



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:54 PM UTC

PDB ID : 1H3Y / pdb_00001h3y
Title : Crystal structure of a human IgG1 Fc-fragment,high salt condition
Authors : Krapp, S.; Mimura, Y.; Jefferis, R.; Huber, R.; Sondermann, P.
Deposited on : 2002-09-19
Resolution : 4.10 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

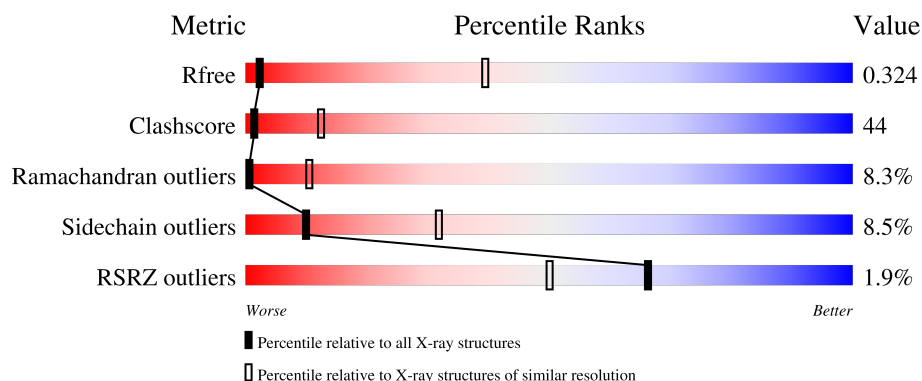
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.10 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	1243 (4.40-3.80)
Clashscore	190562	1293 (4.40-3.80)
Ramachandran outliers	187476	1206 (4.40-3.80)
Sidechain outliers	187428	1193 (4.40-3.80)
RSRZ outliers	180081	1240 (4.40-3.80)

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	223	2% 43% 40% 9% 7%
1	B	223	37% 43% 12% 8%
2	C	9	33% 67%
3	D	9	33% 67%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	C	5	X	-	-	-
2	MAN	C	7	-	-	X	-
2	NAG	C	8	-	-	X	-
2	FUL	C	9	-	-	X	-

2 Entry composition [i](#)

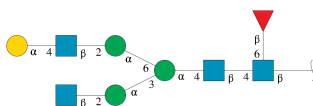
There are 3 unique types of molecules in this entry. The entry contains 3015 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 IG GAMMA-1 CHAIN C REGION.

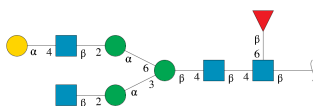
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	1
			1373	848	245	274	6			
1	B	206	Total	C	N	O	S	0	0	0
			1422	886	246	283	7			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	9	Total	C	N	O	0	0	0
			110	62	4	44			

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

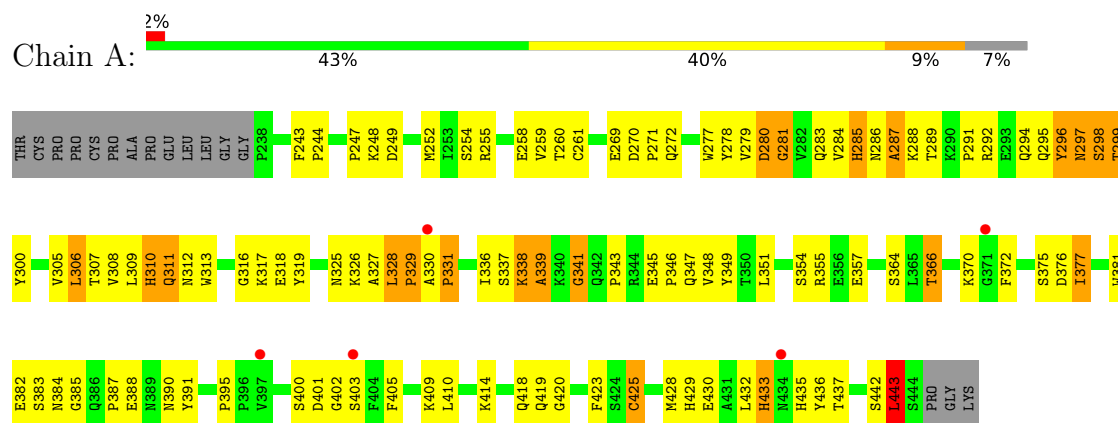


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	9	Total	C	N	O	0	0	0
			110	62	4	44			

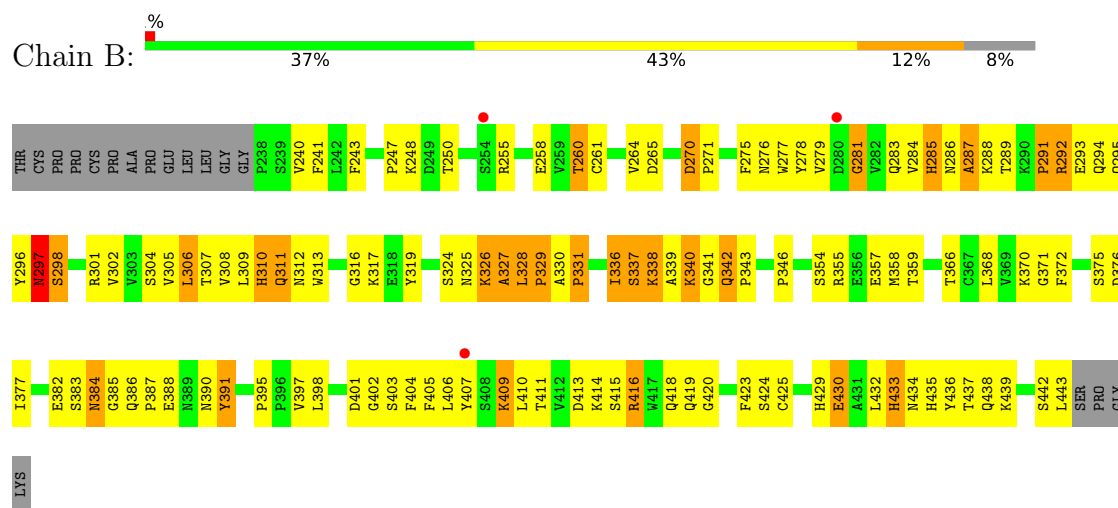
3 Residue-property plots

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: IG GAMMA-1 CHAIN C REGION



• Molecule 1: IG GAMMA-1 CHAIN C REGION



• Molecule 2: alpha-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[beta-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1
MAG2
MAN3
MAN4
NAG5
GLA6
MAN7
MAG8
FUL9

- Molecule 3: alpha-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[beta-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:

33%

67%

MAG1
MAG2
BMA3
MAN4
NAG5
GLA6
MAN7
MAG8
FUL9

4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	152.94Å 152.94Å 116.96Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 4.10 50.00 – 4.13	Depositor EDS
% Data completeness (in resolution range)	80.0 (50.00-4.10) 79.0 (50.00-4.13)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 4.13Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.285 , 0.349 0.269 , 0.324	Depositor DCC
R_{free} test set	285 reflections (5.46%)	wwPDB-VP
Wilson B-factor (Å ²)	59.4	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 271.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	3015	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.48% 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: BMA, MAN, FUL, NAG, GLA

