



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:30 PM UTC

PDB ID : 3HA0 / pdb_00003ha0
Title : Crystal structure of the IgE-Fc3-4 domains
Authors : Wurzburg, B.A.
Deposited on : 2009-04-30
Resolution : 2.80 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

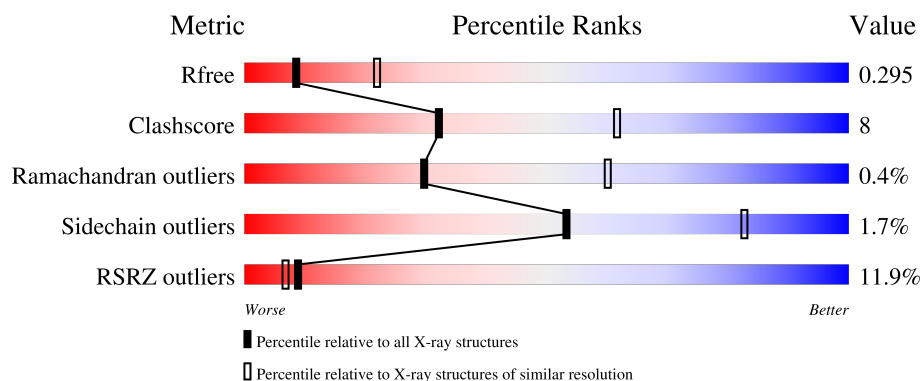
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	3% 82% 12% 6%
1	B	223	4% 80% 13% 6%
1	C	223	10% 77% 17% 6%
1	D	223	9% 75% 18% 6%
1	E	223	27% 72% 20% 6%

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Mol	Chain	Length	Quality of chain
1	F	223	13% 77% 16% • 6%
2	G	5	20% 80%
2	H	5	60% 40%
2	I	5	40% 60%
2	J	5	40% 60%
2	K	5	20% 40% 40%
2	L	5	60% 40%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10276 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 epsilon chain C region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	0	0
			1640	1027	299	308	6			
1	B	209	Total	C	N	O	S	0	0	0
			1656	1037	305	308	6			
1	C	209	Total	C	N	O	S	0	0	0
			1656	1037	305	308	6			
1	D	209	Total	C	N	O	S	0	0	0
			1656	1037	305	308	6			
1	E	209	Total	C	N	O	S	0	0	0
			1656	1037	305	308	6			
1	F	209	Total	C	N	O	S	0	0	0
			1646	1030	302	308	6			

There are 30 discrepancies between the modelled and reference sequences:

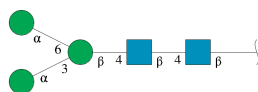
Chain	Residue	Modelled	Actual	Comment	Reference
A	325	ALA	-	expression tag	UNP P01854
A	326	ASP	-	expression tag	UNP P01854
A	327	PRO	-	expression tag	UNP P01854
A	371	GLN	ASN	engineered mutation	UNP P01854
A	383	GLN	ASN	engineered mutation	UNP P01854
B	325	ALA	-	expression tag	UNP P01854
B	326	ASP	-	expression tag	UNP P01854
B	327	PRO	-	expression tag	UNP P01854
B	371	GLN	ASN	engineered mutation	UNP P01854
B	383	GLN	ASN	engineered mutation	UNP P01854
C	325	ALA	-	expression tag	UNP P01854
C	326	ASP	-	expression tag	UNP P01854
C	327	PRO	-	expression tag	UNP P01854
C	371	GLN	ASN	engineered mutation	UNP P01854
C	383	GLN	ASN	engineered mutation	UNP P01854
D	325	ALA	-	expression tag	UNP P01854
D	326	ASP	-	expression tag	UNP P01854

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Chain	Residue	Modelled	Actual	Comment	Reference
D	327	PRO	-	expression tag	UNP P01854
D	371	GLN	ASN	engineered mutation	UNP P01854
D	383	GLN	ASN	engineered mutation	UNP P01854
E	325	ALA	-	expression tag	UNP P01854
E	326	ASP	-	expression tag	UNP P01854
E	327	PRO	-	expression tag	UNP P01854
E	371	GLN	ASN	engineered mutation	UNP P01854
E	383	GLN	ASN	engineered mutation	UNP P01854
F	325	ALA	-	expression tag	UNP P01854
F	326	ASP	-	expression tag	UNP P01854
F	327	PRO	-	expression tag	UNP P01854
F	371	GLN	ASN	engineered mutation	UNP P01854
F	383	GLN	ASN	engineered mutation	UNP P01854

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

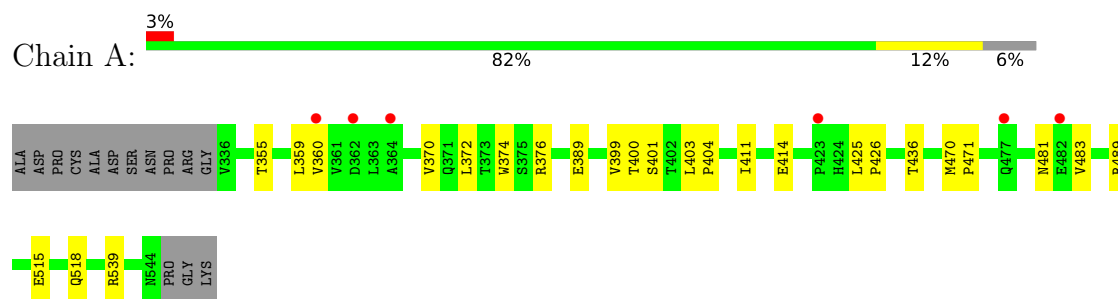


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	H	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	I	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	J	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	K	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	L	5	Total	C	N	O	0	0	0
			61	34	2	25			

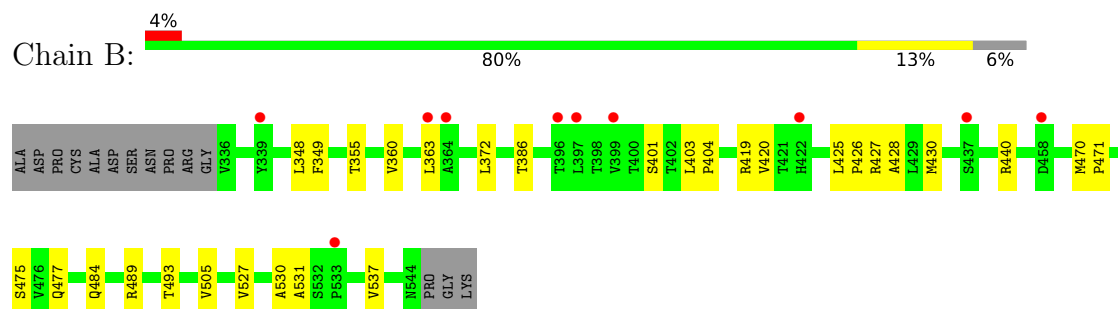
3 Residue-property plots [i](#)

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.

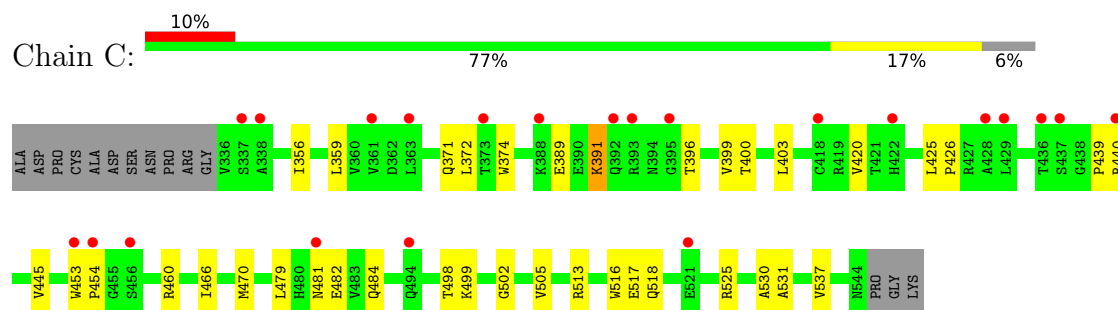
- Molecule 1: Ig epsilon chain C region



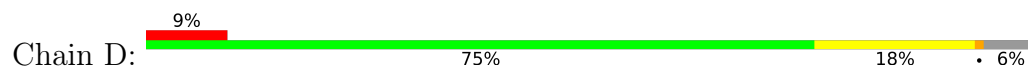
- Molecule 1: Ig epsilon chain C region

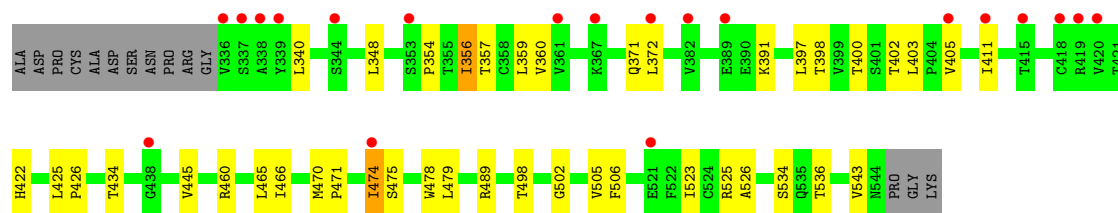


- Molecule 1: Ig epsilon chain C region

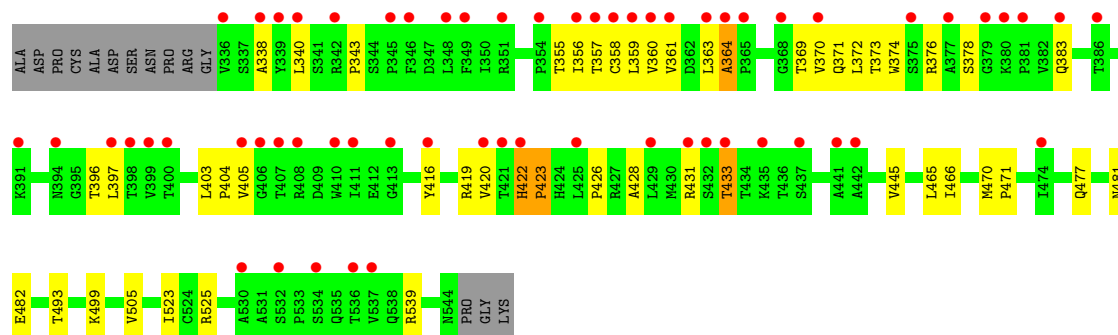
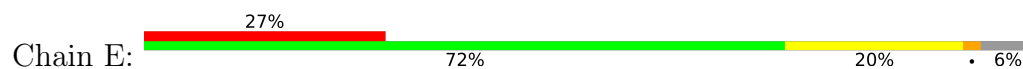


