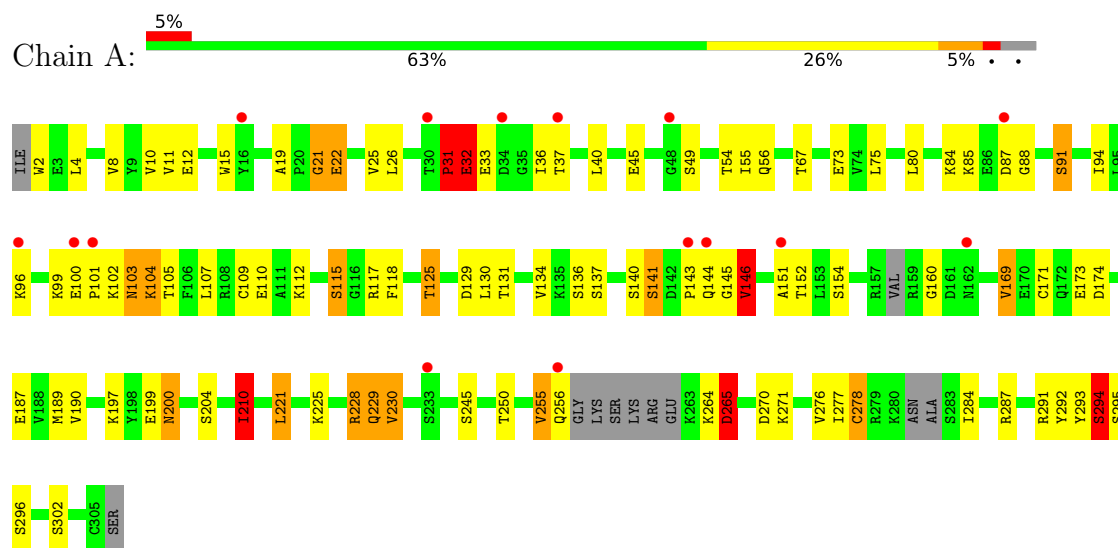
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	O	0	0
			1	1		

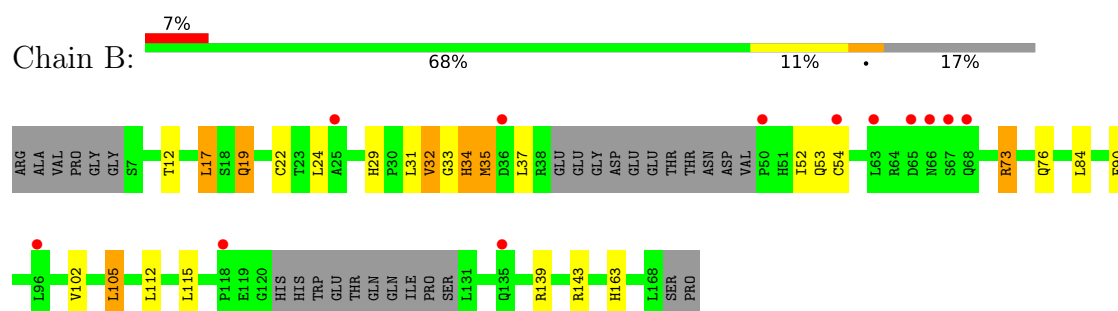
3 Residue-property plots [i](#)

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.

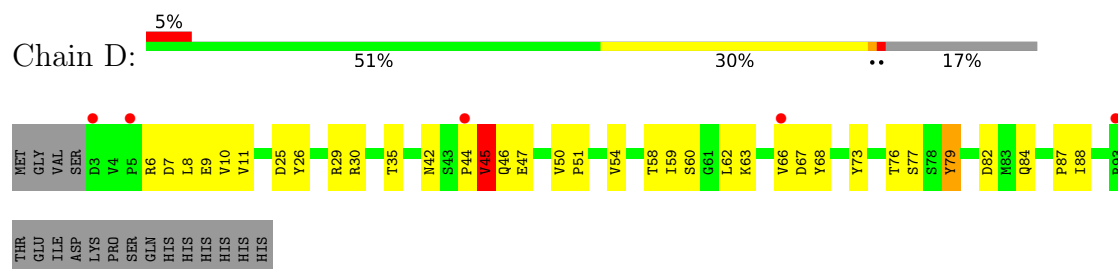
• Molecule 1: Interleukin-12 subunit beta