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	0.70	2/1415 (0.1%)	1.36	23/1927 (1.2%)
1	B	0.71	0/1466	1.38	23/1988 (1.2%)
All	All	0.71	2/2881 (0.1%)	1.37	46/3915 (1.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	299	THR	CA-CB	5.78	1.62	1.53
1	A	443	LEU	C-N	-5.55	1.25	1.33

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	402	GLY	N-CA-C	-12.43	98.02	115.30
1	A	402	GLY	N-CA-C	-11.90	98.65	115.43
1	B	395	PRO	N-CA-C	-9.74	98.81	110.70
1	A	328	LEU	CA-C-N	9.23	131.37	119.84
1	A	328	LEU	C-N-CA	9.23	131.37	119.84
1	A	341	GLY	N-CA-C	8.86	121.08	112.04
1	B	342	GLN	CA-C-N	8.05	128.10	119.89
1	B	342	GLN	C-N-CA	8.05	128.10	119.89
1	A	351	LEU	CA-C-N	7.73	128.35	120.38
1	A	351	LEU	C-N-CA	7.73	128.35	120.38
1	B	298	SER	N-CA-C	-7.71	101.00	113.19
1	B	386	GLN	CA-C-N	7.55	130.44	120.25
1	B	386	GLN	C-N-CA	7.55	130.44	120.25
1	B	328	LEU	CA-C-N	7.50	129.22	119.84
1	B	328	LEU	C-N-CA	7.50	129.22	119.84
1	A	385	GLY	N-CA-C	-7.49	102.63	113.86
1	B	270	ASP	CA-C-N	7.33	127.50	119.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	270	ASP	C-N-CA	7.33	127.50	119.87
1	B	385	GLY	N-CA-C	-7.03	103.32	113.86
1	A	395	PRO	N-CA-C	-6.96	102.21	110.70
1	B	338	LYS	N-CA-C	-6.80	101.40	110.24
1	A	244	PRO	CA-C-N	6.46	126.37	119.85
1	A	244	PRO	C-N-CA	6.46	126.37	119.85
1	A	348	VAL	N-CA-C	6.35	117.74	108.53
1	B	255	ARG	N-CA-C	6.22	118.56	110.53
1	A	347	GLN	N-CA-C	-6.17	99.14	109.07
1	A	327	ALA	N-CA-C	-6.12	104.30	110.97
1	A	337	SER	CA-C-N	-6.11	114.05	122.30
1	A	337	SER	C-N-CA	-6.11	114.05	122.30
1	B	370	LYS	N-CA-C	6.11	118.49	109.69
1	B	305	VAL	N-CA-C	5.91	116.99	108.48
1	B	327	ALA	N-CA-C	-5.83	104.83	111.07
1	B	366	THR	N-CA-C	5.78	117.86	108.26
1	A	305	VAL	N-CA-C	5.75	116.75	108.48
1	A	298	SER	N-CA-C	-5.53	105.15	113.89
1	A	370	LYS	N-CA-C	5.51	117.63	109.69
1	B	388	GLU	N-CA-C	5.46	118.82	110.36
1	B	297	ASN	N-CA-CB	-5.45	101.27	110.49
1	B	366	THR	CB-CA-C	-5.39	101.84	110.14
1	B	384	ASN	N-CA-C	-5.25	99.62	110.80
1	A	388	GLU	N-CA-C	5.22	117.55	110.35
1	A	296	TYR	N-CA-C	5.16	121.78	110.80
1	A	366	THR	N-CA-C	5.15	116.81	108.26
1	B	430	GLU	N-CA-C	5.14	116.88	111.28
1	A	270	ASP	CA-C-N	5.12	125.19	119.87
1	A	270	ASP	C-N-CA	5.12	125.19	119.87

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	1373	0	997	95	0
1	B	1422	0	1100	121	1
2	C	110	0	94	22	0
3	D	110	0	94	13	0
All	All	3015	0	2285	235	1

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 44.