- Molecule 1: Ig epsilon chain C region

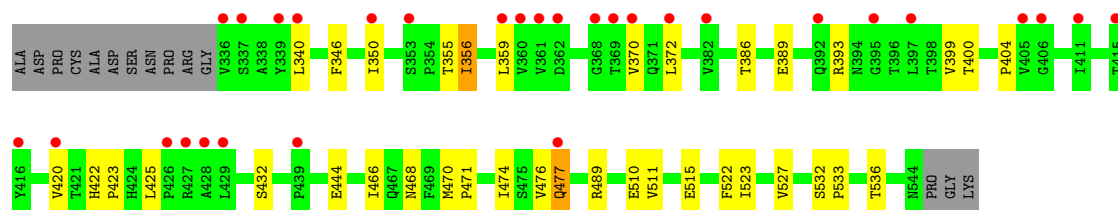
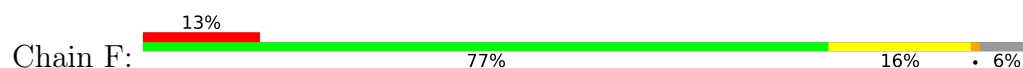




• Molecule 1: Ig epsilon chain C region



• Molecule 1: Ig epsilon chain C region



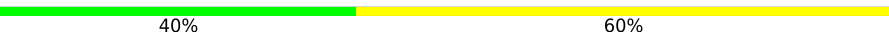
• Molecule 2: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

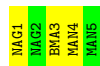


• Molecule 2: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

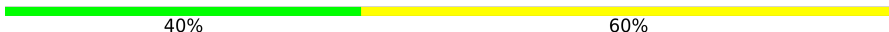


• Molecule 2: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain I: 

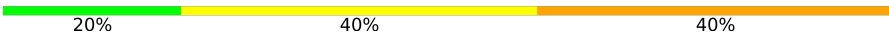


• Molecule 2: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain J: 



• Molecule 2: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain K: 



• Molecule 2: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain L: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.90Å 104.90Å 150.00Å 90.00° 96.20° 90.00°	Depositor
Resolution (Å)	29.83 – 2.80 29.83 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.83-2.80) 99.4 (29.83-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.80Å)	Xtriage
Refinement program	REFMAC, CNS	Depositor
R, R_{free}	0.241 , 0.280 0.255 , 0.295	Depositor DCC
R_{free} test set	1860 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	61.8	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	10276	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.16 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.5892e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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, NAG, 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	0.47	0/1682	0.68	0/2291
1	B	0.44	0/1698	0.68	0/2311
1	C	0.40	0/1698	0.68	0/2311
1	D	0.42	0/1698	0.70	0/2311
1	E	0.35	0/1698	0.65	0/2311
1	F	0.40	0/1688	0.63	0/2298
All	All	0.41	0/10162	0.67	0/13833

There are no bond length outliers.

There are no bond angle outliers.

