



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:19 AM UTC

PDB ID : 1H3V / pdb_00001h3v
Title : CRYSTAL STRUCTURE OF THE HUMAN IGG1 FC-FRAGMENT, GLYC
OFORM (G2F)2, SG P212121
Authors : Krapp, S.; Mimura, Y.; Jefferis, R.; Huber, R.; Sonderrmann, P.
Deposited on : 2002-09-19
Resolution : 3.10 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

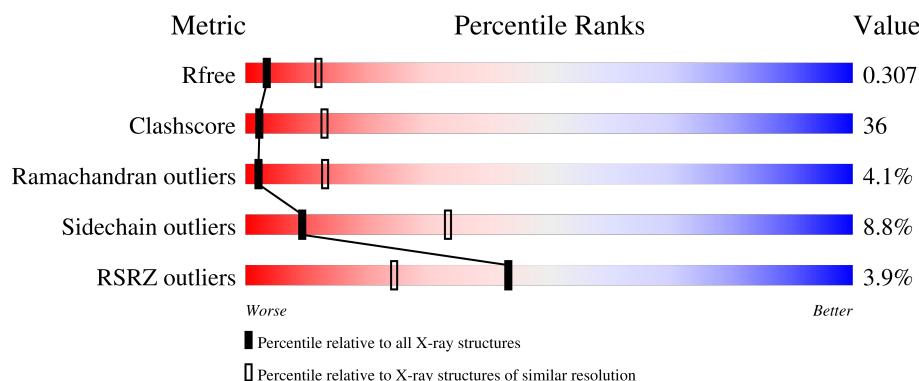
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	
1	B	223	
2	C	8	
3	D	8	

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	GAL	C	6	-	-	X	-
3	NAG	D	1	X	-	-	-

2 Entry composition [i](#)

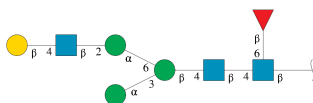
There are 4 unique types of molecules in this entry. The entry contains 3600 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

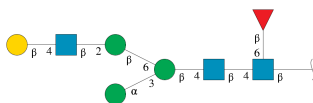
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	45	0	1
			1661	1057	283	314	7			
1	B	207	Total	C	N	O	S	131	0	1
			1656	1054	282	313	7			

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[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
2	C	8	Total	C	N	O	0	0	0
			96	54	3	39			

- Molecule 3 is an oligosaccharide called beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-beta-D-mannopyranose-(1-6)-[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	8	Total	C	N	O	15	0	0
			96	54	3	39			

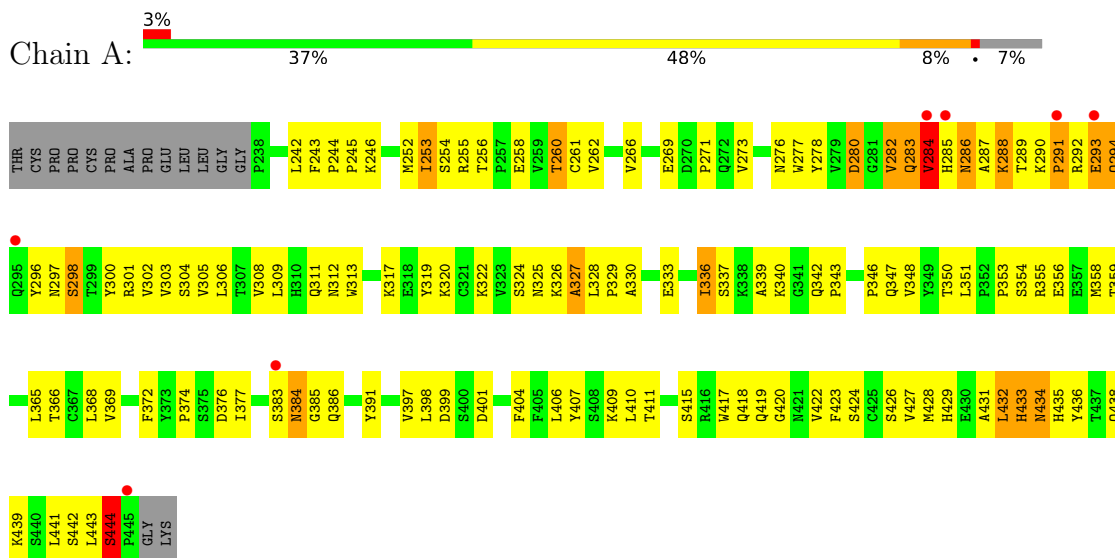
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	51	Total 51	O 51	0	0
4	B	40	Total 40	O 40	0	0

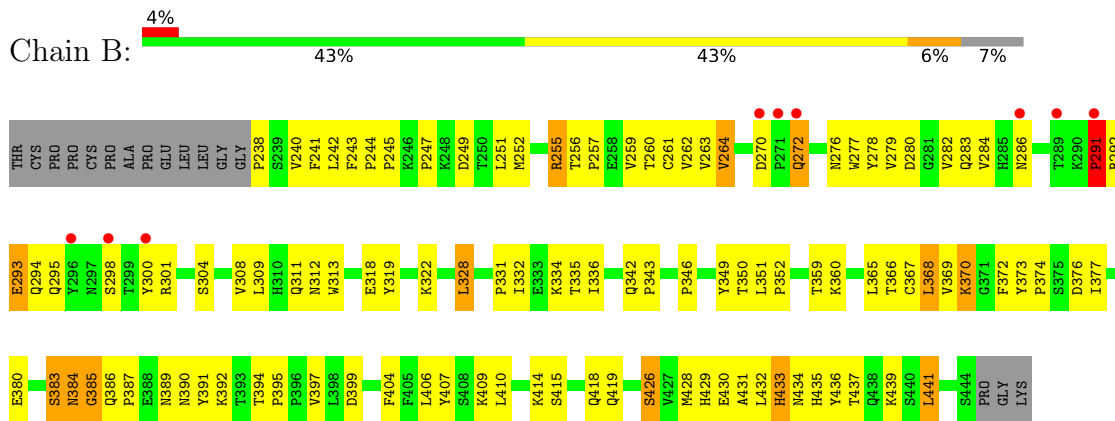
3 Residue-property plots

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.