All (235) 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)
2:C:8:NAG:H83	3:D:9:FUL:H4	1.25	1.18
1:A:346:PRO:HB3	1:A:372:PHE:HB3	1.24	1.10
1:B:325:ASN:ND2	1:B:326:LYS:H	1.52	1.06
1:A:283:GLN:HG2	1:A:285:HIS:H	1.30	0.96
3:D:7:MAN:H2	3:D:8:NAG:N2	1.78	0.96
1:B:308:VAL:HG13	1:B:319:TYR:OH	1.68	0.93
1:B:346:PRO:HB3	1:B:372:PHE:HB3	1.48	0.93
1:B:293:GLU:HA	1:B:302:VAL:HG22	1.52	0.92
1:A:346:PRO:CB	1:A:372:PHE:HB3	2.04	0.88
2:C:8:NAG:H83	3:D:9:FUL:C4	2.02	0.88
1:B:346:PRO:CB	1:B:372:PHE:HB3	2.08	0.84
2:C:8:NAG:C8	3:D:9:FUL:H4	2.09	0.82
3:D:7:MAN:H2	3:D:8:NAG:C7	2.10	0.81
1:A:418:GLN:C	1:A:420:GLY:H	1.90	0.80
1:B:346:PRO:HB3	1:B:372:PHE:CB	2.12	0.79
1:A:277:TRP:HE1	1:A:289:THR:HG23	1.48	0.77
1:B:325:ASN:ND2	1:B:326:LYS:N	2.29	0.77
1:B:241:PHE:HE1	3:D:2:NAG:H4	1.50	0.77
2:C:2:NAG:H83	2:C:9:FUL:H3	1.67	0.77
1:A:280:ASP:OD1	1:A:318:GLU:N	2.15	0.76
1:A:297:ASN:OD1	2:C:9:FUL:H63	1.86	0.76
1:A:390:ASN:O	1:A:410:LEU:HD12	1.84	0.76
1:B:346:PRO:CA	1:B:372:PHE:HB3	2.15	0.76
3:D:7:MAN:C2	3:D:8:NAG:N2	2.50	0.75
1:B:325:ASN:HD22	1:B:326:LYS:H	1.34	0.74
1:B:418:GLN:C	1:B:420:GLY:H	1.94	0.74
1:A:259:VAL:HG23	1:A:308:VAL:HG21	1.70	0.72
1:B:324:SER:OG	1:B:331:PRO:HB3	1.89	0.72
1:A:277:TRP:NE1	1:A:289:THR:HG23	2.04	0.72
1:A:325:ASN:ND2	1:A:326:LYS:H	1.87	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:4:MAN:O5	2:C:5:NAG:H82	1.89	0.71
3:D:5:NAG:O3	3:D:6:GLA:H2	1.90	0.71
1:B:283:GLN:HG2	1:B:285:HIS:H	1.56	0.71
2:C:7:MAN:O6	2:C:8:NAG:H82	1.89	0.71
2:C:7:MAN:H4	2:C:8:NAG:H2	1.72	0.71
1:A:432:LEU:O	1:A:435:HIS:N	2.22	0.69
1:B:401:ASP:HB2	1:B:403:SER:OG	1.93	0.69
1:A:252:MET:SD	1:A:428:MET:HE1	2.32	0.69
1:B:296:TYR:CZ	1:B:301:ARG:HD3	2.30	0.67
1:A:436:TYR:CD1	1:A:437:THR:N	2.63	0.67
1:B:276:ASN:HB3	1:B:278:TYR:HE1	1.59	0.67
1:A:401:ASP:HB2	1:A:403:SER:OG	1.95	0.67
1:B:406:LEU:O	1:B:406:LEU:HD12	1.96	0.66
1:B:432:LEU:O	1:B:435:HIS:N	2.24	0.66
1:A:338:LYS:HD3	1:A:376:ASP:CG	2.21	0.65
1:B:406:LEU:HD12	1:B:406:LEU:C	2.21	0.65
1:A:338:LYS:HD3	1:A:376:ASP:OD2	1.96	0.65
1:B:346:PRO:HA	1:B:372:PHE:HB3	1.78	0.65
1:B:436:TYR:CD1	1:B:437:THR:N	2.64	0.64
1:A:272:GLN:O	1:A:325:ASN:ND2	2.31	0.64
1:B:328:LEU:HD12	1:B:329:PRO:HD2	1.78	0.64
1:B:264:VAL:O	1:B:265:ASP:HB2	1.97	0.63
1:B:418:GLN:O	1:B:420:GLY:N	2.30	0.63
2:C:2:NAG:H83	2:C:9:FUL:C3	2.28	0.63
1:B:405:PHE:CD1	1:B:405:PHE:C	2.75	0.63
1:A:258:GLU:HG3	2:C:6:GLA:H62	1.81	0.63
1:A:382:GLU:HA	1:A:387:PRO:HA	1.80	0.62
1:B:279:VAL:O	1:B:281:GLY:N	2.27	0.62
1:A:313:TRP:CZ2	1:A:338:LYS:HB2	2.34	0.62
1:B:328:LEU:HD12	1:B:329:PRO:CD	2.30	0.62
1:B:401:ASP:HB2	1:B:403:SER:CB	2.29	0.62
1:A:299:THR:OG1	1:A:300:TYR:N	2.31	0.61
1:A:418:GLN:O	1:A:420:GLY:N	2.33	0.61
1:B:413:ASP:O	1:B:416:ARG:HB2	2.01	0.61
1:B:398:LEU:HD13	1:B:404:PHE:CE1	2.36	0.60
1:B:293:GLU:HA	1:B:302:VAL:CG2	2.29	0.60
1:B:383:SER:HA	1:B:423:PHE:HD2	1.66	0.60
1:A:308:VAL:HG13	1:A:319:TYR:OH	2.01	0.60
1:A:308:VAL:HG22	1:A:319:TYR:CE2	2.36	0.60
1:A:277:TRP:CD1	1:A:289:THR:HG1	2.19	0.59
1:A:443:LEU:O	1:A:443:LEU:HD23	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:VAL:O	1:A:281:GLY:N	2.33	0.58
1:A:429:HIS:CG	1:A:430:GLU:N	2.72	0.58
1:B:310:HIS:O	1:B:311:GLN:C	2.47	0.58
1:B:383:SER:HB2	1:B:423:PHE:HE2	1.69	0.58
1:B:276:ASN:HB3	1:B:278:TYR:CE1	2.39	0.58
1:A:410:LEU:O	1:A:410:LEU:HG	2.03	0.57
1:B:308:VAL:HG13	1:B:319:TYR:CZ	2.39	0.57
2:C:2:NAG:H82	2:C:9:FUL:O3	2.04	0.57
1:B:354:SER:O	1:B:355:ARG:C	2.46	0.57
1:B:337:SER:O	1:B:338:LYS:C	2.47	0.56
1:A:418:GLN:C	1:A:420:GLY:N	2.59	0.56
1:B:258:GLU:HG3	3:D:6:GLA:H61	1.86	0.56
1:A:381:TRP:CH2	1:A:425:CYS:HB3	2.41	0.56
1:A:243:PHE:CD2	2:C:5:NAG:H62	2.41	0.56
1:A:346:PRO:HB3	1:A:372:PHE:CB	2.15	0.56
1:A:294:GLN:CB	1:A:296:TYR:HE1	2.18	0.55
1:B:261:CYS:HB2	1:B:277:TRP:CH2	2.41	0.55
1:B:288:LYS:O	1:B:289:THR:C	2.47	0.55
1:A:383:SER:HB2	1:A:423:PHE:CE2	2.41	0.55
1:A:259:VAL:HG23	1:A:308:VAL:CG2	2.37	0.55
1:B:390:ASN:O	1:B:410:LEU:HD12	2.07	0.55
1:A:338:LYS:HD3	1:A:376:ASP:OD1	2.07	0.54
1:A:284:VAL:O	1:A:287:ALA:HB2	2.06	0.54
1:A:249:ASP:HA	1:A:255:ARG:HD3	1.90	0.53
1:B:368:LEU:HB2	1:B:407:TYR:CE2	2.44	0.53
1:B:383:SER:HB2	1:B:423:PHE:CE2	2.43	0.53
1:B:240:VAL:O	1:B:241:PHE:CD2	2.61	0.53
1:B:240:VAL:O	1:B:241:PHE:HD2	1.90	0.53
1:B:429:HIS:CG	1:B:430:GLU:N	2.77	0.53
1:A:338:LYS:NZ	1:A:339:ALA:HB2	2.23	0.53
1:B:339:ALA:O	1:B:340:LYS:O	2.26	0.53
1:A:258:GLU:O	2:C:6:GLA:H61	2.09	0.52
1:A:310:HIS:O	1:A:311:GLN:C	2.52	0.52
1:B:418:GLN:C	1:B:420:GLY:N	2.63	0.52
1:B:346:PRO:HD3	1:B:429:HIS:HD2	1.75	0.52
1:B:295:GLN:C	1:B:296:TYR:HD1	2.17	0.52
1:B:312:ASN:HB3	1:B:317:LYS:HG3	1.91	0.52
1:A:432:LEU:O	1:A:433:HIS:C	2.52	0.51
1:A:354:SER:O	1:A:355:ARG:C	2.53	0.51
1:A:325:ASN:CG	1:A:326:LYS:H	2.18	0.51
1:A:366:THR:HG21	1:B:407:TYR:CE2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:278:TYR:CD1	1:B:278:TYR:N	2.79	0.51
1:A:338:LYS:O	1:A:339:ALA:O	2.29	0.51
1:B:241:PHE:CE1	3:D:2:NAG:H4	2.39	0.50
1:B:339:ALA:C	1:B:340:LYS:O	2.55	0.50
1:B:436:TYR:HD1	1:B:437:THR:N	2.08	0.50