• Molecule 2: Interleukin-23 subunit alpha



• Molecule 3: ADNECTIN



- Molecule 4: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  25% 75%

MAG1
MAG2
BMA3
MAN4

4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.70Å 91.70Å 225.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.48 – 3.25 42.48 – 3.25	Depositor EDS
% Data completeness (in resolution range)	97.7 (42.48-3.25) 97.7 (42.48-3.25)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.27 (at 3.25Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.13.0, BUSTER 2.13.0	Depositor
R, R_{free}	0.234 , 0.264 0.241 , 0.276	Depositor DCC
R_{free} test set	1003 reflections (7.83%)	wwPDB-VP
Wilson B-factor (Å ²)	88.3	Xtriage
Anisotropy	0.932	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 111.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4040	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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, MAN

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	0.84	2/2321 (0.1%)	1.30	24/3168 (0.8%)
2	B	0.78	1/1034 (0.1%)	1.30	3/1412 (0.2%)
3	D	0.82	1/736 (0.1%)	1.33	10/1012 (1.0%)
All	All	0.82	4/4091 (0.1%)	1.31	37/5592 (0.7%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	32	VAL	CA-C	6.88	1.61	1.52
1	A	293	TYR	C-N	-5.89	1.25	1.33
3	D	46	GLN	CA-C	-5.29	1.46	1.52
1	A	102	LYS	C-N	5.26	1.39	1.33

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	33	GLY	CA-C-N	12.71	141.58	122.65
2	B	33	GLY	C-N-CA	12.71	141.58	122.65
3	D	45	VAL	N-CA-CB	10.03	127.78	111.23
3	D	44	PRO	N-CA-C	9.78	126.55	113.84
1	A	11	VAL	N-CA-C	9.61	120.14	110.82
1	A	200	ASN	CA-CB-CG	7.84	120.44	112.60
1	A	31	PRO	CA-C-N	7.73	135.62	121.70
1	A	31	PRO	C-N-CA	7.73	135.62	121.70
1	A	145	GLY	N-CA-C	7.71	120.39	111.36
1	A	115	SER	N-CA-C	-7.09	100.50	110.50
1	A	294	SER	N-CA-C	-6.88	96.14	110.80
1	A	141	SER	CA-C-N	6.74	131.23	123.21
1	A	141	SER	C-N-CA	6.74	131.23	123.21
1	A	146	VAL	N-CA-CB	6.56	121.10	111.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	22	GLU	N-CA-C	6.29	124.20	110.80
3	D	7	ASP	CA-CB-CG	6.13	118.73	112.60
1	A	265	ASP	CA-CB-CG	5.94	118.54	112.60
3	D	79	TYR	CA-C-N	5.91	129.01	120.38
3	D	79	TYR	C-N-CA	5.91	129.01	120.38
1	A	264	LYS	CA-C-N	5.87	132.75	121.54
1	A	264	LYS	C-N-CA	5.87	132.75	121.54
2	B	34	HIS	N-CA-CB	-5.78	100.98	110.52
3	D	51	PRO	CA-C-N	5.75	129.45	120.82
3	D	51	PRO	C-N-CA	5.75	129.45	120.82
3	D	7	ASP	CA-C-N	5.74	131.58	122.74
3	D	7	ASP	C-N-CA	5.74	131.58	122.74
1	A	210	ILE	CA-CB-CG2	5.67	120.14	110.50
1	A	21	GLY	CA-C-N	5.61	132.26	121.54
1	A	21	GLY	C-N-CA	5.61	132.26	121.54
3	D	8	LEU	N-CA-C	-5.45	99.55	108.76
1	A	270	ASP	N-CA-C	-5.39	106.70	113.28
1	A	94	ILE	N-CA-C	5.38	115.58	110.53
1	A	33	GLU	N-CA-C	5.20	118.57	111.39
1	A	174	ASP	N-CA-C	5.16	116.59	111.07
1	A	125	THR	N-CA-C	-5.07	106.25	112.90
1	A	32	GLU	CA-C-N	5.05	129.53	122.36
1	A	32	GLU	C-N-CA	5.05	129.53	122.36

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	A	2266	0	2065	39	0
2	B	1008	0	943	14	0
3	D	715	0	673	20	0
4	C	50	0	43	3	0
5	A	1	0	0	0	0
All	All	4040	0	3724	63	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 8.