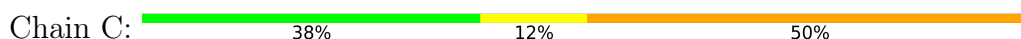
• Molecule 1: IG GAMMA-1 CHAIN C REGION



• Molecule 1: IG GAMMA-1 CHAIN C REGION



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



MAG1	MAG2	EMA3	MAN4	MAG5	GAL6	MAN7	FUL8
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- Molecule 3: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-beta-D-mannopyranose-(1-6)-[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:



MAG1	MAG2	EMA3	MAN4	MAG5	GAL6	MAN7	FUL8
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.47Å 80.85Å 141.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.10 50.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	82.8 (50.00-3.10) 82.6 (50.00-3.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 3.17Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.240 , 0.314 0.233 , 0.307	Depositor DCC
R_{free} test set	445 reflections (5.44%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 59.9	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	0.88	EDS
Total number of atoms	3600	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.82% 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: MAN, NAG, FUL, GAL, BMA

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	0.58	1/1707 (0.1%)	1.04	7/2325 (0.3%)
1	B	0.53	0/1702	1.04	10/2318 (0.4%)
All	All	0.56	1/3409 (0.0%)	1.04	17/4643 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	444	SER	C-N	-5.45	1.25	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	284	VAL	N-CA-C	-9.59	101.52	110.53
1	A	383	SER	N-CA-C	-8.85	95.85	109.65
1	B	383	SER	N-CA-C	-7.91	98.30	109.69
1	A	293	GLU	N-CA-C	7.11	119.16	110.41
1	A	296	TYR	N-CA-C	6.91	118.50	108.34
1	B	395	PRO	N-CA-C	-6.30	103.01	110.70
1	B	394	THR	CA-C-N	-5.96	114.24	120.38
1	B	394	THR	C-N-CA	-5.96	114.24	120.38
1	A	359	THR	N-CA-C	-5.40	106.46	112.57
1	B	433	HIS	N-CA-C	-5.40	99.29	110.80
1	B	342	GLN	CA-C-N	5.27	125.17	119.85
1	B	342	GLN	C-N-CA	5.27	125.17	119.85
1	A	284	VAL	N-CA-C	-5.25	105.42	110.72
1	B	293	GLU	N-CA-C	5.25	117.30	110.43
1	B	256	THR	N-CA-C	5.20	118.05	108.69
1	A	432	LEU	N-CA-C	-5.12	102.58	109.95
1	A	327	ALA	N-CA-C	-5.10	107.22	112.93

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	1661	0	1630	129	0
1	B	1656	0	1628	96	0
2	C	96	0	82	16	0
3	D	96	0	82	11	0
4	A	51	0	0	14	0
4	B	40	0	0	7	0
All	All	3600	0	3422	236	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 36.