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	1640	0	1609	19	0
1	B	1656	0	1641	19	0
1	C	1656	0	1641	30	0
1	D	1656	0	1641	33	0
1	E	1656	0	1643	35	0
1	F	1646	0	1620	34	0
2	G	61	0	52	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	61	0	52	2	0
2	I	61	0	52	1	0
2	J	61	0	52	0	0
2	K	61	0	52	4	0
2	L	61	0	52	0	0
All	All	10276	0	10107	162	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 (162) 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:C:372:LEU:HD22	1:C:420:VAL:HG22	1.41	1.01
1:A:359:LEU:HD13	1:A:400:THR:HG22	1.55	0.89
1:B:363:LEU:HD22	2:H:1:NAG:H83	1.58	0.85
1:F:359:LEU:HD13	1:F:400:THR:HG22	1.61	0.82
1:D:474:ILE:HD11	1:D:526:ALA:HB1	1.62	0.80
1:A:436:THR:HG22	1:C:453:TRP:HZ2	1.47	0.78
1:D:460:ARG:NH1	1:D:543:VAL:HG13	1.98	0.78
1:B:531:ALA:HB1	1:B:537:VAL:HG23	1.66	0.78
1:B:425:LEU:HD12	1:B:425:LEU:O	1.85	0.75
1:E:363:LEU:HD22	2:K:1:NAG:H83	1.68	0.74
1:E:499:LYS:NZ	1:F:510:GLU:HG2	2.03	0.73
1:D:359:LEU:HD11	1:D:398:THR:HG23	1.73	0.70
1:E:355:THR:HG22	1:E:404:PRO:HA	1.74	0.70
1:A:436:THR:HG22	1:C:453:TRP:CZ2	2.27	0.70
1:F:370:VAL:O	1:F:370:VAL:HG23	1.91	0.69
1:E:359:LEU:HD21	2:K:2:NAG:H2	1.74	0.69
1:D:425:LEU:HD12	1:D:425:LEU:O	1.94	0.68
1:F:476:VAL:C	1:F:477:GLN:HE21	2.03	0.67
1:D:371:GLN:O	1:D:372:LEU:HD23	1.97	0.65
1:D:425:LEU:HD12	1:D:425:LEU:C	2.22	0.65
1:B:403:LEU:HD12	1:B:404:PRO:HD2	1.80	0.64
1:B:477:GLN:NE2	1:B:484:GLN:OE1	2.32	0.63
1:C:396:THR:HG21	2:I:1:NAG:H3	1.80	0.62
1:E:499:LYS:HZ2	1:F:510:GLU:HG2	1.63	0.62
1:A:360:VAL:HG21	1:A:372:LEU:HD21	1.81	0.62
1:E:370:VAL:HG22	1:E:422:HIS:CD2	2.35	0.61
1:E:363:LEU:O	1:E:364:ALA:HB2	2.01	0.61
1:A:436:THR:CG2	1:C:453:TRP:HZ2	2.13	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:343:PRO:HD3	1:E:356:ILE:HG22	1.83	0.59
1:E:361:VAL:HG13	1:E:396:THR:HG21	1.84	0.59
1:F:425:LEU:HD12	1:F:425:LEU:C	2.27	0.59
1:F:372:LEU:HD22	1:F:420:VAL:HG22	1.85	0.58
1:A:411:ILE:HG22	1:D:534:SER:HB3	1.86	0.58
1:D:445:VAL:HG13	1:D:466:ILE:HD13	1.86	0.57
1:B:363:LEU:CD2	2:H:1:NAG:H83	2.32	0.57
1:E:356:ILE:HG23	1:E:405:VAL:HG11	1.86	0.57
1:E:445:VAL:HG22	1:E:466:ILE:HG13	1.87	0.56
1:E:376:ARG:HD2	1:E:378:SER:OG	2.05	0.56
1:B:348:LEU:HD23	1:B:349:PHE:CE1	2.41	0.56
1:C:425:LEU:HD12	1:C:426:PRO:HD2	1.88	0.56
1:C:481:ASN:HD21	1:C:518:GLN:CD	2.13	0.56
1:A:389:GLU:HG2	1:A:399:VAL:HG22	1.88	0.56
1:D:354:PRO:HG2	1:D:405:VAL:HG23	1.88	0.56
1:A:359:LEU:CD1	1:A:400:THR:HG22	2.34	0.55
1:B:372:LEU:HD22	1:B:420:VAL:HG22	1.88	0.55
1:D:359:LEU:HD13	1:D:400:THR:HG22	1.87	0.55
1:F:372:LEU:CD2	1:F:420:VAL:HG13	2.37	0.54
1:E:482:GLU:HA	1:E:482:GLU:OE1	2.07	0.54
1:B:360:VAL:HG21	1:B:372:LEU:HD21	1.89	0.54
1:E:383:GLN:O	1:E:403:LEU:HD11	2.07	0.54
1:E:416:TYR:O	1:E:433:THR:HG22	2.07	0.54
1:D:391:LYS:HG2	1:D:397:LEU:HD13	1.89	0.54
1:F:466:ILE:HG21	1:F:474:ILE:HD11	1.89	0.54
1:B:419:ARG:NE	1:B:430:MET:HE2	2.23	0.53
1:B:425:LEU:HD13	1:B:427:ARG:O	2.08	0.53
1:E:477:GLN:HE21	1:E:525:ARG:HD2	1.72	0.53
1:F:425:LEU:HD12	1:F:425:LEU:O	2.08	0.53
1:D:474:ILE:HD13	1:D:475:SER:N	2.23	0.53
1:D:478:TRP:C	1:D:479:LEU:HD12	2.33	0.53
1:C:371:GLN:O	1:C:372:LEU:HD23	2.09	0.53
1:D:460:ARG:CZ	1:D:543:VAL:HG13	2.39	0.53
1:D:340:LEU:HD12	1:D:357:THR:O	2.09	0.52
1:D:445:VAL:HG13	1:D:466:ILE:CD1	2.40	0.52
1:A:370:VAL:O	1:A:370:VAL:HG23	2.10	0.52
1:F:489:ARG:NH2	1:F:515:GLU:OE1	2.43	0.52
1:F:510:GLU:CD	1:F:510:GLU:N	2.69	0.51
1:C:481:ASN:O	1:C:482:GLU:C	2.53	0.51
1:B:425:LEU:HD12	1:B:425:LEU:C	2.35	0.51
1:C:440:ARG:HB3	1:C:530:ALA:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:356:ILE:H	1:F:356:ILE:HD13	1.75	0.50
1:A:411:ILE:CG2	1:D:534:SER:HB3	2.42	0.49
1:F:372:LEU:HD23	1:F:420:VAL:HG13	1.94	0.49
1:E:374:TRP:CE3	1:E:403:LEU:HD22	2.48	0.49
1:B:493:THR:O	1:B:505:VAL:HG23	2.12	0.48
1:E:338:ALA:HB1	1:E:431:ARG:CZ	2.43	0.48
1:C:513:ARG:NH1	1:C:517:GLU:OE2	2.46	0.48
1:F:350:ILE:HD11	1:F:527:VAL:HG21	1.95	0.48
1:C:479:LEU:HD11	1:C:525:ARG:CZ	2.43	0.48
1:D:479:LEU:HD21	1:D:525:ARG:NH2	2.28	0.48
1:E:340:LEU:HD12	1:E:357:THR:O	2.12	0.48
1:C:499:LYS:HZ1	1:D:489:ARG:HH11	1.61	0.47
1:C:531:ALA:HB1	1:C:537:VAL:HG23	1.96	0.47
1:E:363:LEU:HD22	2:K:1:NAG:C8	2.42	0.47
1:D:466:ILE:HB	1:D:505:VAL:HG13	1.96	0.47
1:C:453:TRP:CG	1:C:454:PRO:HD2	2.50	0.47
1:A:425:LEU:HD12	1:A:426:PRO:HD2	1.97	0.47
1:E:416:TYR:H	1:E:433:THR:HG22	1.80	0.46
1:F:355:THR:HG22	1:F:404:PRO:HA	1.97	0.46
1:D:422:HIS:N	1:D:425:LEU:HD23	2.31	0.46
1:F:523:ILE:HD12	1:F:523:ILE:N	2.30	0.46
1:C:466:ILE:HB	1:C:505:VAL:HG12	1.97	0.46
1:C:374:TRP:CD2	1:C:403:LEU:HD22	2.51	0.46
1:A:374:TRP:CD2	1:A:403:LEU:HD22	2.50	0.46
1:A:355:THR:HG22	1:A:404:PRO:HA	1.97	0.46
1:D:359:LEU:CD1	1:D:398:THR:HG23	2.44	0.46
1:E:396:THR:HG22	1:E:397:LEU:N	2.31	0.46
1:F:350:ILE:HD11	1:F:527:VAL:CG2	2.46	0.45
1:D:465:LEU:HD13	1:D:506:PHE:CZ	2.50	0.45
1:D:479:LEU:HD12	1:D:479:LEU:N	2.31	0.45
1:E:369:THR:HG21	1:E:371:GLN:NE2	2.31	0.45
1:F:470:MET:HA	1:F:471:PRO:C	2.41	0.45
1:C:479:LEU:HD23	1:C:484:GLN:HA	1.98	0.45
1:D:356:ILE:HD11	1:D:403:LEU:HD23	1.98	0.45
1:F:511:VAL:HG11	1:F:522:PHE:CE2	2.51	0.45
1:A:470:MET:HA	1:A:471:PRO:C	2.42	0.45
1:C:479:LEU:HD11	1:C:525:ARG:NE	2.32	0.45
1:E:358:CYS:SG	1:E:359:LEU:N	2.89	0.45
1:C:445:VAL:HG13	1:C:466:ILE:CD1	2.47	0.45
1:F:422:HIS:CG	1:F:423:PRO:HD2	2.52	0.45
1:C:479:LEU:HD23	1:C:484:GLN:CA	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:489:ARG:NH2	1:A:515:GLU:OE1	2.50	0.44
1:D:359:LEU:CD1	1:D:400:THR:HG22	2.48	0.44
1:F:527:VAL:HG22	1:F:536:THR:HG22	2.00	0.44
1:C:359:LEU:HD13	1:C:400:THR:HG22	1.98	0.44
1:E:363:LEU:CD2	2:K:1:NAG:H83	2.43	0.44
1:F:389:GLU:HG3	1:F:399:VAL:HG22	2.00	0.44
1:B:355:THR:HG22	1:B:404:PRO:HA	2.00	0.44
1:F:425:LEU:C	1:F:425:LEU:CD1	2.90	0.44
1:E:372:LEU:CD2	1:E:420:VAL:HG22	2.48	0.44
1:D:348:LEU:HD11	1:D:411:ILE:CD1	2.48	0.43
1:C:389:GLU:HG2	1:C:399:VAL:HG13	2.00	0.43
1:F:340:LEU:HD23	1:F:432:SER:C	2.44	0.43
1:F:477:GLN:HE21	1:F:477:GLN:N	2.16	0.43
1:C:479:LEU:HD23	1:C:484:GLN:N	2.34	0.43
1:F:370:VAL:O	1:F:370:VAL:CG2	2.63	0.43
1:F:444:GLU:CD	1:F:468:ASN:HD22	2.27	0.43
1:B:470:MET:HA	1:B:471:PRO:C	2.44	0.42
1:F:477:GLN:N	1:F:477:GLN:NE2	2.67	0.42
1:E:470:MET:HA	1:E:471:PRO:C	2.44	0.42
1:F:466:ILE:HG21	1:F:474:ILE:CD1	2.49	0.42
1:A:370:VAL:O	1:A:370:VAL:CG2	2.66	0.42
1:D:391:LYS:HB3	1:D:391:LYS:NZ	2.34	0.42
1:F:346:PHE:HD1	1:F:350:ILE:HD12	1.85	0.42
1:A:481:ASN:HD21	1:A:518:GLN:NE2	2.17	0.42
1:A:372:LEU:HD12	1:A:401:SER:HB2	2.02	0.42
1:B:420:VAL:O	1:B:428:ALA:HB1	2.19	0.42
1:D:425:LEU:C	1:D:425:LEU:CD1	2.90	0.42
1:A:376:ARG:HD3	1:A:414:GLU:HG2	2.00	0.42
1:C:498:THR:HG23	1:C:502:GLY:O	2.20	0.42
1:E:369:THR:HG21	1:E:371:GLN:HE21	1.83	0.42
1:B:386:THR:O	1:B:401:SER:HA	2.20	0.42
1:D:359:LEU:HD12	1:D:360:VAL:H	1.85	0.42
1:D:523:ILE:N	1:D:523:ILE:HD12	2.34	0.42
1:E:419:ARG:NH2	1:E:428:ALA:HB2	2.35	0.42
1:C:359:LEU:CD1	1:C:400:THR:HG22	2.50	0.41
1:B:440:ARG:HB3	1:B:530:ALA:HB2	2.02	0.41
1:E:422:HIS:ND1	1:E:423:PRO:HD2	2.34	0.41
1:E:465:LEU:HD12	1:E:466:ILE:N	2.36	0.41
1:C:372:LEU:CD2	1:C:420:VAL:HG13	2.50	0.41
1:E:373:THR:HB	1:E:419:ARG:HB3	2.02	0.41
1:D:470:MET:HA	1:D:471:PRO:C	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:475:SER:HB2	1:B:527:VAL:HB	2.03	0.41
1:C:439:PRO:HB2	1:C:470:MET:HE2	2.02	0.41
1:E:523:ILE:HD12	1:E:523:ILE:N	2.36	0.41
1:C:513:ARG:HA	1:C:516:TRP:CH2	2.56	0.41
1:E:360:VAL:HG21	1:E:420:VAL:HG21	2.03	0.41
1:D:498:THR:HG23	1:D:502:GLY:O	2.21	0.40
1:F:372:LEU:CD2	1:F:420:VAL:HG22	2.51	0.40
1:F:532:SER:HB2	1:F:533:PRO:HA	2.03	0.40
1:F:350:ILE:CD1	1:F:527:VAL:HG21	2.52	0.40
1:E:493:THR:OG1	1:E:505:VAL:HG23	2.22	0.40
1:C:391:LYS:HD3	1:C:391:LYS: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	207/223 (93%)	200 (97%)	7 (3%)	0	100	100
1	B	207/223 (93%)	198 (96%)	8 (4%)	1 (0%)	24	55
1	C	207/223 (93%)	199 (96%)	8 (4%)	0	100	100
1	D	207/223 (93%)	197 (95%)	9 (4%)	1 (0%)	24	55
1	E	207/223 (93%)	175 (84%)	29 (14%)	3 (1%)	9	30
1	F	207/223 (93%)	197 (95%)	10 (5%)	0	100	100
All	All	1242/1338 (93%)	1166 (94%)	71 (6%)	5 (0%)	30	60

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	364	ALA

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Mol	Chain	Res	Type
1	D	426	PRO
1	E	423	PRO
1	E	426	PRO
1	B	426	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	182/195 (93%)	180 (99%)	2 (1%)	65	87
1	B	185/195 (95%)	184 (100%)	1 (0%)	81	93
1	C	185/195 (95%)	182 (98%)	3 (2%)	55	83
1	D	185/195 (95%)	180 (97%)	5 (3%)	39	74
1	E	185/195 (95%)	181 (98%)	4 (2%)	45	78
1	F	183/195 (94%)	179 (98%)	4 (2%)	45	78
All	All	1105/1170 (94%)	1086 (98%)	19 (2%)	53	83