2:C:2:NAG:C8	2:C:9:FUL:H3	2.40	0.50
1:B:250:THR:O	1:B:310:HIS:HD2	1.95	0.50
1:A:381:TRP:CZ3	1:A:425:CYS:HB3	2.47	0.50
1:B:375:SER:O	1:B:376:ASP:C	2.55	0.50
2:C:2:NAG:C8	2:C:9:FUL:C3	2.90	0.49
3:D:7:MAN:O3	3:D:8:NAG:C1	2.60	0.49
1:A:308:VAL:HG22	1:A:319:TYR:HE2	1.77	0.49
1:B:325:ASN:HD22	1:B:326:LYS:N	2.00	0.49
1:A:283:GLN:HG2	1:A:285:HIS:N	2.12	0.49
1:A:316:GLY:O	1:A:317:LYS:C	2.55	0.49
1:A:375:SER:O	1:A:377:ILE:HG22	2.13	0.49
1:B:316:GLY:O	1:B:317:LYS:C	2.55	0.49
1:B:436:TYR:HE1	1:B:438:GLN:HB2	1.78	0.48
1:A:288:LYS:O	1:A:289:THR:C	2.55	0.48
1:B:247:PRO:O	1:B:250:THR:OG1	2.31	0.48
1:B:339:ALA:O	1:B:340:LYS:C	2.55	0.48
1:B:306:LEU:HD12	1:B:307:THR:N	2.28	0.48
1:B:414:LYS:O	1:B:415:SER:C	2.56	0.48
1:B:243:PHE:HB2	1:B:260:THR:HB	1.95	0.48
1:B:330:ALA:O	1:B:331:PRO:C	2.56	0.48
1:B:416:ARG:HA	1:B:416:ARG:CZ	2.44	0.48
2:C:7:MAN:C4	2:C:8:NAG:H2	2.43	0.48
1:B:286:ASN:O	1:B:287:ALA:C	2.56	0.48
1:A:375:SER:O	1:A:376:ASP:C	2.56	0.48
1:B:283:GLN:CG	1:B:285:HIS:H	2.26	0.48
1:B:295:GLN:C	1:B:296:TYR:CD1	2.92	0.48
1:B:336:ILE:HG23	1:B:336:ILE:O	2.14	0.48
1:A:330:ALA:O	1:A:331:PRO:C	2.57	0.47
1:A:401:ASP:HB2	1:A:403:SER:CB	2.43	0.47
1:B:432:LEU:O	1:B:433:HIS:C	2.57	0.47
1:A:280:ASP:N	1:A:318:GLU:O	2.44	0.47
1:A:429:HIS:ND1	1:A:430:GLU:N	2.57	0.47
1:B:284:VAL:O	1:B:287:ALA:HB2	2.15	0.47
2:C:7:MAN:H4	2:C:8:NAG:C7	2.43	0.47
1:A:278:TYR:HA	1:A:283:GLN:HA	1.97	0.47
1:B:401:ASP:HB2	1:B:403:SER:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:328:LEU:CD1	1:B:329:PRO:HD2	2.43	0.47
1:B:391:TYR:C	1:B:391:TYR:CD2	2.92	0.47
1:A:286:ASN:O	1:A:287:ALA:C	2.57	0.47
1:A:295:GLN:C	1:A:296:TYR:CD1	2.93	0.47
1:A:341:GLY:O	1:A:343:PRO:HD3	2.14	0.47
1:B:382:GLU:HA	1:B:387:PRO:HA	1.97	0.47
1:B:424:SER:HB3	1:B:439:LYS:O	2.15	0.47
1:B:310:HIS:O	1:B:313:TRP:N	2.48	0.47
1:A:261:CYS:HB2	1:A:277:TRP:CH2	2.50	0.47
1:A:306:LEU:HD12	1:A:307:THR:N	2.29	0.47
1:A:325:ASN:CG	1:A:326:LYS:N	2.72	0.47
2:C:7:MAN:H4	2:C:8:NAG:C2	2.42	0.46
1:B:247:PRO:O	1:B:248:LYS:C	2.59	0.46
1:B:296:TYR:CE2	1:B:301:ARG:HD3	2.50	0.46
1:B:250:THR:O	1:B:310:HIS:CD2	2.68	0.46
1:B:294:GLN:CB	1:B:296:TYR:HE1	2.28	0.46
1:B:325:ASN:CG	1:B:326:LYS:N	2.74	0.46
1:B:243:PHE:N	1:B:260:THR:O	2.45	0.45
1:A:297:ASN:HB2	1:A:298:SER:H	1.58	0.45
1:B:286:ASN:O	1:B:288:LYS:N	2.49	0.45
1:A:286:ASN:O	1:A:288:LYS:N	2.50	0.45
1:A:295:GLN:C	1:A:296:TYR:HD1	2.25	0.45
1:B:326:LYS:HE2	1:B:326:LYS:HB2	1.77	0.45
1:A:338:LYS:HE3	1:A:338:LYS:C	2.42	0.45
2:C:4:MAN:O3	2:C:5:NAG:C1	2.65	0.45
1:B:283:GLN:HG2	1:B:285:HIS:N	2.28	0.45
1:B:354:SER:O	1:B:357:GLU:N	2.50	0.45
1:B:401:ASP:HB2	1:B:403:SER:HG	1.79	0.45
1:A:405:PHE:CE1	1:B:409:LYS:NZ	2.85	0.45
1:B:313:TRP:CH2	1:B:338:LYS:N	2.85	0.45
1:B:309:LEU:HD12	1:B:309:LEU:HA	1.83	0.44
2:C:7:MAN:O6	2:C:8:NAG:C8	2.63	0.44
1:B:328:LEU:HA	1:B:329:PRO:HD3	1.73	0.44
1:B:309:LEU:O	1:B:310:HIS:C	2.57	0.44
1:A:319:TYR:CD1	1:A:319:TYR:N	2.85	0.44
1:B:270:ASP:HB2	1:B:327:ALA:HB2	1.99	0.44
1:B:291:PRO:HD3	1:B:304:SER:OG	2.17	0.44
1:B:297:ASN:HB3	1:B:298:SER:H	1.28	0.44
1:B:301:ARG:O	1:B:301:ARG:HG2	2.17	0.44
2:C:2:NAG:C8	2:C:9:FUL:O3	2.65	0.44
1:A:338:LYS:HZ1	1:A:339:ALA:HB2	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:275:PHE:CE2	1:B:304:SER:HB3	2.53	0.43
1:A:247:PRO:O	1:A:248:LYS:C	2.60	0.43
1:A:259:VAL:CG2	1:A:308:VAL:HG21	2.42	0.43
1:A:345:GLU:O	1:A:346:PRO:C	2.61	0.43
1:B:355:ARG:O	1:B:358:MET:HG2	2.18	0.43
1:A:277:TRP:CD1	1:A:289:THR:CG2	3.02	0.43
1:B:243:PHE:CD1	3:D:5:NAG:H5	2.54	0.43
1:B:443:LEU:HD12	1:B:443:LEU:HA	1.55	0.43
1:A:255:ARG:NH1	1:A:255:ARG:HG2	2.33	0.43
1:B:241:PHE:CE1	3:D:3:BMA:C1	3.01	0.43
1:B:382:GLU:O	1:B:424:SER:N	2.47	0.43
1:B:341:GLY:O	1:B:342:GLN:C	2.61	0.43
1:A:391:TYR:HA	1:A:409:LYS:O	2.18	0.43
1:A:258:GLU:HG3	2:C:6:GLA:C6	2.46	0.43
1:A:309:LEU:O	1:A:310:HIS:C	2.62	0.43
1:B:429:HIS:CG	1:B:430:GLU:H	2.33	0.43
1:A:308:VAL:HG12	1:A:309:LEU:N	2.33	0.43
1:A:381:TRP:O	1:A:387:PRO:HA	2.18	0.43
1:A:442:SER:O	1:A:443:LEU:O	2.37	0.43
1:B:291:PRO:HB2	1:B:292:ARG:H	1.62	0.43
1:A:243:PHE:N	1:A:260:THR:O	2.52	0.42
1:A:391:TYR:HB2	1:A:409:LYS:O	2.19	0.42
1:B:397:VAL:HG12	1:B:398:LEU:N	2.34	0.42
1:B:391:TYR:HB2	1:B:409:LYS:O	2.19	0.42
1:B:433:HIS:HD1	1:B:434:ASN:CG	2.26	0.42
1:B:409:LYS:HZ2	1:B:409:LYS:CB	2.32	0.42
1:A:308:VAL:CG1	1:A:309:LEU:N	2.82	0.42
1:B:279:VAL:C	1:B:281:GLY:N	2.77	0.42
1:A:405:PHE:C	1:A:405:PHE:CD1	2.98	0.41
1:A:255:ARG:HG2	1:A:255:ARG:HH11	1.84	0.41
1:A:269:GLU:C	1:A:271:PRO:HD3	2.44	0.41
1:A:357:GLU:OE1	1:A:364:SER:N	2.52	0.41
1:B:391:TYR:HA	1:B:409:LYS:O	2.21	0.41
1:B:409:LYS:NZ	1:B:409:LYS:CB	2.83	0.41
1:A:328:LEU:HA	1:A:329:PRO:HD3	1.63	0.41
1:A:349:TYR:CE2	1:B:357:GLU:HG3	2.56	0.41
1:A:354:SER:O	1:A:357:GLU:N	2.54	0.41
1:B:372:PHE:CD1	1:B:372:PHE:C	2.99	0.41
1:B:341:GLY:O	1:B:343:PRO:N	2.54	0.40
1:B:371:GLY:HA2	1:B:403:SER:HB2	2.03	0.40
1:A:277:TRP:NE1	1:A:289:THR:CG2	2.80	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:309:LEU:CD1	1:B:309:LEU:CD1[9_765]	1.59	0.61