All (63) 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:134:VAL:HG21	1:A:169:VAL:HG11	1.51	0.92
3:D:35:THR:HG22	3:D:47:GLU:HB3	1.69	0.74
3:D:59:ILE:HG22	3:D:62:LEU:HD21	1.72	0.71
1:A:229:GLN:HB3	1:A:277:ILE:HG13	1.75	0.69
1:A:245:SER:OG	2:B:163:HIS:ND1	2.25	0.69
3:D:30:ARG:HD3	3:D:76:THR:O	1.92	0.68
2:B:35:MET:HE2	3:D:30:ARG:HH12	1.61	0.64
1:A:25:VAL:HG22	1:A:54:THR:HG22	1.80	0.64
1:A:255:VAL:HG23	1:A:265:ASP:HB3	1.79	0.64
1:A:256:GLN:HE22	3:D:6:ARG:H	1.45	0.64
1:A:2:TRP:CZ2	4:C:1:NAG:H4	2.33	0.64
3:D:50:VAL:CG1	3:D:54:VAL:HB	2.29	0.63
1:A:230:VAL:CG2	1:A:278:CYS:HB2	2.29	0.62
3:D:73:TYR:CE2	3:D:87:PRO:HB3	2.34	0.62
1:A:136:SER:HB3	1:A:146:VAL:HG22	1.83	0.60
1:A:109:CYS:O	1:A:210:ILE:HG12	2.01	0.60
2:B:17:LEU:HD22	2:B:112:LEU:HB2	1.85	0.59
1:A:250:THR:OG1	1:A:291:ARG:HG3	2.04	0.57
2:B:29:HIS:ND1	3:D:30:ARG:HD2	2.20	0.57
1:A:31:PRO:HA	1:A:32:GLU:CB	2.37	0.55
2:B:35:MET:HE2	3:D:30:ARG:NH1	2.22	0.55
3:D:62:LEU:HB3	3:D:68:TYR:CE2	2.42	0.54
1:A:278:CYS:SG	1:A:284:ILE:HD11	2.48	0.54
1:A:292:TYR:HE1	2:B:19:GLN:HG2	1.75	0.51
2:B:35:MET:HE2	3:D:77:SER:HA	1.93	0.51
1:A:85:LYS:HE3	1:A:88:GLY:O	2.11	0.51
1:A:4:LEU:O	1:A:8:VAL:HB	2.10	0.51
3:D:79:TYR:HB3	3:D:82:ASP:HB2	1.92	0.50
1:A:255:VAL:HG12	1:A:284:ILE:HG12	1.94	0.50
1:A:137:SER:HA	1:A:146:VAL:HG13	1.94	0.50
1:A:101:PRO:HD2	1:A:105:THR:HG21	1.93	0.49
3:D:50:VAL:HG12	3:D:54:VAL:HB	1.94	0.48
3:D:73:TYR:CD2	3:D:87:PRO:HB3	2.48	0.48
1:A:140:SER:O	1:A:143:PRO:HD3	2.13	0.48
2:B:139:ARG:O	2:B:143:ARG:HG3	2.13	0.48
1:A:287:ARG:HH22	3:D:25:ASP:CG	2.21	0.48
3:D:63:LYS:O	3:D:66:VAL:HB	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:LEU:HD21	1:A:284:ILE:HD12	1.97	0.46
2:B:35:MET:CE	3:D:30:ARG:HH12	2.27	0.46
1:A:146:VAL:HB	1:A:173:GLU:HA	1.98	0.46
2:B:24:LEU:HB3	2:B:105:LEU:HD13	1.98	0.45
1:A:292:TYR:HB3	2:B:22:CYS:SG	2.57	0.45
2:B:52:ILE:CD1	2:B:73:ARG:HB3	2.47	0.45
2:B:34:HIS:O	3:D:30:ARG:NH2	2.50	0.44
3:D:26:TYR:O	3:D:29:ARG:HG2	2.16	0.44
1:A:103:ASN:O	1:A:104:LYS:C	2.61	0.44
1:A:151:ALA:HA	1:A:169:VAL:HG12	2.00	0.44
1:A:228:ARG:NE	1:A:228:ARG:HA	2.34	0.43
2:B:90:PHE:HD1	2:B:102:VAL:HG11	1.82	0.43
1:A:134:VAL:HG22	1:A:190:VAL:HG22	2.01	0.42
4:C:2:NAG:H61	4:C:3:BMA:O2	2.20	0.42
1:A:115:SER:C	1:A:117:ARG:H	2.27	0.42
1:A:294:SER:O	1:A:295:SER:OG	2.29	0.42
1:A:15:TRP:CE2	1:A:84:LYS:HD2	2.55	0.42
1:A:118:PHE:CZ	1:A:171:CYS:HB2	2.55	0.42
1:A:26:LEU:HD13	1:A:55:ILE:HD12	2.02	0.42
1:A:91:SER:HB3	1:A:199:GLU:OE1	2.20	0.42
1:A:294:SER:OG	1:A:295:SER:N	2.53	0.42
1:A:221:LEU:HD11	1:A:284:ILE:HD12	2.03	0.41
1:A:230:VAL:HG23	1:A:278:CYS:HB2	1.99	0.40
1:A:12:GLU:OE2	4:C:2:NAG:H83	2.22	0.40
1:A:36:ILE:O	1:A:49:SER:HA	2.21	0.40
3:D:9:GLU:HG2	3:D:10:VAL:N	2.36	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	287/306 (94%)	258 (90%)	18 (6%)	11 (4%)	2	15
2	B	135/170 (79%)	128 (95%)	6 (4%)	1 (1%)	18	48
3	D	89/109 (82%)	80 (90%)	8 (9%)	1 (1%)	11	38
All	All	511/585 (87%)	466 (91%)	32 (6%)	13 (2%)	4	22