All (236) 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)
3:D:1:NAG:H62	3:D:8:FUL:C6	1.82	1.10
1:A:351:LEU:HB2	1:A:366:THR:HB	1.33	1.10
1:A:433:HIS:O	1:A:435:HIS:N	1.85	1.07
1:B:433:HIS:O	1:B:435:HIS:N	1.88	1.07
3:D:1:NAG:C6	3:D:8:FUL:H63	1.89	1.01
1:B:238:PRO:HB2	4:B:2001:HOH:O	1.60	1.01
1:A:350:THR:HB	1:A:441:LEU:HG	1.46	0.97
1:B:365:LEU:HD12	1:B:410:LEU:HD23	1.44	0.96
3:D:1:NAG:H61	3:D:8:FUL:O3	1.65	0.96
1:A:253:ILE:HD12	1:A:253:ILE:H	1.29	0.94
1:B:346:PRO:HB3	1:B:372:PHE:HB3	1.55	0.88
3:D:1:NAG:H62	3:D:8:FUL:H63	0.92	0.86
1:A:292:ARG:HG2	1:A:293:GLU:H	1.41	0.85
1:B:351:LEU:HB2	1:B:366:THR:HB	1.59	0.84
1:B:260:THR:HG21	4:B:2040:HOH:O	1.79	0.82
1:A:309:LEU:HB2	1:A:312:ASN:HD22	1.43	0.82
1:A:320:LYS:HD3	1:A:333:GLU:HG2	1.61	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:350:THR:HB	1:B:441:LEU:HG	1.60	0.81
1:A:260:THR:HG23	1:A:305:VAL:HG22	1.65	0.77
1:B:312:ASN:HB3	1:B:319:TYR:OH	1.83	0.77
1:A:276:ASN:HB2	1:A:322:LYS:HB3	1.66	0.76
1:A:253:ILE:H	1:A:253:ILE:CD1	1.97	0.76
1:B:249:ASP:O	1:B:257:PRO:HG3	1.87	0.75
1:A:244:PRO:HB3	1:A:336:ILE:HD11	1.68	0.74
1:A:433:HIS:C	1:A:435:HIS:H	1.93	0.74
1:A:245:PRO:HA	2:C:6:GAL:H61	1.69	0.74
1:A:253:ILE:HD12	1:A:253:ILE:N	2.03	0.73
1:A:292:ARG:HG2	1:A:293:GLU:N	2.01	0.73
1:A:297:ASN:O	1:A:298:SER:HB2	1.88	0.72
1:A:346:PRO:HB3	1:A:372:PHE:HB3	1.71	0.71
1:B:272:GLN:OE1	1:B:272:GLN:HA	1.90	0.71
1:B:238:PRO:HB2	1:B:328:LEU:HD11	1.71	0.70
1:A:245:PRO:HA	2:C:6:GAL:C6	2.21	0.70
1:A:391:TYR:HB3	1:A:410:LEU:HD12	1.75	0.69
1:A:273:VAL:HG21	1:A:302:VAL:HG21	1.74	0.68
1:A:423:PHE:HB2	4:A:2047:HOH:O	1.92	0.67
1:A:358:MET:HE1	1:A:417:TRP:CD1	2.28	0.67
1:A:353:PRO:HD3	1:A:365:LEU:HD23	1.77	0.67
1:A:309:LEU:HB2	1:A:312:ASN:ND2	2.08	0.66
1:B:359:THR:O	1:B:359:THR:HG22	1.94	0.66
1:B:261:CYS:HB2	1:B:277:TRP:CZ2	2.31	0.65
2:C:2:NAG:H81	2:C:4:MAN:O4	1.96	0.65
4:A:2048:HOH:O	2:C:1:NAG:H82	1.97	0.65
1:B:311:GLN:OE1	1:B:311:GLN:N	2.28	0.64
2:C:5:NAG:H4	2:C:6:GAL:O2	1.96	0.64
1:B:291:PRO:HB3	1:B:304:SER:HA	1.79	0.64
1:B:278:TYR:HA	1:B:282:VAL:O	1.97	0.64
1:A:284:VAL:HG13	1:A:285:HIS:ND1	2.13	0.64
1:B:261:CYS:HB2	1:B:277:TRP:CH2	2.34	0.62
1:B:241:PHE:HE1	3:D:2:NAG:H4	1.64	0.62
1:A:441:LEU:CB	4:A:2047:HOH:O	2.47	0.62
1:B:433:HIS:C	1:B:435:HIS:H	2.07	0.62
1:B:432:LEU:C	1:B:433:HIS:O	2.38	0.61
1:B:360:LYS:HB3	4:B:2021:HOH:O	1.99	0.61
1:B:247:PRO:O	1:B:251:LEU:HG	2.01	0.60
1:B:241:PHE:CE1	3:D:2:NAG:H4	2.36	0.60
1:A:398:LEU:HD12	1:A:399:ASP:N	2.16	0.60
2:C:1:NAG:H61	2:C:2:NAG:HN2	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:332:ILE:HG23	1:B:334:LYS:NZ	2.17	0.59
1:B:383:SER:O	1:B:384:ASN:C	2.46	0.59
1:B:383:SER:O	1:B:385:GLY:N	2.35	0.59
1:A:242:LEU:HD13	1:A:336:ILE:HG22	1.83	0.59
1:B:247:PRO:HB3	1:B:376:ASP:HB3	1.84	0.59
1:A:289:THR:HG22	1:A:290:LYS:H	1.68	0.59
1:B:259:VAL:HG23	1:B:308:VAL:HG21	1.85	0.58
1:B:335:THR:HG22	1:B:336:ILE:N	2.18	0.58
1:A:266:VAL:HB	1:A:300:TYR:HB2	1.84	0.58
1:A:301:ARG:HH12	1:A:303:VAL:HG21	1.69	0.58
1:A:252:MET:HE3	1:A:428:MET:HE1	1.86	0.57
1:B:276:ASN:HB2	1:B:322:LYS:HB3	1.85	0.57
1:A:436:TYR:CE1	1:A:438:GLN:HG3	2.40	0.57
1:A:325:ASN:HD21	1:A:327:ALA:HB3	1.70	0.57
1:A:347:GLN:NE2	1:B:360:LYS:HE3	2.19	0.57
1:A:289:THR:HG22	1:A:290:LYS:N	2.20	0.57
1:A:391:TYR:HB3	1:A:410:LEU:CD1	2.35	0.56
1:B:346:PRO:HB3	1:B:372:PHE:CB	2.31	0.56
3:D:1:NAG:C6	3:D:8:FUL:O3	2.48	0.56
1:A:325:ASN:ND2	1:A:327:ALA:HB3	2.21	0.55
1:A:280:ASP:OD2	1:A:317:LYS:HD3	2.07	0.55
1:B:346:PRO:CB	1:B:372:PHE:HB3	2.32	0.54
1:A:441:LEU:O	4:A:2047:HOH:O	2.18	0.54
1:B:238:PRO:CB	1:B:328:LEU:HD11	2.38	0.54
1:A:432:LEU:C	1:A:433:HIS:O	2.50	0.54
1:B:243:PHE:O	1:B:259:VAL:HG13	2.07	0.54
1:B:415:SER:O	1:B:419:GLN:N	2.40	0.54
1:A:353:PRO:HD3	1:A:365:LEU:CD2	2.37	0.53
1:B:373:TYR:CD1	1:B:374:PRO:HA	2.44	0.53
1:A:273:VAL:CG2	1:A:302:VAL:HG21	2.38	0.53
1:A:443:LEU:O	1:A:444:SER:CB	2.56	0.53
1:A:287:ALA:N	4:A:2008:HOH:O	2.39	0.53
1:B:259:VAL:HG23	1:B:308:VAL:CG2	2.38	0.53
1:B:368:LEU:HD11	1:B:370:LYS:HB3	1.91	0.53
1:B:433:HIS:C	1:B:435:HIS:N	2.66	0.53
1:A:311:GLN:HA	1:A:311:GLN:NE2	2.24	0.52
2:C:1:NAG:H61	2:C:2:NAG:N2	2.24	0.52
1:B:244:PRO:HD3	1:B:336:ILE:HD11	1.90	0.52
1:B:429:HIS:HB3	1:B:432:LEU:HG	1.91	0.52
1:A:258:GLU:O	2:C:6:GAL:H61	2.10	0.52
1:B:309:LEU:O	1:B:312:ASN:HB2	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:GLU:HG3	1:A:294:GLN:H	1.75	0.51
1:A:312:ASN:O	1:A:313:TRP:C	2.52	0.51
1:B:279:VAL:O	1:B:282:VAL:HG22	2.10	0.51
1:B:252:MET:SD	1:B:428:MET:HE1	2.51	0.51
1:A:271:PRO:O	1:A:273:VAL:HG23	2.10	0.51
1:B:322:LYS:NZ	1:B:331:PRO:HG3	2.26	0.51
1:A:258:GLU:O	2:C:6:GAL:C6	2.59	0.51
1:A:308:VAL:HG12	1:A:319:TYR:CE2	2.46	0.51
4:A:2048:HOH:O	2:C:1:NAG:H2	2.10	0.51
1:B:349:TYR:O	1:B:367:CYS:HA	2.10	0.51
1:A:427:VAL:O	1:A:436:TYR:HA	2.11	0.51
1:B:397:VAL:O	1:B:404:PHE:HA	2.10	0.50
1:A:252:MET:HB2	1:A:255:ARG:HD3	1.92	0.50
1:B:415:SER:O	1:B:419:GLN:HG3	2.10	0.50
1:A:242:LEU:HD22	1:A:336:ILE:HG21	1.93	0.50
1:B:308:VAL:HG12	1:B:309:LEU:N	2.25	0.50
1:A:293:GLU:HG3	1:A:294:GLN:N	2.27	0.50
1:A:406:LEU:HD12	1:A:406:LEU:C	2.36	0.50
1:B:332:ILE:HG23	1:B:334:LYS:HZ2	1.77	0.50
1:A:301:ARG:HG2	1:A:301:ARG:HH11	1.77	0.49