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	205/223 (92%)	165 (80%)	22 (11%)	18 (9%)	0	10
1	B	204/223 (92%)	159 (78%)	29 (14%)	16 (8%)	1	12
All	All	409/446 (92%)	324 (79%)	51 (12%)	34 (8%)	0	10

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	339	ALA
1	A	377	ILE
1	A	443	LEU
1	B	297	ASN
1	B	377	ILE
1	A	281	GLY
1	A	287	ALA
1	A	292	ARG
1	A	419	GLN
1	A	433	HIS
1	B	287	ALA
1	B	292	ARG
1	B	340	LYS
1	B	419	GLN
1	A	291	PRO
1	A	336	ILE
1	B	281	GLY

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Mol	Chain	Res	Type
1	B	291	PRO
1	B	311	GLN
1	A	297	ASN
1	A	310	HIS
1	A	311	GLN
1	A	331	PRO
1	A	384	ASN
1	B	384	ASN
1	A	285	HIS
1	A	329	PRO
1	A	414	LYS
1	B	285	HIS
1	B	310	HIS
1	B	329	PRO
1	B	336	ILE
1	B	331	PRO
1	B	433	HIS

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	108/206 (52%)	100 (93%)	8 (7%)	13	35
1	B	126/206 (61%)	114 (90%)	12 (10%)	8	27
All	All	234/412 (57%)	214 (92%)	20 (8%)	10	32

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	254	SER
1	A	280	ASP
1	A	306	LEU
1	A	312	ASN
1	A	338	LYS
1	A	400	SER

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Mol	Chain	Res	Type
1	A	425	CYS
1	A	443	LEU
1	B	260	THR
1	B	271	PRO
1	B	306	LEU
1	B	326	LYS
1	B	337	SER
1	B	359	THR
1	B	391	TYR
1	B	409	LYS
1	B	411	THR
1	B	416	ARG
1	B	425	CYS
1	B	442	SER

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	295	GLN
1	A	325	ASN
1	A	418	GLN
1	A	434	ASN
1	B	283	GLN
1	B	312	ASN
1	B	325	ASN
1	B	342	GLN
1	B	438	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 ⓘ

18 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	C	1	2,1	14,14,15	1.14	2 (14%)	17,19,21	1.51	3 (17%)
2	NAG	C	2	2	14,14,15	1.00	1 (7%)	17,19,21	1.52	3 (17%)
2	MAN	C	3	2	11,11,12	0.69	0	15,15,17	1.57	2 (13%)
2	MAN	C	4	2	11,11,12	1.01	1 (9%)	15,15,17	0.75	0
2	NAG	C	5	2	14,14,15	1.35	3 (21%)	17,19,21	2.14	5 (29%)
2	GLA	C	6	2	11,11,12	1.36	1 (9%)	15,15,17	0.94	0
2	MAN	C	7	2	11,11,12	0.91	0	15,15,17	1.21	2 (13%)
2	NAG	C	8	2	14,14,15	1.29	2 (14%)	17,19,21	1.21	2 (11%)
2	FUL	C	9	2	10,10,11	0.66	0	14,14,16	0.60	0
3	NAG	D	1	3,1	14,14,15	0.79	0	17,19,21	1.66	3 (17%)
3	NAG	D	2	3	14,14,15	0.75	0	17,19,21	1.28	2 (11%)
3	BMA	D	3	3	11,11,12	1.13	1 (9%)	15,15,17	0.70	0
3	MAN	D	4	3	11,11,12	1.05	1 (9%)	15,15,17	1.28	1 (6%)
3	NAG	D	5	3	14,14,15	0.98	0	17,19,21	1.75	4 (23%)
3	GLA	D	6	3	11,11,12	0.82	0	15,15,17	1.04	2 (13%)
3	MAN	D	7	3	11,11,12	1.09	1 (9%)	15,15,17	2.06	3 (20%)
3	NAG	D	8	3	14,14,15	0.89	1 (7%)	17,19,21	1.11	1 (5%)
3	FUL	D	9	3	10,10,11	0.97	0	14,14,16	0.74	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	C	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	C	2	2	-	3/6/23/26	0/1/1/1
2	MAN	C	3	2	-	2/2/19/22	0/1/1/1
2	MAN	C	4	2	-	2/2/19/22	0/1/1/1
2	NAG	C	5	2	1/1/5/7	3/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLA	C	6	2	-	1/2/19/22	0/1/1/1
2	MAN	C	7	2	-	2/2/19/22	0/1/1/1
2	NAG	C	8	2	-	2/6/23/26	0/1/1/1
2	FUL	C	9	2	-	-	0/1/1/1
3	NAG	D	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
3	BMA	D	3	3	-	2/2/19/22	0/1/1/1
3	MAN	D	4	3	-	2/2/19/22	0/1/1/1
3	NAG	D	5	3	-	2/6/23/26	0/1/1/1
3	GLA	D	6	3	-	2/2/19/22	0/1/1/1
3	MAN	D	7	3	-	0/2/19/22	0/1/1/1
3	NAG	D	8	3	-	4/6/23/26	0/1/1/1
3	FUL	D	9	3	-	-	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	5	NAG	C1-C2	3.34	1.56	1.52
2	C	8	NAG	C1-C2	3.31	1.56	1.52
3	D	3	BMA	C2-C3	3.18	1.57	1.52
3	D	7	MAN	C2-C3	3.07	1.57	1.52
2	C	6	GLA	C1-C2	2.76	1.58	1.52
3	D	4	MAN	C1-C2	2.68	1.58	1.52
2	C	1	NAG	C4-C5	2.51	1.58	1.53
2	C	5	NAG	O5-C5	2.40	1.48	1.43
2	C	5	NAG	O5-C1	2.35	1.47	1.43
2	C	1	NAG	C1-C2	-2.22	1.49	1.52
2	C	8	NAG	O5-C1	2.12	1.47	1.43
2	C	4	MAN	O2-C2	2.11	1.47	1.43
2	C	2	NAG	O4-C4	2.10	1.48	1.43
3	D	8	NAG	C1-C2	2.02	1.55	1.52