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	19	ALA
1	A	22	GLU
1	A	225	LYS
1	A	265	ASP
1	A	294	SER
2	B	32	VAL
1	A	21	GLY
1	A	32	GLU
1	A	104	LYS
1	A	160	GLY
1	A	31	PRO
3	D	45	VAL
1	A	100	GLU

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	241/277 (87%)	198 (82%)	43 (18%)	2	7
2	B	96/144 (67%)	83 (86%)	13 (14%)	4	16
3	D	77/99 (78%)	69 (90%)	8 (10%)	7	25
All	All	414/520 (80%)	350 (84%)	64 (16%)	2	12

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	A	37	THR
1	A	40	LEU
1	A	45	GLU
1	A	56	GLN
1	A	67	THR
1	A	73	GLU
1	A	75	LEU
1	A	80	LEU
1	A	87	ASP
1	A	91	SER
1	A	96	LYS
1	A	99	LYS
1	A	103	ASN
1	A	107	LEU
1	A	110	GLU
1	A	112	LYS
1	A	125	THR
1	A	129	ASP
1	A	130	LEU
1	A	131	THR
1	A	141	SER
1	A	144	GLN
1	A	146	VAL
1	A	152	THR
1	A	154	SER
1	A	169	VAL
1	A	187	GLU
1	A	189	MET
1	A	197	LYS
1	A	200	ASN
1	A	204	SER
1	A	210	ILE
1	A	221	LEU
1	A	228	ARG
1	A	229	GLN
1	A	230	VAL
1	A	255	VAL
1	A	271	LYS
1	A	276	VAL
1	A	278	CYS
1	A	296	SER
1	A	302	SER

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Mol	Chain	Res	Type
2	B	12	THR
2	B	17	LEU
2	B	19	GLN
2	B	31	LEU
2	B	35	MET
2	B	37	LEU
2	B	53	GLN
2	B	54	CYS
2	B	73	ARG
2	B	76	GLN
2	B	84	LEU
2	B	105	LEU
2	B	115	LEU
3	D	11	VAL
3	D	42	ASN
3	D	45	VAL
3	D	58	THR
3	D	60	SER
3	D	67	ASP
3	D	84	GLN
3	D	88	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	144	GLN
1	A	194	HIS
1	A	229	GLN
1	A	256	GLN
2	B	13	GLN
2	B	51	HIS
2	B	75	HIS
3	D	46	GLN

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	C	1	1,4	14,14,15	0.57	0	17,19,21	1.40	2 (11%)
4	NAG	C	2	4	14,14,15	0.44	0	17,19,21	1.53	2 (11%)
4	BMA	C	3	4	11,11,12	0.59	0	15,15,17	2.05	2 (13%)
4	MAN	C	4	4	11,11,12	0.48	0	15,15,17	1.13	1 (6%)

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	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
4	BMA	C	3	4	-	0/2/19/22	0/1/1/1
4	MAN	C	4	4	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	3	BMA	C1-O5-C5	6.07	120.32	112.19
4	C	2	NAG	C1-O5-C5	4.74	118.54	112.19
4	C	1	NAG	C1-O5-C5	4.41	118.10	112.19
4	C	3	BMA	C1-C2-C3	3.57	114.84	109.64
4	C	4	MAN	C1-O5-C5	3.38	116.72	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1	NAG	C1-C2-N2	2.74	114.75	110.43
4	C	2	NAG	C1-C2-N2	2.14	113.80	110.43

There are no chirality outliers.