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	483	VAL
1	A	539	ARG
1	B	489	ARG
1	C	356	ILE
1	C	391	LYS
1	C	460	ARG
1	D	356	ILE
1	D	402	THR
1	D	434	THR
1	D	474	ILE
1	D	536	THR
1	E	422	HIS
1	E	433	THR
1	E	481	ASN

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Mol	Chain	Res	Type
1	E	539	ARG
1	F	356	ILE
1	F	386	THR
1	F	393	ARG
1	F	477	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	417	GLN
1	A	422	HIS
1	A	468	ASN
1	A	481	ASN
1	A	484	GLN
1	A	490	HIS
1	A	494	GLN
1	A	538	GLN
1	B	477	GLN
1	B	481	ASN
1	B	535	GLN
1	C	422	HIS
1	C	424	HIS
1	C	477	GLN
1	C	481	ASN
1	C	490	HIS
1	C	535	GLN
1	C	538	GLN
1	D	477	GLN
1	D	481	ASN
1	D	484	GLN
1	D	490	HIS
1	D	518	GLN
1	E	371	GLN
1	E	424	HIS
1	F	468	ASN
1	F	477	GLN
1	F	484	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 ⓘ

30 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	G	1	1,2	14,14,15	0.45	0	17,19,21	0.89	1 (5%)
2	NAG	G	2	2	14,14,15	0.54	0	17,19,21	0.88	1 (5%)
2	BMA	G	3	2	11,11,12	0.56	0	15,15,17	0.97	1 (6%)
2	MAN	G	4	2	11,11,12	0.58	0	15,15,17	0.87	1 (6%)
2	MAN	G	5	2	11,11,12	0.57	0	15,15,17	0.70	0
2	NAG	H	1	1,2	14,14,15	0.48	0	17,19,21	0.83	0
2	NAG	H	2	2	14,14,15	0.54	0	17,19,21	0.89	0
2	BMA	H	3	2	11,11,12	0.62	0	15,15,17	0.98	0
2	MAN	H	4	2	11,11,12	0.54	0	15,15,17	1.03	1 (6%)
2	MAN	H	5	2	11,11,12	0.56	0	15,15,17	0.73	0
2	NAG	I	1	1,2	14,14,15	0.55	0	17,19,21	0.71	0
2	NAG	I	2	2	14,14,15	0.51	0	17,19,21	0.81	0
2	BMA	I	3	2	11,11,12	0.63	0	15,15,17	1.20	1 (6%)
2	MAN	I	4	2	11,11,12	0.55	0	15,15,17	1.06	1 (6%)
2	MAN	I	5	2	11,11,12	0.65	0	15,15,17	1.10	0
2	NAG	J	1	1,2	14,14,15	0.48	0	17,19,21	0.99	1 (5%)
2	NAG	J	2	2	14,14,15	0.55	0	17,19,21	1.04	0
2	BMA	J	3	2	11,11,12	0.61	0	15,15,17	1.32	1 (6%)
2	MAN	J	4	2	11,11,12	0.55	0	15,15,17	0.95	1 (6%)
2	MAN	J	5	2	11,11,12	0.62	0	15,15,17	0.84	0
2	NAG	K	1	1,2	14,14,15	0.44	0	17,19,21	0.90	1 (5%)
2	NAG	K	2	2	14,14,15	0.51	0	17,19,21	0.86	1 (5%)
2	BMA	K	3	2	11,11,12	0.61	0	15,15,17	0.78	1 (6%)
2	MAN	K	4	2	11,11,12	0.56	0	15,15,17	1.03	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAN	K	5	2	11,11,12	0.60	0	15,15,17	0.58	0
2	NAG	L	1	1,2	14,14,15	0.48	0	17,19,21	0.74	0
2	NAG	L	2	2	14,14,15	0.49	0	17,19,21	0.70	0
2	BMA	L	3	2	11,11,12	0.64	0	15,15,17	0.92	0
2	MAN	L	4	2	11,11,12	0.57	0	15,15,17	0.81	1 (6%)
2	MAN	L	5	2	11,11,12	0.50	0	15,15,17	0.86	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
2	NAG	G	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
2	BMA	G	3	2	-	0/2/19/22	0/1/1/1
2	MAN	G	4	2	-	1/2/19/22	0/1/1/1
2	MAN	G	5	2	-	2/2/19/22	0/1/1/1
2	NAG	H	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1
2	BMA	H	3	2	-	1/2/19/22	0/1/1/1
2	MAN	H	4	2	-	2/2/19/22	0/1/1/1
2	MAN	H	5	2	-	2/2/19/22	0/1/1/1
2	NAG	I	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	BMA	I	3	2	-	0/2/19/22	0/1/1/1
2	MAN	I	4	2	-	1/2/19/22	0/1/1/1
2	MAN	I	5	2	-	1/2/19/22	0/1/1/1
2	NAG	J	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	BMA	J	3	2	-	2/2/19/22	0/1/1/1
2	MAN	J	4	2	-	2/2/19/22	0/1/1/1
2	MAN	J	5	2	-	2/2/19/22	0/1/1/1
2	NAG	K	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	K	2	2	-	0/6/23/26	0/1/1/1
2	BMA	K	3	2	-	2/2/19/22	0/1/1/1
2	MAN	K	4	2	-	0/2/19/22	0/1/1/1
2	MAN	K	5	2	-	0/2/19/22	0/1/1/1
2	NAG	L	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	L	2	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BMA	L	3	2	-	2/2/19/22	0/1/1/1
2	MAN	L	4	2	-	0/2/19/22	0/1/1/1
2	MAN	L	5	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	4	MAN	C1-O5-C5	3.20	116.48	112.19
2	H	4	MAN	C1-O5-C5	2.97	116.17	112.19
2	I	4	MAN	C1-O5-C5	2.77	115.90	112.19
2	J	4	MAN	C1-O5-C5	2.72	115.83	112.19
2	G	3	BMA	O3-C3-C4	2.48	116.21	110.38
2	K	1	NAG	C1-O5-C5	2.47	115.50	112.19
2	G	4	MAN	C1-O5-C5	2.45	115.47	112.19
2	J	3	BMA	O3-C3-C4	2.42	116.08	110.38
2	L	5	MAN	C1-O5-C5	2.38	115.38	112.19
2	L	4	MAN	C1-O5-C5	2.34	115.32	112.19
2	J	1	NAG	O5-C1-C2	-2.31	107.71	111.29
2	G	1	NAG	O5-C5-C6	2.16	111.87	107.66
2	K	3	BMA	C1-O5-C5	2.16	115.08	112.19
2	K	2	NAG	C1-O5-C5	2.16	115.08	112.19
2	I	3	BMA	O3-C3-C4	2.04	115.18	110.38
2	G	2	NAG	C1-O5-C5	2.02	114.90	112.19

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	5	MAN	O5-C5-C6-O6
2	J	4	MAN	O5-C5-C6-O6
2	J	5	MAN	O5-C5-C6-O6
2	L	3	BMA	O5-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
2	H	5	MAN	O5-C5-C6-O6
2	J	5	MAN	C4-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
2	J	4	MAN	C4-C5-C6-O6
2	H	4	MAN	C4-C5-C6-O6
2	H	5	MAN	C4-C5-C6-O6
2	H	4	MAN	O5-C5-C6-O6

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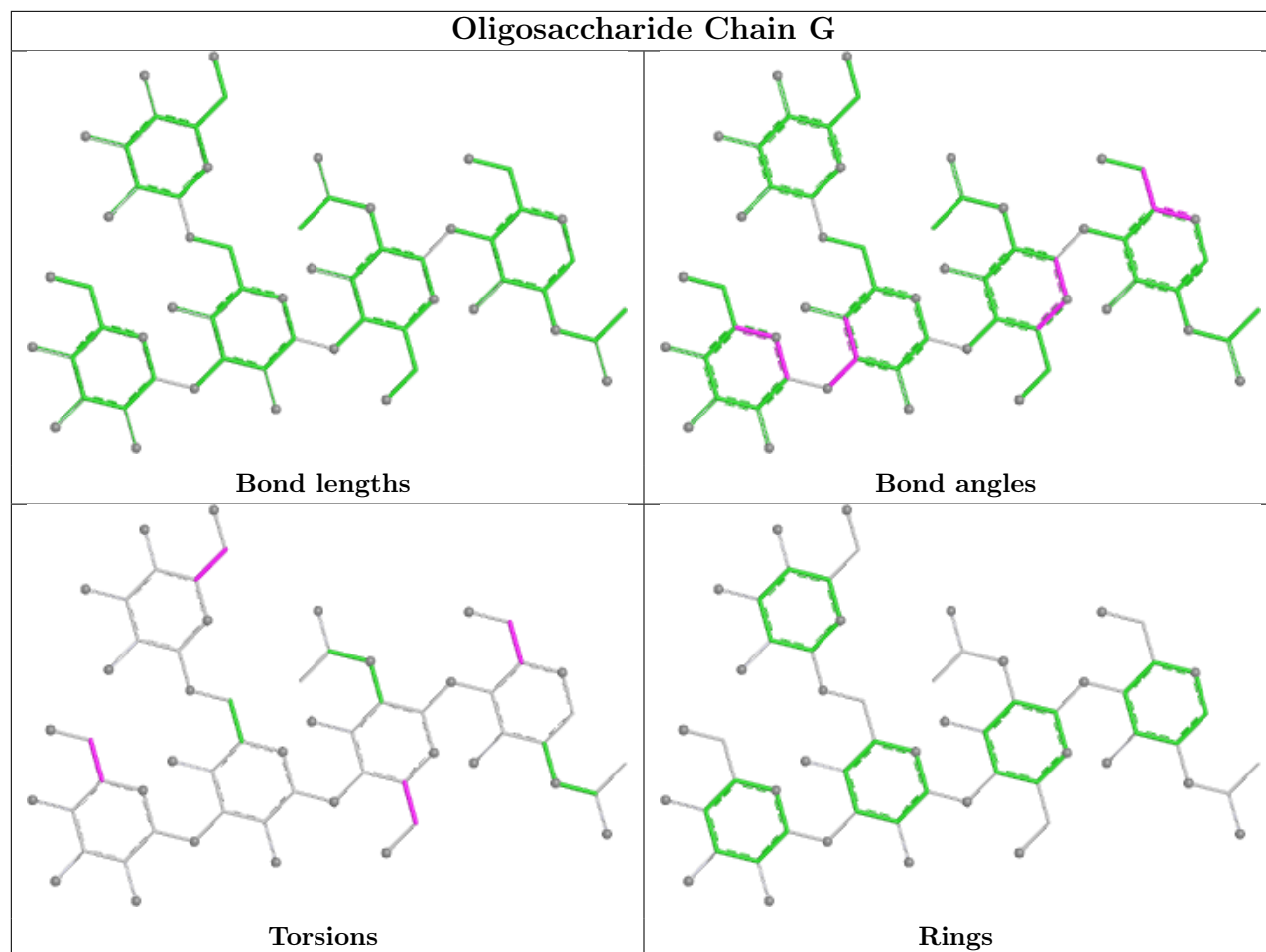
Mol	Chain	Res	Type	Atoms
2	G	5	MAN	C4-C5-C6-O6
2	L	3	BMA	C4-C5-C6-O6
2	J	3	BMA	O5-C5-C6-O6
2	J	1	NAG	C8-C7-N2-C2
2	J	1	NAG	O7-C7-N2-C2
2	J	2	NAG	C4-C5-C6-O6
2	L	2	NAG	C4-C5-C6-O6
2	I	1	NAG	O5-C5-C6-O6
2	J	3	BMA	C4-C5-C6-O6
2	K	3	BMA	C4-C5-C6-O6
2	L	2	NAG	O5-C5-C6-O6
2	I	4	MAN	O5-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
2	I	5	MAN	C4-C5-C6-O6
2	K	3	BMA	O5-C5-C6-O6
2	H	3	BMA	O5-C5-C6-O6
2	G	4	MAN	C4-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6