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	7	MAN	C1-C2-C3	6.28	118.79	109.64
2	C	5	NAG	C4-C3-C2	-5.56	102.86	111.02
2	C	3	MAN	C1-C2-C3	4.64	116.41	109.64
3	D	1	NAG	C2-N2-C7	-4.20	117.27	122.90
2	C	5	NAG	C1-O5-C5	3.79	117.26	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	5	NAG	C4-C3-C2	-3.74	105.53	111.02
3	D	1	NAG	C4-C3-C2	-3.74	105.53	111.02
3	D	7	MAN	C1-O5-C5	3.72	117.18	112.19
3	D	4	MAN	C1-O5-C5	3.50	116.87	112.19
2	C	5	NAG	C2-N2-C7	-3.37	118.39	122.90
3	D	2	NAG	C2-N2-C7	-3.36	118.40	122.90
2	C	1	NAG	C2-N2-C7	-3.29	118.50	122.90
2	C	2	NAG	C4-C3-C2	-3.24	106.27	111.02
2	C	7	MAN	C1-C2-C3	-3.21	104.97	109.64
3	D	5	NAG	C2-N2-C7	-3.17	118.65	122.90
3	D	5	NAG	O5-C1-C2	-3.15	106.42	111.29
2	C	2	NAG	C2-N2-C7	-3.12	118.72	122.90
3	D	1	NAG	C1-O5-C5	3.02	116.24	112.19
2	C	1	NAG	C3-C4-C5	2.87	115.44	110.23
2	C	5	NAG	C3-C4-C5	-2.84	105.08	110.23
2	C	8	NAG	C2-N2-C7	-2.76	119.20	122.90
3	D	8	NAG	C3-C4-C5	2.57	114.89	110.23
3	D	6	GLA	C1-C2-C3	2.56	113.37	109.64
2	C	3	MAN	C2-C3-C4	2.54	115.32	110.86
2	C	1	NAG	O5-C1-C2	-2.51	107.41	111.29
2	C	2	NAG	C1-O5-C5	2.45	115.47	112.19
2	C	7	MAN	C1-O5-C5	2.42	115.43	112.19
3	D	6	GLA	C1-O5-C5	2.28	115.24	112.19
3	D	5	NAG	C1-O5-C5	2.19	115.12	112.19
2	C	8	NAG	O7-C7-C8	-2.08	118.36	122.05
2	C	5	NAG	O4-C4-C5	2.07	114.43	109.32
3	D	7	MAN	C3-C4-C5	-2.04	106.53	110.23
3	D	2	NAG	O5-C1-C2	-2.04	108.13	111.29

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	5	NAG	C1

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
3	D	1	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	D	1	NAG	O7-C7-N2-C2
2	C	5	NAG	C8-C7-N2-C2
3	D	6	GLA	C4-C5-C6-O6
2	C	5	NAG	O7-C7-N2-C2
3	D	1	NAG	O5-C5-C6-O6
3	D	3	BMA	O5-C5-C6-O6
3	D	8	NAG	O5-C5-C6-O6
2	C	8	NAG	C4-C5-C6-O6
3	D	5	NAG	C4-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
3	D	3	BMA	C4-C5-C6-O6
3	D	6	GLA	O5-C5-C6-O6
2	C	3	MAN	C4-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
3	D	8	NAG	C4-C5-C6-O6
2	C	8	NAG	O5-C5-C6-O6
3	D	5	NAG	O5-C5-C6-O6
2	C	4	MAN	C4-C5-C6-O6
2	C	7	MAN	O5-C5-C6-O6
2	C	3	MAN	O5-C5-C6-O6
2	C	7	MAN	C4-C5-C6-O6
3	D	4	MAN	O5-C5-C6-O6
3	D	8	NAG	C8-C7-N2-C2
3	D	8	NAG	O7-C7-N2-C2
2	C	4	MAN	O5-C5-C6-O6
3	D	4	MAN	C4-C5-C6-O6
2	C	6	GLA	O5-C5-C6-O6
2	C	5	NAG	C1-C2-N2-C7
2	C	2	NAG	C4-C5-C6-O6

There are no ring outliers.

14 monomers are involved in 32 short contacts:

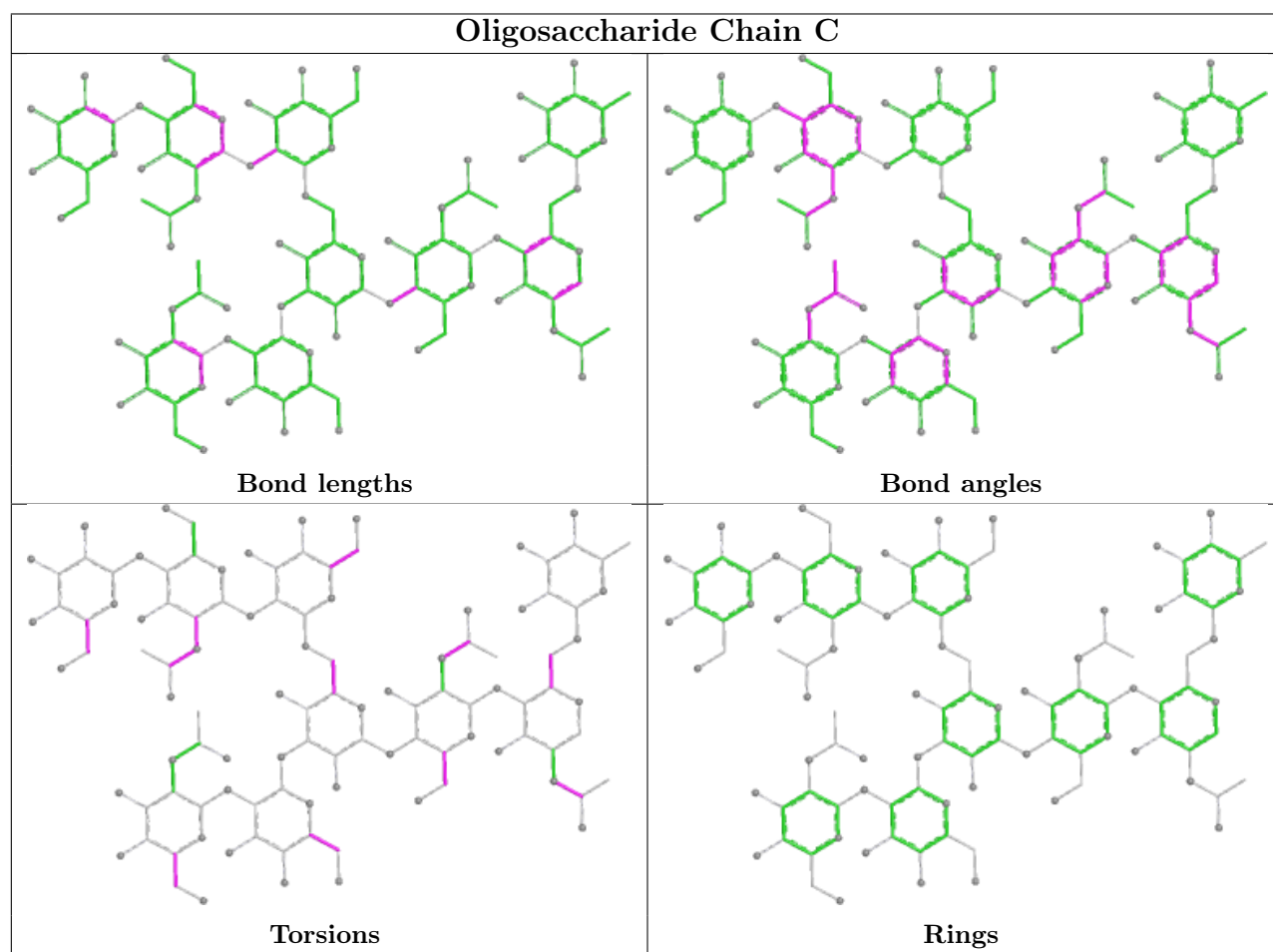
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	NAG	6	0
2	C	9	FUL	7	0
3	D	6	GLA	2	0
2	C	6	GLA	3	0
3	D	2	NAG	2	0
3	D	7	MAN	4	0
3	D	9	FUL	3	0

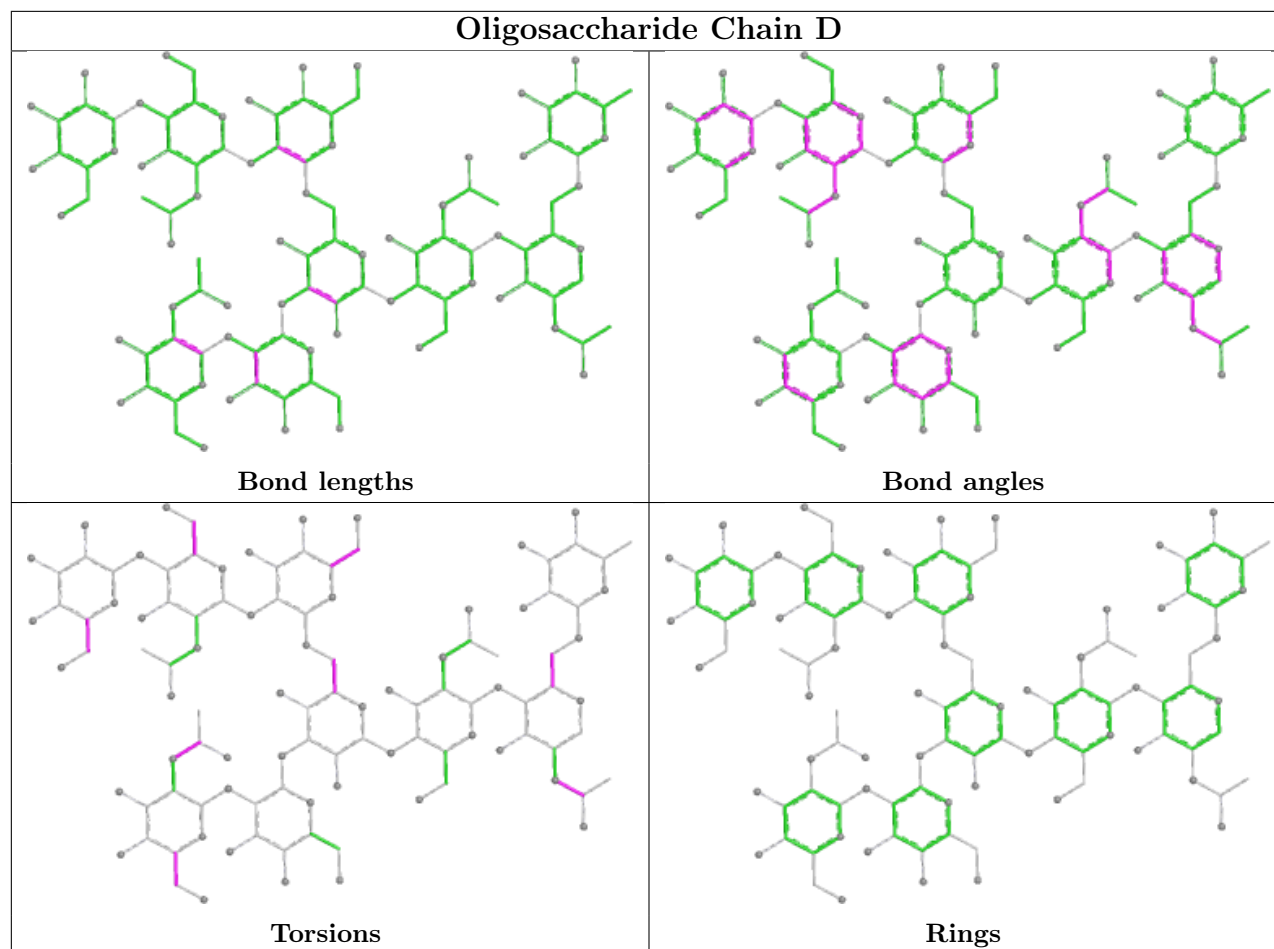
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	3	BMA	1	0
2	C	7	MAN	6	0
3	D	5	NAG	2	0
3	D	8	NAG	4	0
2	C	4	MAN	2	0
2	C	5	NAG	3	0
2	C	8	NAG	9	0

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	A	207/223 (92%)	0.60	5 (2%) 59 45	28, 32, 47, 47	0
1	B	206/223 (92%)	0.38	3 (1%) 72 54	14, 17, 22, 23	0
All	All	413/446 (92%)	0.49	8 (1%) 66 50	14, 23, 46, 47	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	254	SER	2.6
1	A	330	ALA	2.4
1	A	397	VAL	2.4
1	B	280	ASP	2.4
1	A	403	SER	2.3
1	B	407	TYR	2.2
1	A	371	GLY	2.1
1	A	434	ASN	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	C	8	14/15	0.73	0.16	71,73,75,76	0