1:A:278:TYR:HA	1:A:282:VAL:O	2.12	0.49
1:A:384:ASN:C	1:A:386:GLN:H	2.20	0.49
1:B:368:LEU:C	1:B:368:LEU:HD12	2.36	0.49
1:B:376:ASP:HB2	4:B:2031:HOH:O	2.13	0.49
1:A:415:SER:O	1:A:419:GLN:N	2.45	0.49
1:A:432:LEU:O	1:A:433:HIS:O	2.30	0.49
1:A:422:VAL:HA	1:A:442:SER:HG	1.77	0.49
1:A:429:HIS:CD2	1:A:431:ALA:H	2.31	0.49
1:A:434:ASN:O	1:A:435:HIS:HB2	2.12	0.49
1:A:283:GLN:NE2	4:A:2008:HOH:O	2.29	0.49
1:A:397:VAL:O	1:A:404:PHE:HA	2.13	0.49
1:B:414:LYS:O	1:B:418:GLN:CG	2.61	0.48
1:A:426:SER:HB3	1:A:436:TYR:CE2	2.48	0.48
1:B:308:VAL:HG13	1:B:319:TYR:OH	2.13	0.48
1:A:278:TYR:HB2	1:A:320:LYS:HB3	1.94	0.48
1:A:340:LYS:HA	4:A:2025:HOH:O	2.14	0.48
1:A:358:MET:HE1	1:A:417:TRP:NE1	2.26	0.48
1:A:266:VAL:O	1:A:300:TYR:HB2	2.14	0.48
1:A:336:ILE:HG13	1:A:337:SER:N	2.28	0.48
3:D:5:NAG:H4	3:D:6:GAL:O2	2.14	0.48
1:B:311:GLN:H	1:B:311:GLN:CD	2.17	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:ARG:HG3	1:B:292:ARG:HH11	1.79	0.48
1:B:359:THR:O	1:B:360:LYS:HD3	2.14	0.47
2:C:5:NAG:C4	2:C:6:GAL:O2	2.62	0.47
1:A:301:ARG:NH1	1:A:303:VAL:HG21	2.29	0.47
1:B:391:TYR:O	1:B:392:LYS:HD3	2.14	0.47
1:B:283:GLN:HG2	4:B:2007:HOH:O	2.13	0.47
1:B:380:GLU:HG2	4:B:2023:HOH:O	2.14	0.47
1:A:288:LYS:HD3	1:A:306:LEU:HD11	1.97	0.47
1:A:399:ASP:OD2	1:B:409:LYS:NZ	2.47	0.47
1:B:391:TYR:HA	1:B:409:LYS:O	2.15	0.47
1:B:439:LYS:HA	1:B:439:LYS:HD2	1.72	0.47
1:B:406:LEU:C	1:B:406:LEU:HD12	2.40	0.47
1:A:245:PRO:HA	2:C:6:GAL:O6	2.15	0.46
1:A:328:LEU:HG	1:A:330:ALA:O	2.15	0.46
1:B:245:PRO:HD3	1:B:259:VAL:HG22	1.97	0.46
1:A:311:GLN:NE2	1:A:311:GLN:CA	2.77	0.46
1:A:284:VAL:HG13	1:A:285:HIS:HD1	1.81	0.46
1:A:365:LEU:HD21	1:A:417:TRP:CE3	2.51	0.46
1:B:251:LEU:HD13	1:B:429:HIS:C	2.41	0.46
1:A:368:LEU:HD13	1:A:407:TYR:CZ	2.51	0.45
1:B:247:PRO:HG2	1:B:377:ILE:H	1.81	0.45
1:A:255:ARG:HG2	1:A:255:ARG:HH11	1.80	0.45
1:A:391:TYR:HA	1:A:409:LYS:O	2.17	0.45
1:A:261:CYS:HB2	1:A:277:TRP:CH2	2.52	0.45
1:A:293:GLU:OE2	4:A:2013:HOH:O	2.21	0.45
1:A:409:LYS:HB2	1:B:407:TYR:OH	2.16	0.45
1:A:418:GLN:C	1:A:420:GLY:H	2.25	0.45
1:B:322:LYS:HZ1	1:B:331:PRO:HG3	1.81	0.45
1:A:246:LYS:HE2	2:C:6:GAL:H3	1.99	0.45
1:B:426:SER:HB3	1:B:436:TYR:CE2	2.52	0.45
1:A:368:LEU:HD13	1:A:407:TYR:CE1	2.52	0.45
1:A:285:HIS:O	1:A:286:ASN:O	2.35	0.45
1:B:335:THR:CG2	1:B:336:ILE:N	2.80	0.45
1:A:243:PHE:CD2	2:C:5:NAG:H5	2.52	0.44
1:B:292:ARG:HH11	1:B:293:GLU:H	1.65	0.44
1:A:423:PHE:CB	4:A:2047:HOH:O	2.57	0.44
1:A:441:LEU:HB2	4:A:2047:HOH:O	2.15	0.44
1:B:264:VAL:HG11	3:D:1:NAG:O4	2.17	0.44
1:B:368:LEU:HD12	1:B:369:VAL:N	2.31	0.44
1:B:429:HIS:O	1:B:435:HIS:HA	2.17	0.44
1:A:246:LYS:HE2	2:C:6:GAL:H5	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:THR:OG1	3:D:5:NAG:H61	2.18	0.44
1:B:365:LEU:HB2	1:B:410:LEU:HB3	1.99	0.44
1:A:423:PHE:N	1:A:423:PHE:CD1	2.85	0.44
1:B:441:LEU:HD23	4:B:2020:HOH:O	2.17	0.44
1:A:288:LYS:CD	1:A:306:LEU:HD11	2.48	0.44
1:A:384:ASN:O	1:A:386:GLN:N	2.50	0.44
1:B:263:VAL:O	1:B:301:ARG:HA	2.18	0.44
1:A:369:VAL:HG12	1:A:372:PHE:CD2	2.53	0.44
1:B:430:GLU:CG	1:B:431:ALA:N	2.81	0.44
1:A:286:ASN:HA	4:A:2008:HOH:O	2.18	0.43
1:A:354:SER:O	1:A:355:ARG:C	2.60	0.43
1:A:376:ASP:O	1:A:377:ILE:HB	2.18	0.43
1:B:308:VAL:HG11	1:B:313:TRP:HB2	2.00	0.43
1:B:292:ARG:HG3	1:B:292:ARG:NH1	2.34	0.43
1:B:312:ASN:HB3	1:B:319:TYR:HH	1.78	0.43
1:A:417:TRP:CZ2	1:A:443:LEU:HB2	2.53	0.43
1:A:436:TYR:CD1	1:A:436:TYR:C	2.96	0.43
1:A:429:HIS:O	1:A:435:HIS:HA	2.19	0.43
1:B:240:VAL:HG23	1:B:332:ILE:HD12	2.01	0.43
1:B:309:LEU:HG	1:B:311:GLN:OE1	2.19	0.43
1:A:277:TRP:NE1	1:A:289:THR:HG21	2.34	0.42
1:A:435:HIS:O	1:A:436:TYR:HB2	2.19	0.42
1:B:262:VAL:HG21	3:D:5:NAG:H62	2.00	0.42
1:A:339:ALA:HB3	1:A:374:PRO:HB3	2.00	0.42
1:A:242:LEU:HD13	1:A:336:ILE:CG2	2.48	0.42
1:A:291:PRO:HB3	1:A:304:SER:HA	2.00	0.42
1:A:348:VAL:HG22	1:A:427:VAL:HG21	2.01	0.42
1:A:284:VAL:HG13	1:A:285:HIS:CE1	2.55	0.42
1:A:328:LEU:HA	1:A:329:PRO:HD3	1.93	0.42
1:A:297:ASN:O	1:A:298:SER:CB	2.62	0.41
1:B:383:SER:O	1:B:386:GLN:N	2.51	0.41
1:A:428:MET:HG2	1:A:436:TYR:HD2	1.86	0.41
1:A:433:HIS:C	1:A:435:HIS:N	2.60	0.41
1:B:255:ARG:CZ	1:B:255:ARG:HB3	2.50	0.41
1:B:384:ASN:C	1:B:386:GLN:H	2.28	0.41
1:A:276:ASN:HB3	1:A:278:TYR:HE1	1.86	0.41
1:A:301:ARG:HG2	1:A:301:ARG:NH1	2.35	0.41
1:A:342:GLN:HA	1:A:343:PRO:HD3	1.89	0.41
1:A:384:ASN:C	1:A:386:GLN:N	2.79	0.41
1:B:352:PRO:HG3	1:B:441:LEU:HD21	2.03	0.41
1:A:398:LEU:HD12	1:A:398:LEU:C	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:4:MAN:O2	2:C:5:NAG:H82	2.21	0.41
1:A:297:ASN:HB3	4:A:2049:HOH:O	2.21	0.41
1:B:244:PRO:CD	1:B:336:ILE:HD11	2.51	0.41
1:B:308:VAL:CG1	1:B:309:LEU:N	2.84	0.41
1:A:424:SER:HA	1:A:439:LYS:O	2.21	0.41
1:A:292:ARG:CG	1:A:293:GLU:N	2.80	0.40
1:B:322:LYS:HZ2	1:B:331:PRO:CG	2.33	0.40
1:A:326:LYS:C	1:A:328:LEU:H	2.29	0.40
1:B:343:PRO:HA	1:B:373:TYR:O	2.22	0.40
1:A:266:VAL:HB	1:A:300:TYR:CB	2.49	0.40
1:A:439:LYS:HA	1:A:439:LYS:HD3	1.76	0.40
1:A:441:LEU:HB3	4:A:2047:HOH:O	2.18	0.40
1:A:243:PHE:HE2	1:A:262:VAL:HG21	1.86	0.40
1:B:311:GLN:O	1:B:312:ASN:C	2.65	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	206/223 (92%)	171 (83%)	26 (13%)	9 (4%)	2	12
1	B	205/223 (92%)	176 (86%)	21 (10%)	8 (4%)	2	14
All	All	411/446 (92%)	347 (84%)	47 (11%)	17 (4%)	2	13