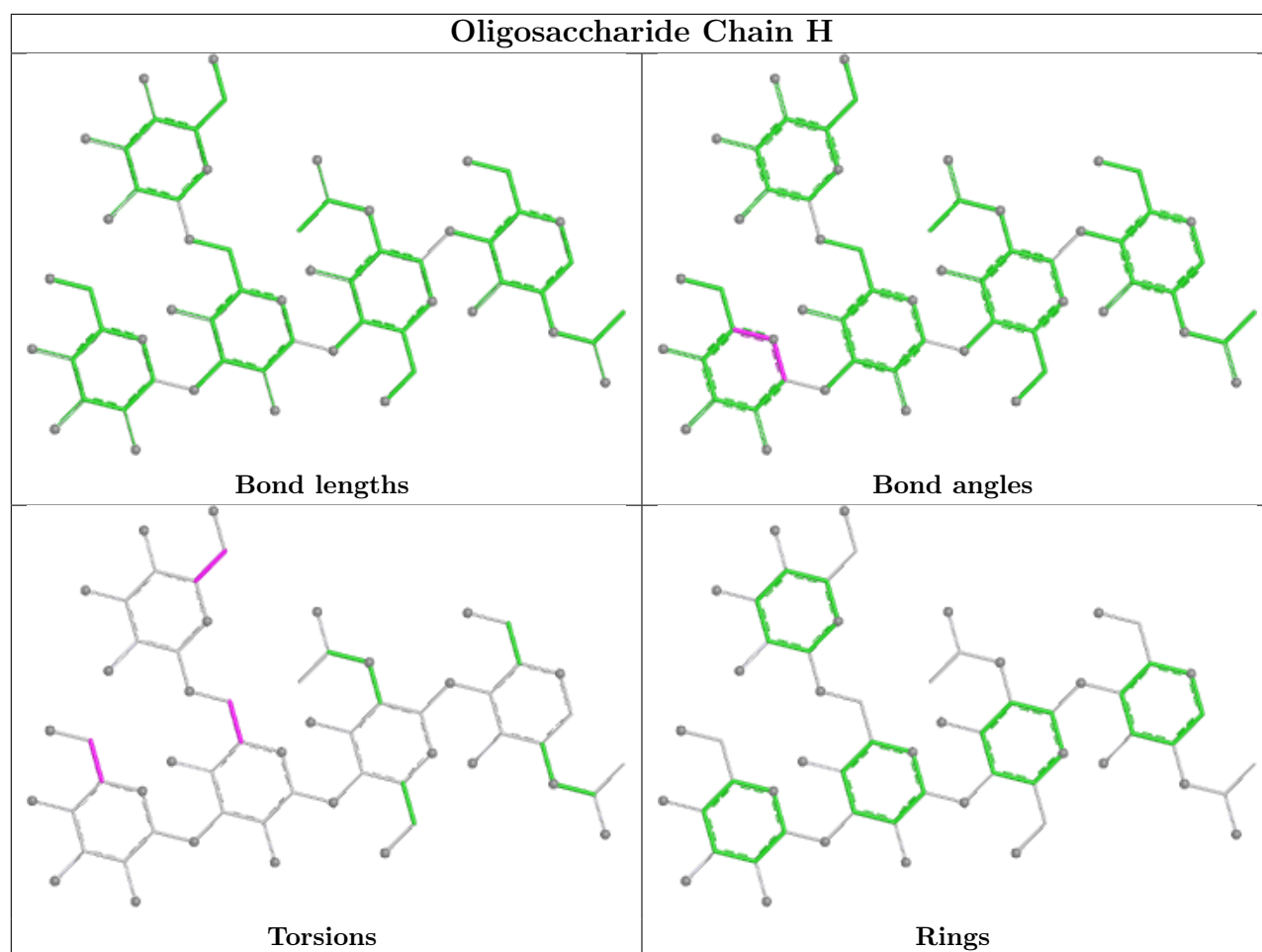
There are no ring outliers.

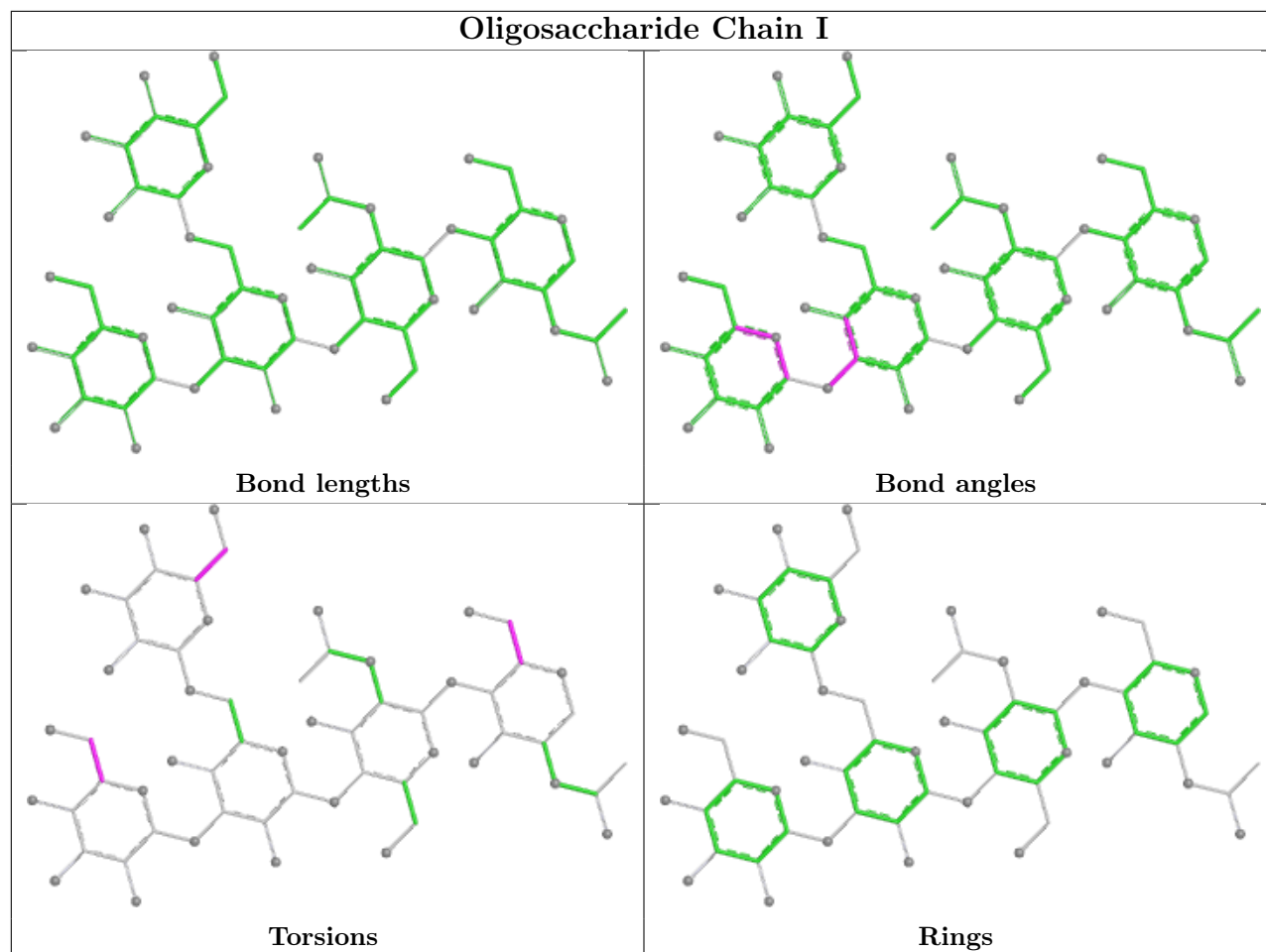
4 monomers are involved in 7 short contacts:

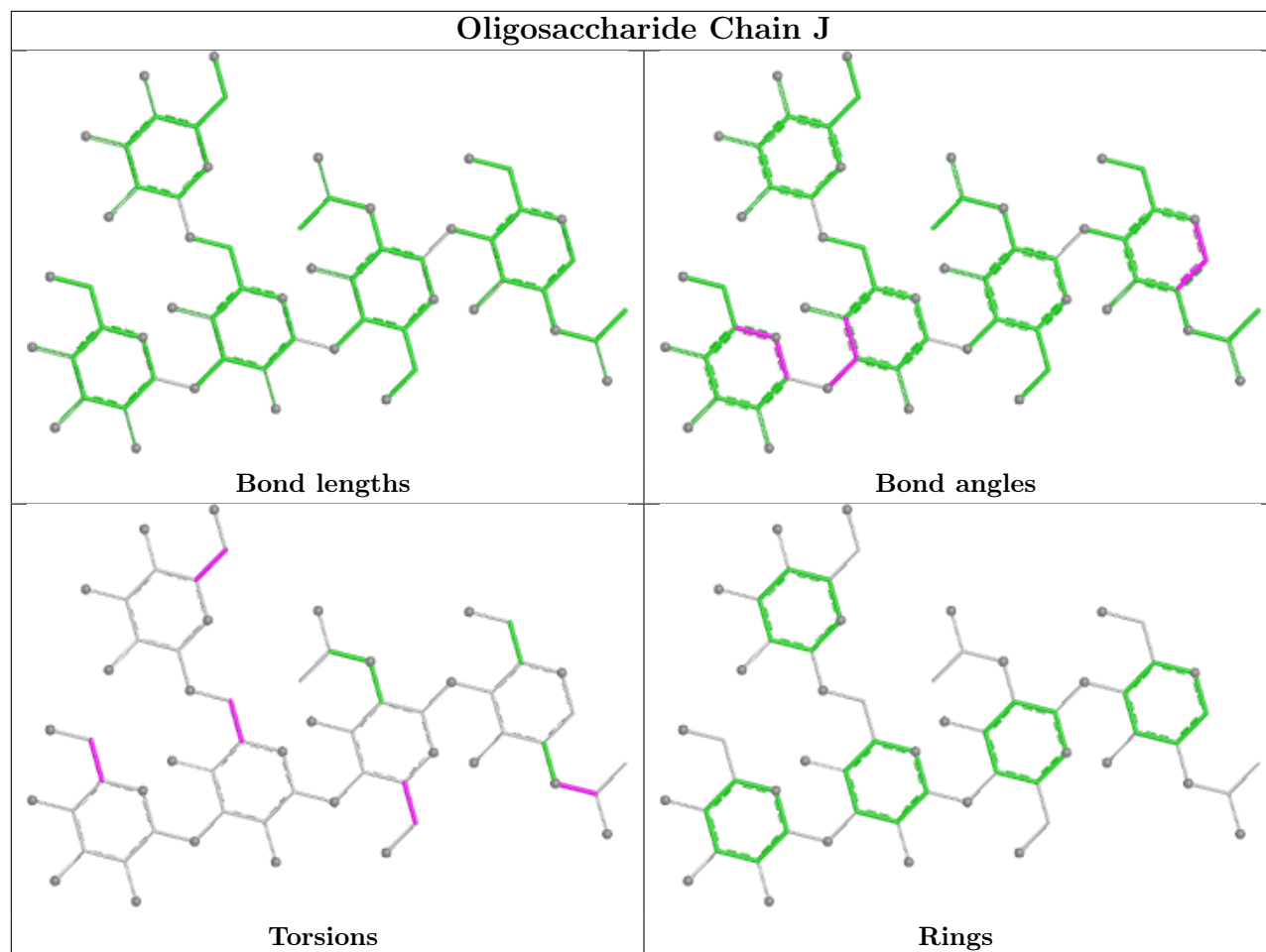
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	1	NAG	2	0
2	K	1	NAG	3	0
2	K	2	NAG	1	0
2	I	1	NAG	1	0

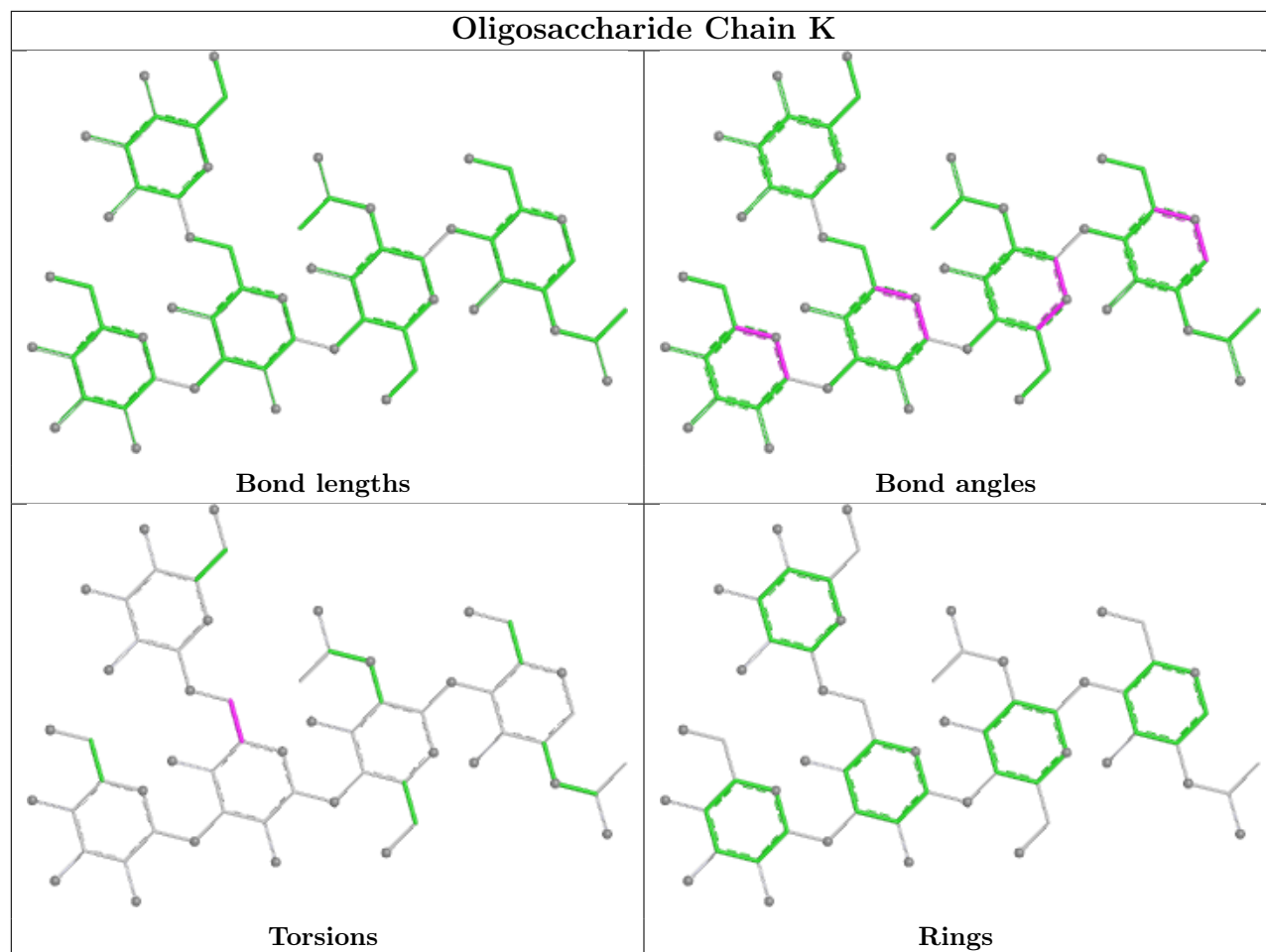
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