All (2) torsion outliers are listed below:

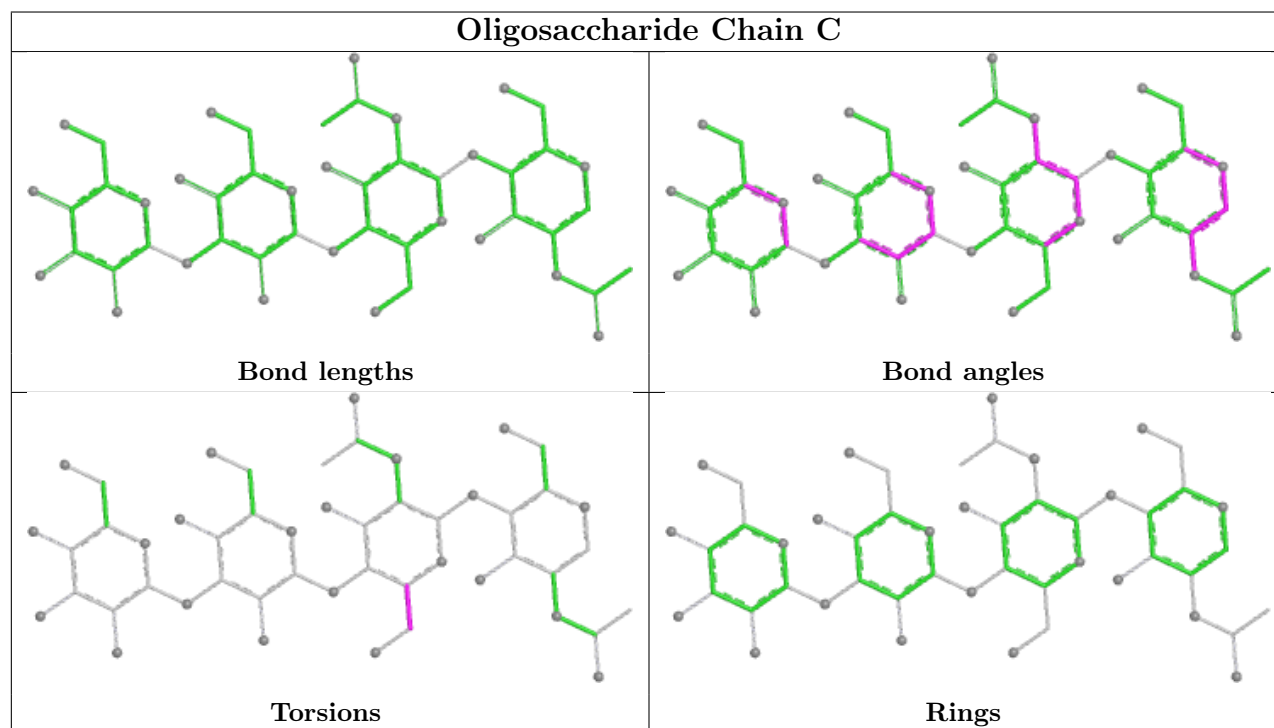
Mol	Chain	Res	Type	Atoms
4	C	2	NAG	C4-C5-C6-O6
4	C	2	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	3	BMA	1	0
4	C	2	NAG	2	0
4	C	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

There are no ligands in this entry.

5.7 Other polymers

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 [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	295/306 (96%)	0.54	15 (5%) 33 22	69, 95, 133, 151	0
2	B	141/170 (82%)	0.78	12 (8%) 16 13	74, 101, 141, 165	0
3	D	91/109 (83%)	0.74	5 (5%) 30 21	67, 99, 143, 156	0
All	All	527/585 (90%)	0.64	32 (6%) 27 19	67, 98, 139, 165	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	5	PRO	4.9
1	A	162	ASN	3.8
2	B	135	GLN	3.8
1	A	87	ASP	3.2
1	A	101	PRO	3.2
2	B	67	SER	3.2
2	B	68	GLN	3.0
2	B	63	LEU	2.9
1	A	100	GLU	2.8
2	B	54	CYS	2.8
2	B	50	PRO	2.6
3	D	93	ARG	2.6
2	B	96	LEU	2.5
3	D	3	ASP	2.4
2	B	36	ASP	2.4
1	A	233	SER	2.3
3	D	44	PRO	2.2
1	A	144	GLN	2.2
2	B	66	ASN	2.2
1	A	151	ALA	2.2
1	A	143	PRO	2.2
1	A	37	THR	2.1
1	A	256	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	48	GLY	2.1
2	B	65	ASP	2.1
1	A	96	LYS	2.1
1	A	30	THR	2.1
1	A	16	TYR	2.1
3	D	66	VAL	2.1
2	B	25	ALA	2.0
2	B	118	PRO	2.0
1	A	34	ASP	2.0