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	286	ASN
1	A	298	SER
1	A	384	ASN
1	A	433	HIS

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Mol	Chain	Res	Type
1	A	434	ASN
1	B	286	ASN
1	B	298	SER
1	B	384	ASN
1	B	434	ASN
1	A	385	GLY
1	A	444	SER
1	A	291	PRO
1	B	291	PRO
1	B	270	ASP
1	A	356	GLU
1	B	387	PRO
1	B	385	GLY

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	193/206 (94%)	178 (92%)	15 (8%)	11	38
1	B	193/206 (94%)	174 (90%)	19 (10%)	7	29
All	All	386/412 (94%)	352 (91%)	34 (9%)	9	33

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	253	ILE
1	A	254	SER
1	A	256	THR
1	A	260	THR
1	A	269	GLU
1	A	280	ASP
1	A	282	VAL
1	A	283	GLN
1	A	284	VAL
1	A	288	LYS

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Mol	Chain	Res	Type
1	A	294	GLN
1	A	324	SER
1	A	336	ILE
1	A	401	ASP
1	A	411	THR
1	B	242	LEU
1	B	255	ARG
1	B	264	VAL
1	B	272	GLN
1	B	280	ASP
1	B	291	PRO
1	B	294	GLN
1	B	295	GLN
1	B	300	TYR
1	B	318	GLU
1	B	328	LEU
1	B	368	LEU
1	B	370	LYS
1	B	389	ASN
1	B	390	ASN
1	B	399	ASP
1	B	426	SER
1	B	437	THR
1	B	441	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	HIS
1	A	276	ASN
1	A	311	GLN
1	A	312	ASN
1	A	325	ASN
1	A	361	ASN
1	A	362	GLN
1	A	389	ASN
1	A	418	GLN
1	A	419	GLN
1	A	421	ASN
1	A	429	HIS
1	A	433	HIS
1	A	438	GLN

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Mol	Chain	Res	Type
1	B	347	GLN
1	B	361	ASN
1	B	390	ASN
1	B	418	GLN
1	B	419	GLN
1	B	429	HIS
1	B	434	ASN
1	B	435	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