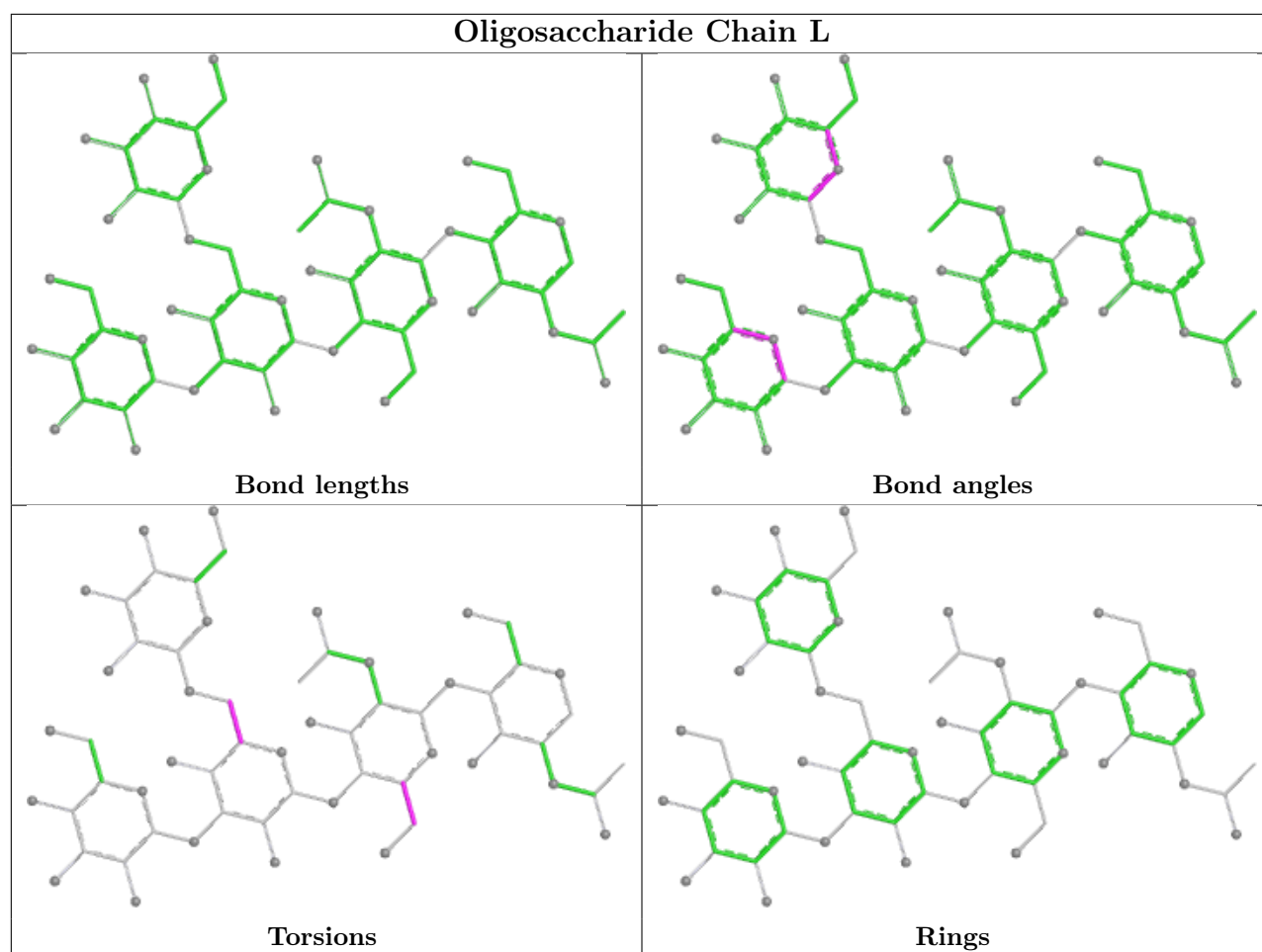












5.6 Ligand geometry [i](#)

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 [i](#)

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	209/223 (93%)	0.34	6 (2%) 53 43	37, 54, 76, 82	0
1	B	209/223 (93%)	0.48	10 (4%) 35 28	45, 58, 86, 92	0
1	C	209/223 (93%)	0.98	22 (10%) 11 8	55, 74, 98, 104	0
1	D	209/223 (93%)	0.97	20 (9%) 13 10	44, 70, 94, 101	0
1	E	209/223 (93%)	1.53	61 (29%) 1 1	65, 93, 124, 126	0
1	F	209/223 (93%)	1.06	30 (14%) 6 5	54, 72, 101, 103	0
All	All	1254/1338 (93%)	0.89	149 (11%) 9 7	37, 69, 115, 126	0

All (149) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	420	VAL	6.7
1	C	395	GLY	5.4
1	C	393	ARG	5.1
1	F	426	PRO	4.3
1	F	361	VAL	4.2
1	D	372	LEU	4.0
1	F	427	ARG	4.0
1	E	416	TYR	4.0
1	E	349	PHE	4.0
1	E	340	LEU	3.9
1	E	432	SER	3.9
1	C	337	SER	3.9
1	C	338	ALA	3.8
1	E	339	TYR	3.8
1	E	431	ARG	3.8
1	E	413	GLY	3.8
1	E	370	VAL	3.6
1	D	418	CYS	3.6
1	B	363	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	E	380	LYS	3.5
1	C	453	TRP	3.5
1	E	363	LEU	3.4
1	A	477	GLN	3.4
1	E	429	LEU	3.3
1	E	364	ALA	3.3
1	E	397	LEU	3.3
1	C	392	GLN	3.3
1	F	420	VAL	3.2
1	E	411	ILE	3.2
1	E	441	ALA	3.1
1	E	400	THR	3.1
1	F	397	LEU	3.1
1	E	425	LEU	3.1
1	F	416	TYR	3.1
1	D	361	VAL	3.1
1	E	407	THR	3.1
1	E	354	PRO	3.0
1	E	532	SER	3.0
1	E	338	ALA	3.0
1	E	377	ALA	3.0
1	E	379	GLY	3.0
1	B	399	VAL	3.0
1	D	420	VAL	3.0
1	E	365	PRO	3.0
1	B	397	LEU	2.9
1	E	381	PRO	2.9
1	E	348	LEU	2.9
1	D	338	ALA	2.9
1	F	337	SER	2.9
1	A	362	ASP	2.8
1	F	362	ASP	2.8
1	E	360	VAL	2.8
1	E	361	VAL	2.8
1	E	356	ILE	2.8
1	E	421	THR	2.8
1	F	360	VAL	2.8
1	B	364	ALA	2.8
1	F	428	ALA	2.8
1	B	422	HIS	2.8
1	C	440	ARG	2.7
1	E	534	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	405	VAL	2.7
1	C	363	LEU	2.7
1	C	422	HIS	2.7
1	E	422	HIS	2.7
1	E	433	THR	2.7
1	D	411	ILE	2.6
1	E	359	LEU	2.6
1	F	339	TYR	2.6
1	C	428	ALA	2.6
1	E	437	SER	2.6
1	F	395	GLY	2.6
1	D	521	GLU	2.6
1	B	437	SER	2.5
1	F	359	LEU	2.5
1	F	350	ILE	2.5
1	A	360	VAL	2.5
1	A	482	GLU	2.5
1	C	521	GLU	2.5
1	D	419	ARG	2.5
1	F	353	SER	2.5
1	D	474	ILE	2.5
1	C	429	LEU	2.5
1	B	339	TYR	2.5
1	C	436	THR	2.5
1	F	336	VAL	2.4
1	E	410	TRP	2.4
1	C	481	ASN	2.4
1	A	423	PRO	2.4
1	D	336	VAL	2.4
1	D	353	SER	2.4
1	F	392	GLN	2.4
1	E	383	GLN	2.4
1	F	477	GLN	2.4
1	D	415	THR	2.4
1	E	398	THR	2.4
1	F	372	LEU	2.4
1	E	399	VAL	2.3
1	E	406	GLY	2.3
1	E	336	VAL	2.3
1	E	394	ASN	2.3
1	D	389	GLU	2.3
1	E	358	CYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	442	ALA	2.3
1	C	494	GLN	2.3
1	D	382	VAL	2.3
1	E	405	VAL	2.3
1	C	373	THR	2.3
1	E	357	THR	2.3
1	B	533	PRO	2.3
1	D	405	VAL	2.3
1	E	346	PHE	2.2
1	C	388	LYS	2.2
1	B	458	ASP	2.2
1	E	408	ARG	2.2
1	F	415	THR	2.2
1	C	418	CYS	2.2
1	C	454	PRO	2.2
1	E	386	THR	2.2
1	F	369	THR	2.2
1	D	337	SER	2.2
1	E	530	ALA	2.2
1	E	537	VAL	2.2
1	B	396	THR	2.2
1	E	536	THR	2.2
1	C	456	SER	2.2
1	D	339	TYR	2.2
1	E	474	ILE	2.1
1	F	370	VAL	2.1
1	F	439	PRO	2.1
1	F	406	GLY	2.1
1	D	438	GLY	2.1
1	E	351	ARG	2.1
1	D	344	SER	2.1
1	E	375	SER	2.1
1	C	361	VAL	2.1
1	E	435	LYS	2.1
1	F	382	VAL	2.1
1	F	429	LEU	2.1
1	C	437	SER	2.0
1	E	345	PRO	2.0
1	E	342	ARG	2.0
1	E	368	GLY	2.0
1	F	411	ILE	2.0
1	E	391	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	340	LEU	2.0
1	F	368	GLY	2.0
1	A	364	ALA	2.0
1	D	367	LYS	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.