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

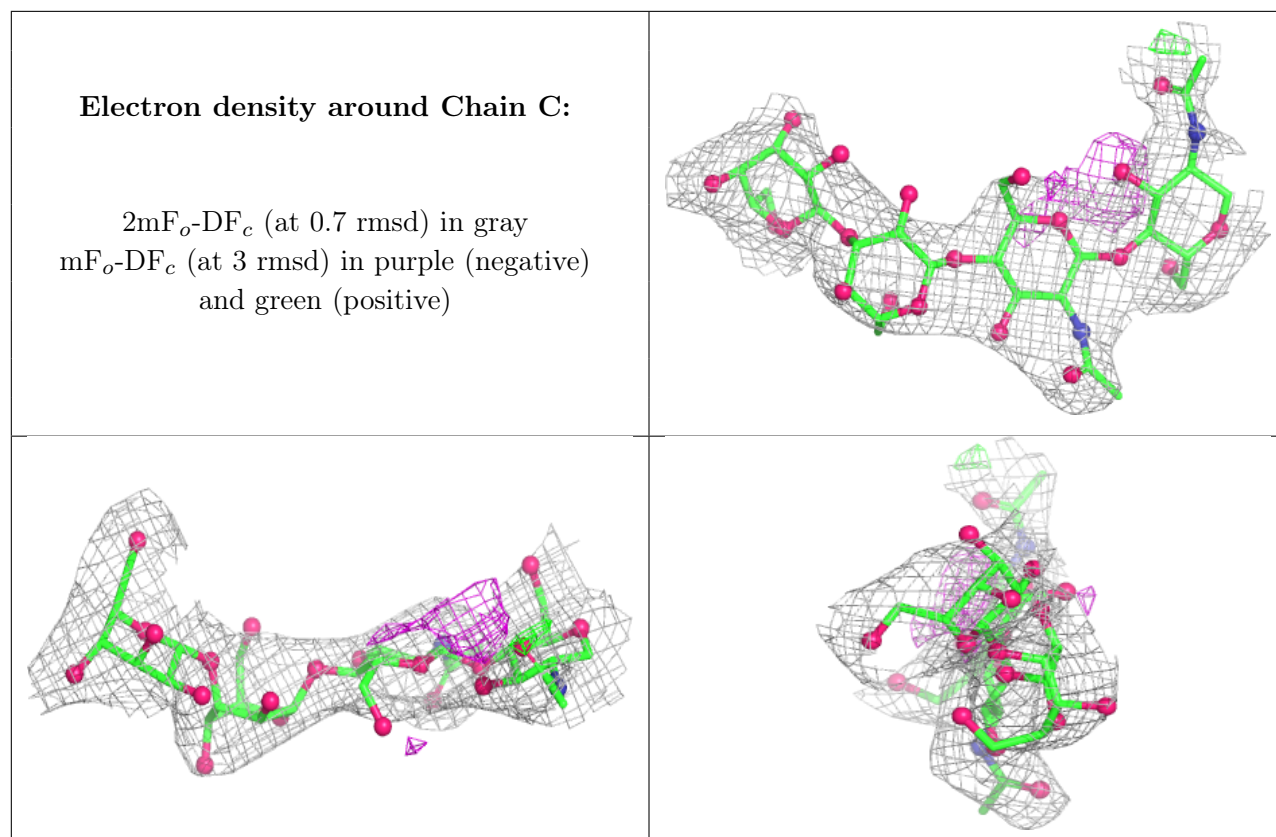
There are no non-standard protein/DNA/RNA residues in this entry.

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
4	NAG	C	1	14/15	-	-	91,98,109,109	0
4	NAG	C	2	14/15	-	-	91,102,110,114	0
4	BMA	C	3	11/12	-	-	123,130,135,136	0
4	MAN	C	4	11/12	-	-	132,134,136,137	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.