16 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	0.63	0	17,19,21	0.76	1 (5%)
2	NAG	C	2	2	14,14,15	0.62	0	17,19,21	0.76	1 (5%)
2	BMA	C	3	2	11,11,12	0.49	0	15,15,17	0.50	0
2	MAN	C	4	2	11,11,12	0.36	0	15,15,17	0.91	1 (6%)
2	NAG	C	5	2	14,14,15	0.61	0	17,19,21	0.87	1 (5%)
2	GAL	C	6	2	11,11,12	0.42	0	15,15,17	0.45	0
2	MAN	C	7	2	11,11,12	0.47	0	15,15,17	0.45	0
2	FUL	C	8	2	10,10,11	0.81	0	14,14,16	0.55	0
3	NAG	D	1	3,1	14,14,15	0.87	0	17,19,21	0.86	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	D	2	3	14,14,15	0.68	0	17,19,21	0.73	0
3	BMA	D	3	3	11,11,12	0.77	0	15,15,17	0.92	1 (6%)
3	BMA	D	4	3	11,11,12	0.63	0	15,15,17	0.84	1 (6%)
3	NAG	D	5	3	14,14,15	0.51	0	17,19,21	0.68	0
3	GAL	D	6	3	11,11,12	0.52	0	15,15,17	0.52	0
3	MAN	D	7	3	11,11,12	0.54	0	15,15,17	0.77	1 (6%)
3	FUL	D	8	3	10,10,11	0.68	0	14,14,16	1.51	2 (14%)

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	-	5/6/23/26	0/1/1/1
2	NAG	C	2	2	-	3/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
2	MAN	C	4	2	-	2/2/19/22	0/1/1/1
2	NAG	C	5	2	-	2/6/23/26	0/1/1/1
2	GAL	C	6	2	-	2/2/19/22	0/1/1/1
2	MAN	C	7	2	-	0/2/19/22	0/1/1/1
2	FUL	C	8	2	-	-	0/1/1/1
3	NAG	D	1	3,1	1/1/5/7	4/6/23/26	0/1/1/1
3	NAG	D	2	3	-	4/6/23/26	0/1/1/1
3	BMA	D	3	3	-	2/2/19/22	0/1/1/1
3	BMA	D	4	3	-	2/2/19/22	0/1/1/1
3	NAG	D	5	3	-	4/6/23/26	0/1/1/1
3	GAL	D	6	3	-	1/2/19/22	0/1/1/1
3	MAN	D	7	3	-	0/2/19/22	0/1/1/1
3	FUL	D	8	3	-	-	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	8	FUL	C3-C4-C5	3.62	115.32	109.81
3	D	3	BMA	C1-C2-C3	2.87	113.83	109.64
3	D	1	NAG	C2-N2-C7	-2.58	119.45	122.90
2	C	4	MAN	C1-O5-C5	2.46	115.49	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	7	MAN	C1-O5-C5	2.45	115.47	112.19
2	C	5	NAG	C2-N2-C7	-2.36	119.74	122.90
2	C	2	NAG	C1-C2-N2	-2.29	106.83	110.43
3	D	8	FUL	O5-C5-C4	2.26	113.62	109.55
2	C	1	NAG	C2-N2-C7	-2.12	120.06	122.90
3	D	4	BMA	C1-C2-C3	-2.05	106.66	109.64

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	D	1	NAG	C1

All (31) 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	O7-C7-N2-C2
2	C	5	NAG	C8-C7-N2-C2
2	C	5	NAG	O7-C7-N2-C2
3	D	1	NAG	C3-C2-N2-C7
3	D	1	NAG	C8-C7-N2-C2
3	D	1	NAG	O7-C7-N2-C2
3	D	2	NAG	C8-C7-N2-C2
3	D	2	NAG	O7-C7-N2-C2
2	C	2	NAG	C8-C7-N2-C2
2	C	1	NAG	O5-C5-C6-O6
3	D	4	BMA	O5-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
3	D	5	NAG	C8-C7-N2-C2
3	D	5	NAG	O7-C7-N2-C2
3	D	4	BMA	C4-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	C	4	MAN	O5-C5-C6-O6
3	D	5	NAG	C4-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
2	C	4	MAN	C4-C5-C6-O6
2	C	6	GAL	O5-C5-C6-O6
3	D	5	NAG	O5-C5-C6-O6
3	D	6	GAL	O5-C5-C6-O6
3	D	3	BMA	C4-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	D	3	BMA	O5-C5-C6-O6
2	C	1	NAG	C3-C2-N2-C7
2	C	2	NAG	C3-C2-N2-C7
2	C	6	GAL	C4-C5-C6-O6

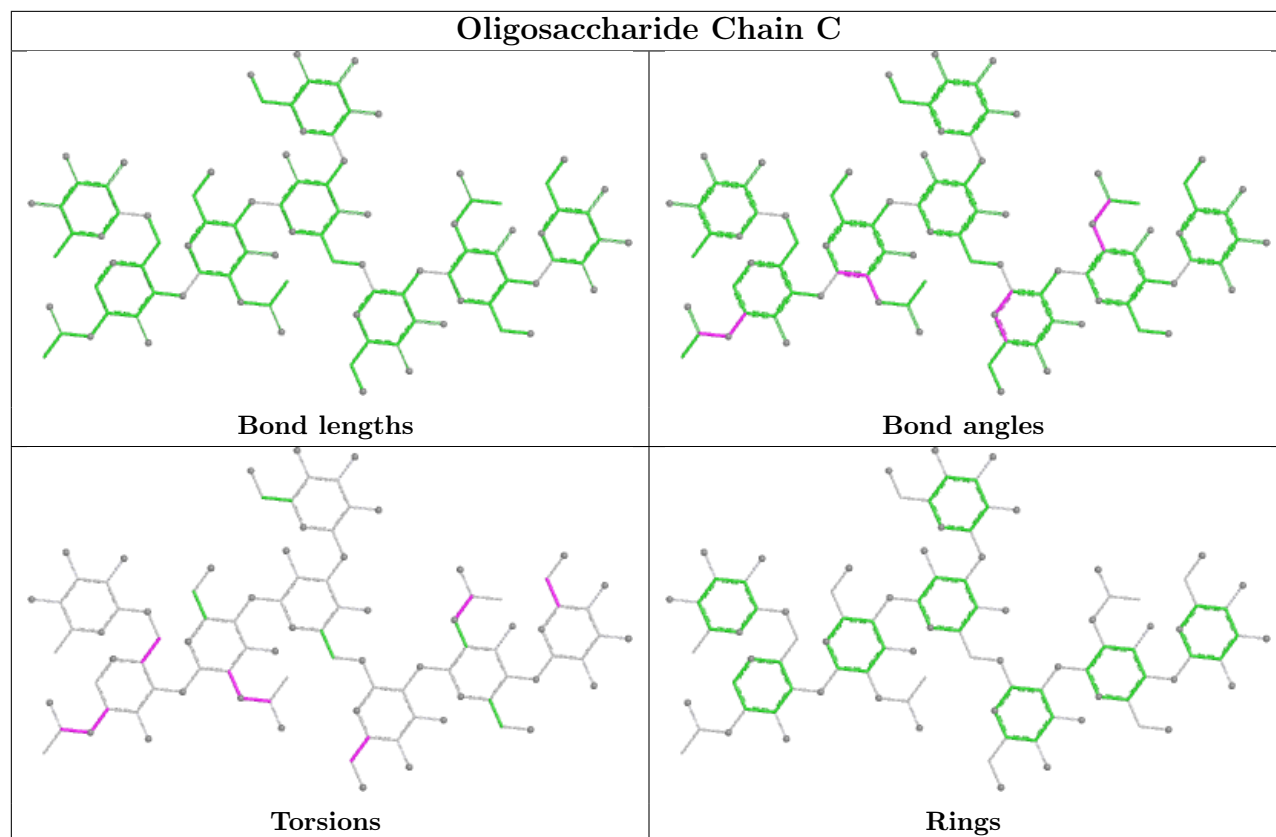
There are no ring outliers.