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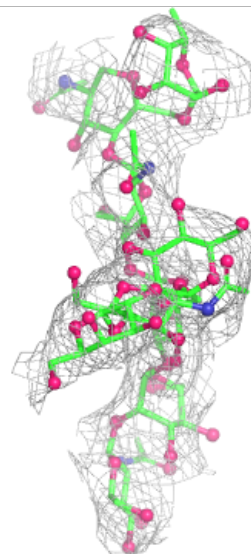
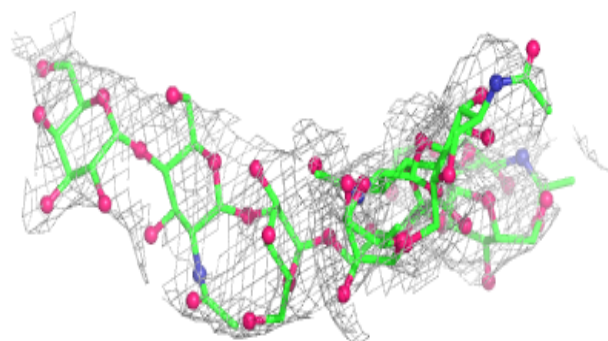
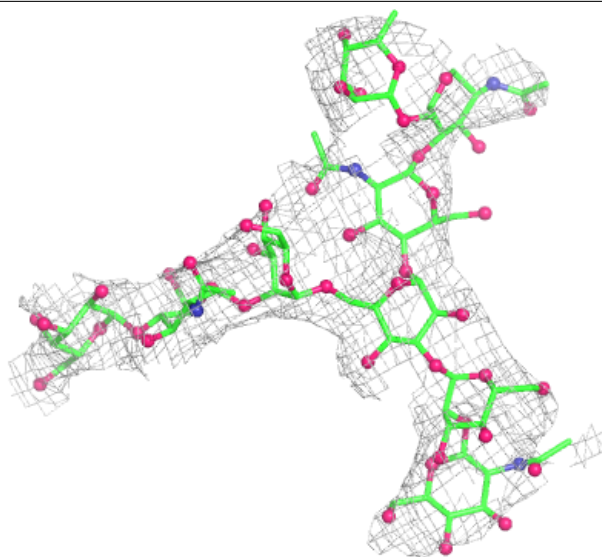
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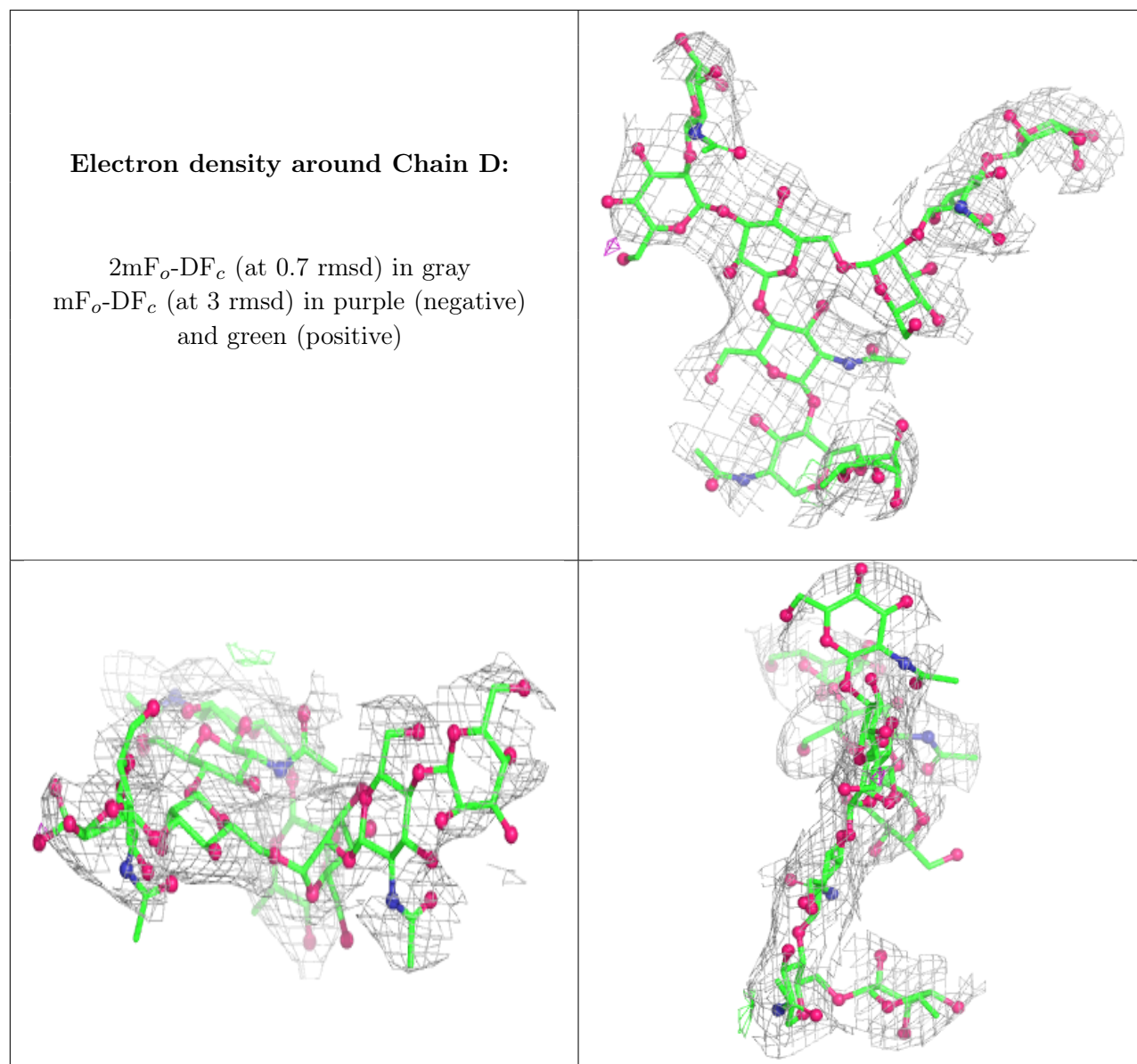
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MAN	D	7	11/12	0.73	0.19	33,35,38,38	0
2	NAG	C	1	14/15	0.75	0.20	71,73,76,76	0
2	NAG	C	2	14/15	0.78	0.18	71,74,75,76	0
2	FUL	C	9	10/11	0.81	0.15	70,73,75,76	0
2	MAN	C	3	11/12	0.81	0.17	71,73,76,76	0
3	NAG	D	8	14/15	0.81	0.16	33,36,39,39	0
2	GLA	C	6	11/12	0.82	0.16	71,72,76,76	0
3	GLA	D	6	11/12	0.82	0.18	33,34,38,39	0
3	NAG	D	1	14/15	0.83	0.16	33,35,39,39	0
3	MAN	D	4	11/12	0.85	0.16	33,35,38,38	0
2	MAN	C	4	11/12	0.85	0.15	71,73,76,76	0
3	NAG	D	5	14/15	0.87	0.13	33,35,38,38	0
2	MAN	C	7	11/12	0.87	0.15	73,75,76,76	0
3	BMA	D	3	11/12	0.88	0.13	33,36,38,39	0
2	NAG	C	5	14/15	0.89	0.12	71,74,76,76	0
3	FUL	D	9	10/11	0.89	0.13	33,36,38,39	0
3	NAG	D	2	14/15	0.92	0.13	33,35,38,39	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.

Electron density around Chain C:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.