10 monomers are involved in 27 short contacts:

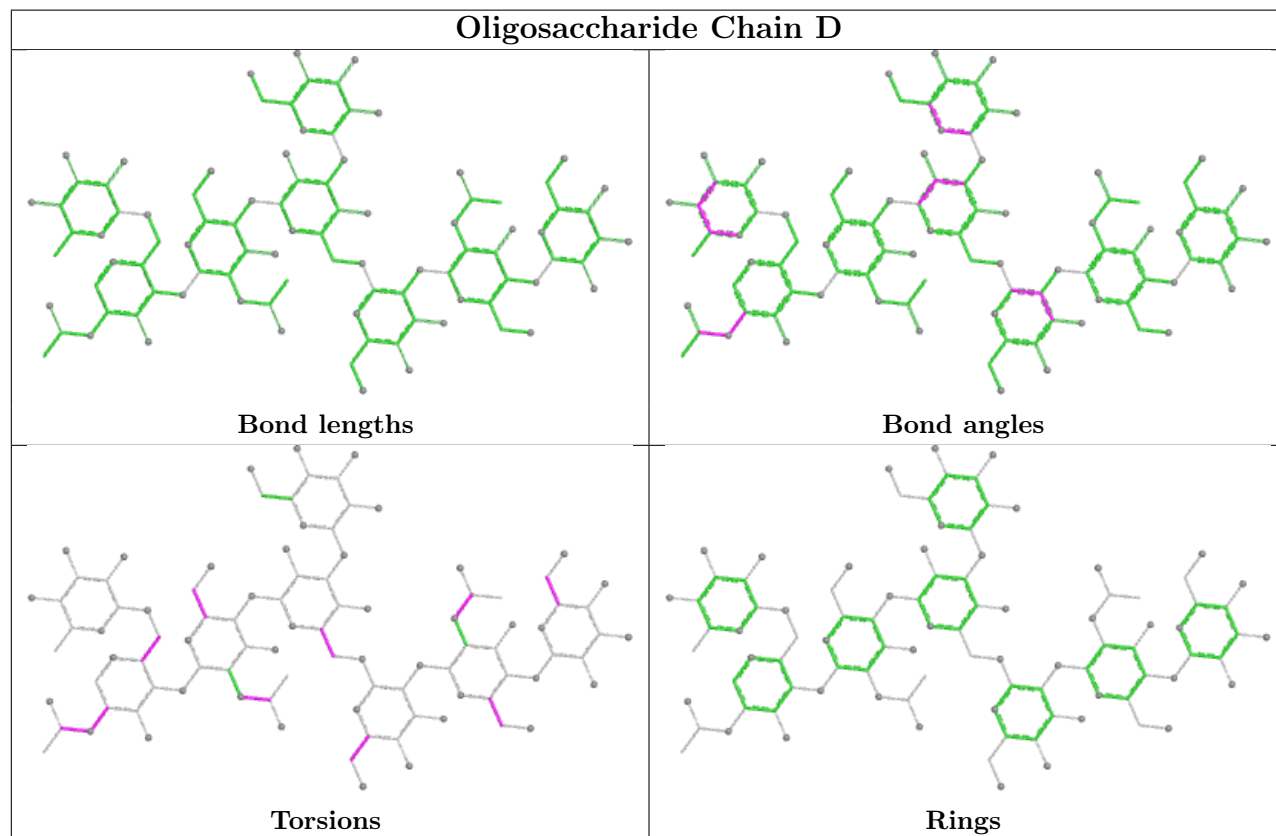
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	4	MAN	2	0
2	C	2	NAG	3	0
2	C	1	NAG	4	0
2	C	5	NAG	4	0
3	D	2	NAG	2	0
3	D	5	NAG	3	0
3	D	6	GAL	1	0
2	C	6	GAL	9	0
3	D	8	FUL	5	0
3	D	1	NAG	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain C



Oligosaccharide Chain D



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	208/223 (93%)	0.02	7 (3%)	48	28	2, 21, 50, 65	15 (7%)
1	B	205/223 (91%)	0.28	9 (4%)	39	21	2, 28, 64, 87	37 (18%)
All	All	413/446 (92%)	0.15	16 (3%)	43	24	2, 24, 57, 87	52 (12%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	286	ASN	4.5
1	B	291	PRO	4.1
1	B	296	TYR	3.3
1	B	270	ASP	2.8
1	A	383	SER	2.8
1	A	285	HIS	2.7
1	A	295	GLN	2.7
1	B	300	TYR	2.6
1	A	291	PRO	2.6
1	B	298	SER	2.5
1	A	284	VAL	2.3
1	A	293	GLU	2.2
1	B	289	THR	2.1
1	B	271	PRO	2.1
1	B	272	GLN	2.0
1	A	445	PRO	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

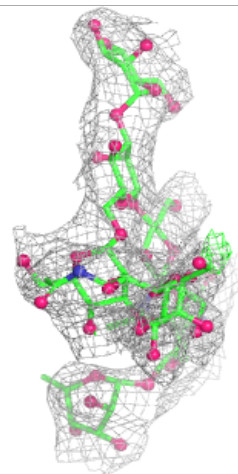
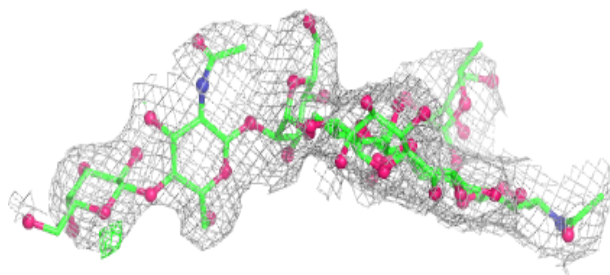
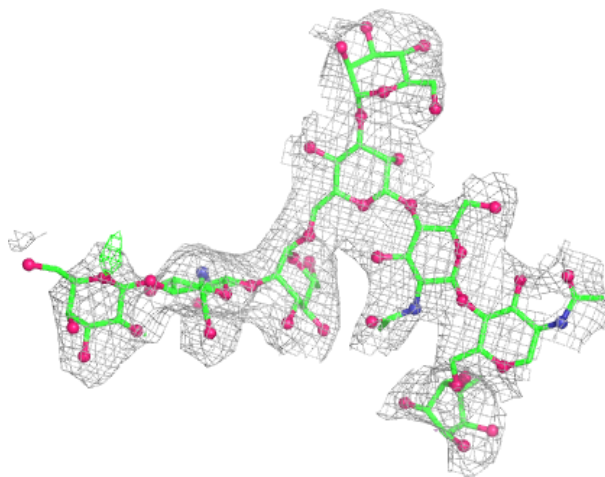
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

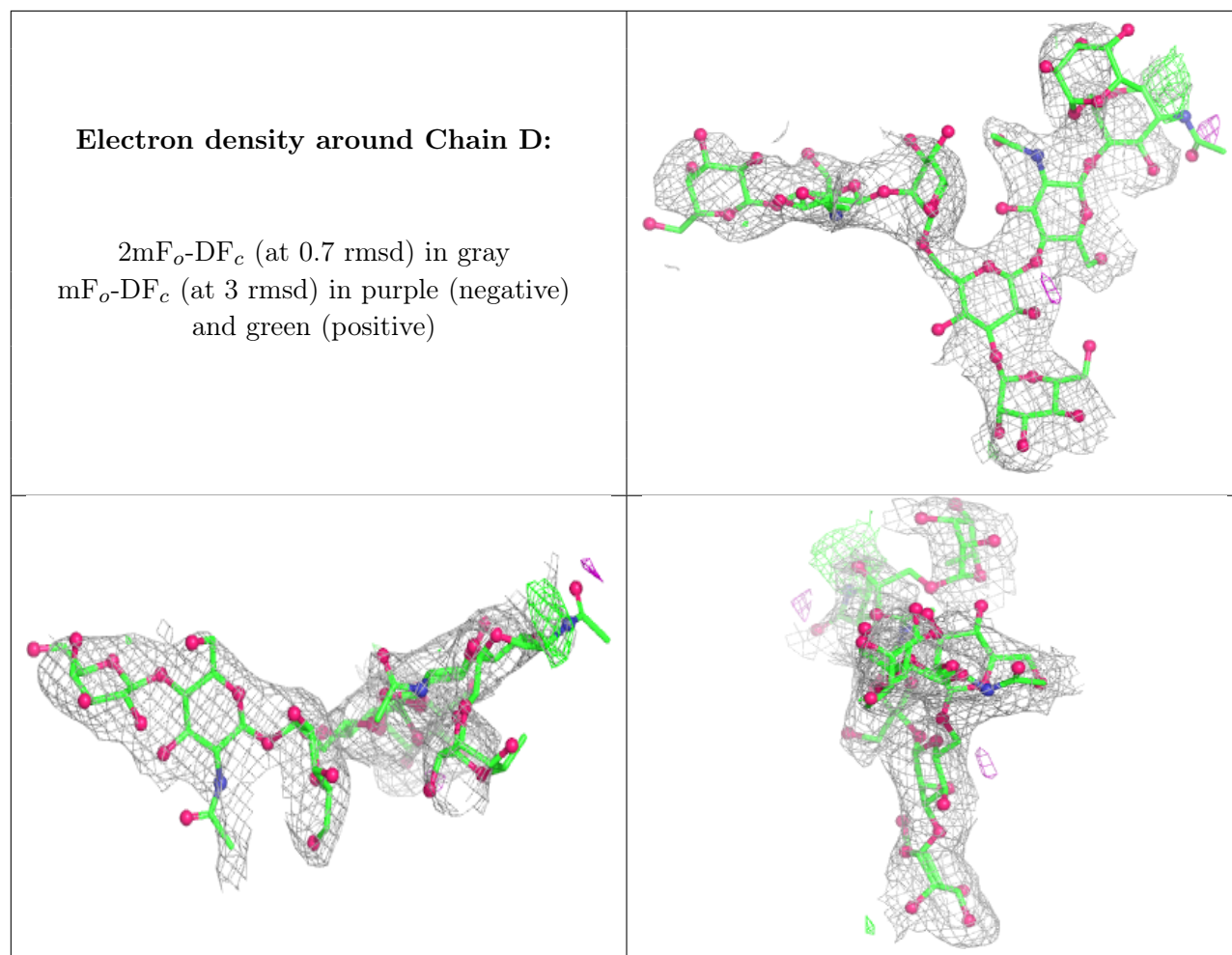
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	D	1	14/15	0.60	0.20	88,93,98,102	6
3	FUL	D	8	10/11	0.71	0.28	106,109,110,110	0
3	MAN	D	7	11/12	0.73	0.15	71,73,75,76	0
2	FUL	C	8	10/11	0.80	0.14	58,61,61,63	0
3	BMA	D	3	11/12	0.81	0.14	72,75,76,76	1
2	GAL	C	6	11/12	0.83	0.20	52,53,56,59	0
2	NAG	C	5	14/15	0.84	0.12	42,47,52,55	0
3	GAL	D	6	11/12	0.84	0.18	66,68,69,71	2
3	NAG	D	2	14/15	0.85	0.16	78,82,84,85	1
3	BMA	D	4	11/12	0.86	0.12	65,68,69,69	2
2	NAG	C	1	14/15	0.86	0.10	42,44,50,56	0
2	MAN	C	7	11/12	0.87	0.09	54,58,60,63	0
3	NAG	D	5	14/15	0.91	0.12	63,64,65,66	3
2	MAN	C	4	11/12	0.91	0.07	37,38,39,41	0
2	NAG	C	2	14/15	0.92	0.09	39,41,42,42	0
2	BMA	C	3	11/12	0.93	0.07	37,42,45,48	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.