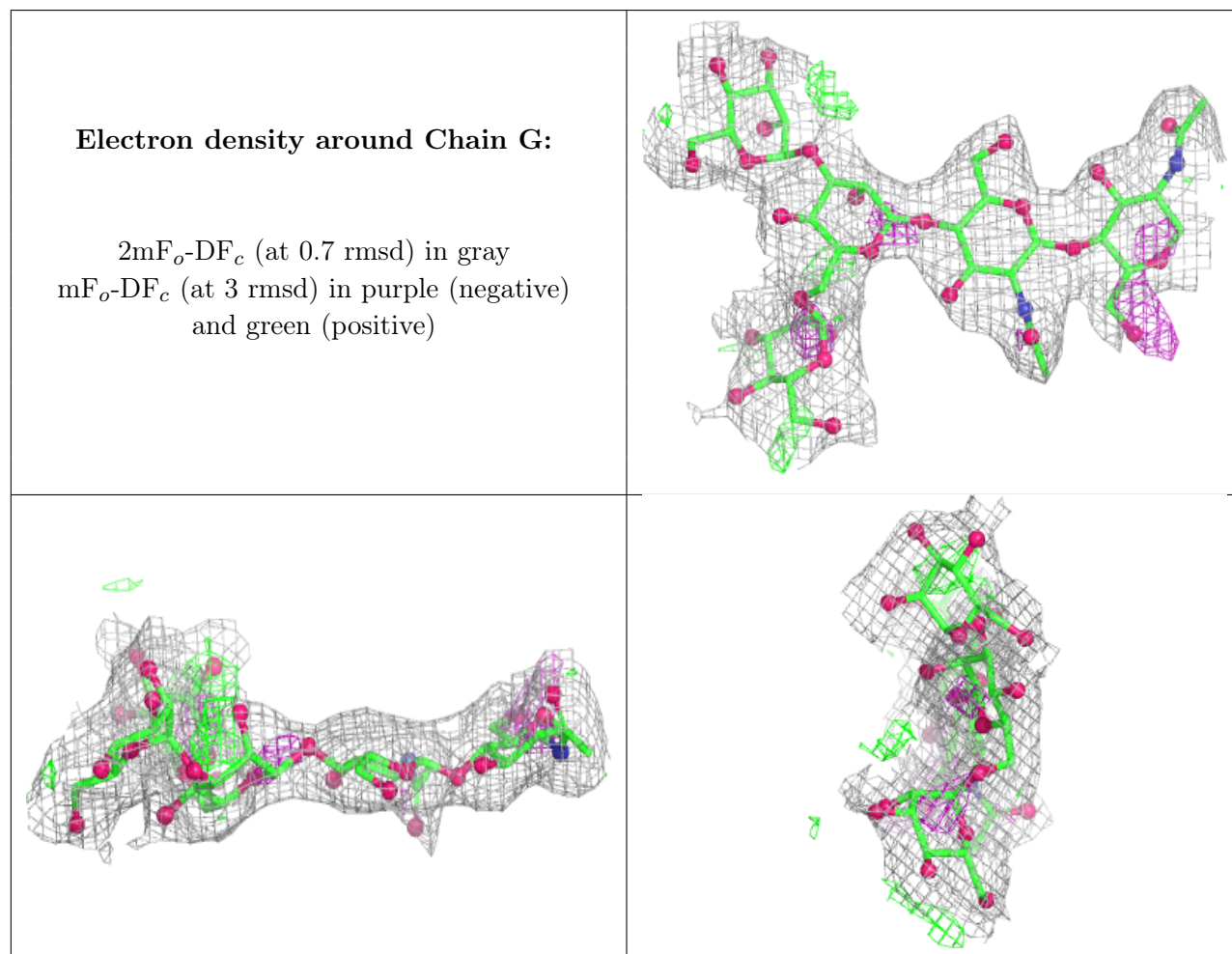
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	L	1	14/15	-0.04	0.26	100,101,101,102	0
2	BMA	K	3	11/12	-0.01	0.19	109,109,110,110	0
2	MAN	L	5	11/12	0.10	0.20	107,107,107,107	0
2	MAN	K	4	11/12	0.13	0.19	109,109,109,109	0
2	MAN	I	4	11/12	0.14	0.21	100,100,101,101	0
2	BMA	L	3	11/12	0.15	0.19	104,105,106,106	0
2	NAG	K	2	14/15	0.17	0.21	110,110,111,111	0
2	MAN	L	4	11/12	0.19	0.17	104,105,105,105	0
2	MAN	H	4	11/12	0.23	0.18	99,99,99,99	0
2	NAG	K	1	14/15	0.24	0.18	111,112,112,112	0
2	MAN	G	5	11/12	0.30	0.18	94,94,94,94	0
2	MAN	K	5	11/12	0.30	0.19	109,109,109,109	0
2	MAN	H	5	11/12	0.30	0.19	100,100,101,101	0
2	NAG	J	1	14/15	0.32	0.18	97,97,98,99	0
2	NAG	L	2	14/15	0.33	0.22	102,103,103,104	0
2	MAN	G	4	11/12	0.38	0.15	90,91,91,91	0
2	MAN	J	5	11/12	0.39	0.18	105,105,106,106	0
2	MAN	J	4	11/12	0.42	0.14	102,103,103,103	0
2	NAG	H	1	14/15	0.47	0.17	93,94,94,95	0
2	NAG	J	2	14/15	0.49	0.20	100,101,101,102	0
2	BMA	J	3	11/12	0.52	0.15	102,102,103,104	0
2	MAN	I	5	11/12	0.54	0.15	95,96,96,96	0
2	BMA	H	3	11/12	0.54	0.16	98,98,99,100	0
2	BMA	I	3	11/12	0.59	0.16	96,97,98,99	0

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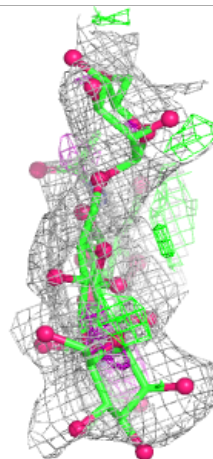
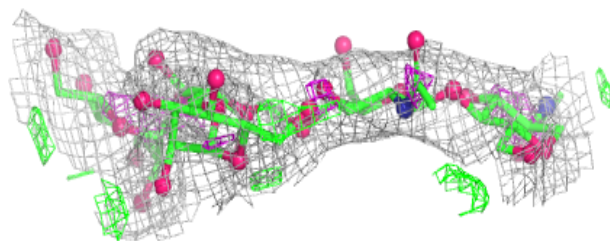
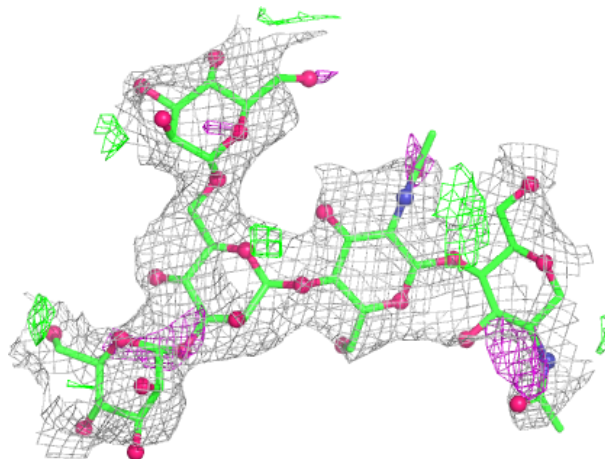
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	G	1	14/15	0.65	0.16	76,77,79,80	0
2	BMA	G	3	11/12	0.65	0.15	88,89,91,93	0
2	NAG	H	2	14/15	0.66	0.17	96,96,96,97	0
2	NAG	I	1	14/15	0.66	0.14	94,95,95,95	0
2	NAG	I	2	14/15	0.68	0.17	95,96,96,97	0
2	NAG	G	2	14/15	0.81	0.12	81,82,84,86	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.



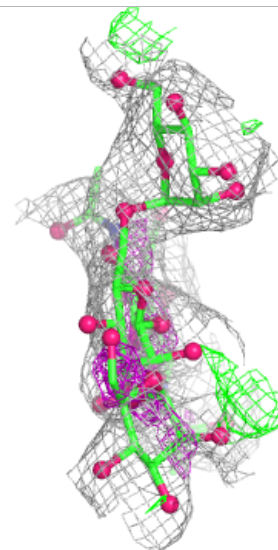
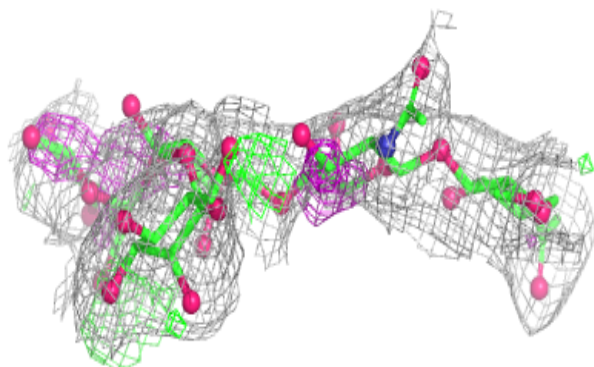
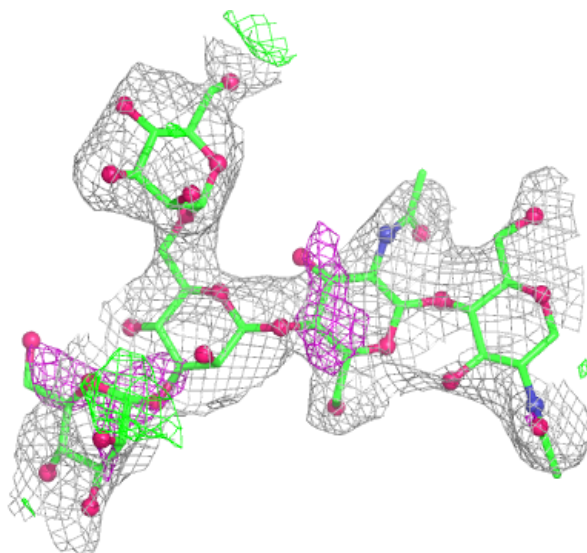
Electron density around Chain H:

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



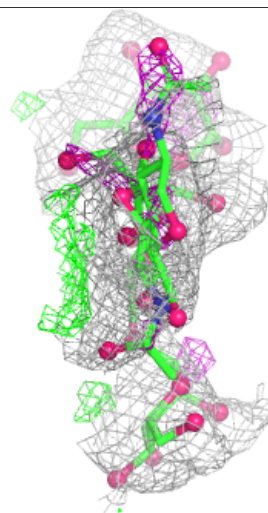
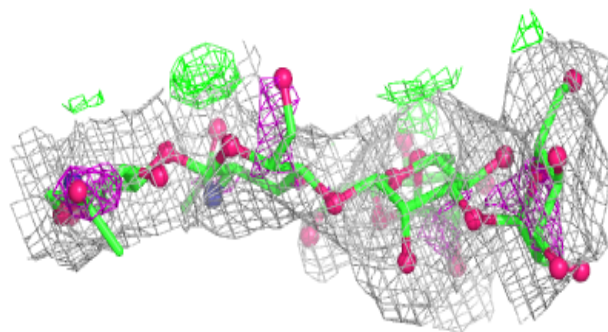
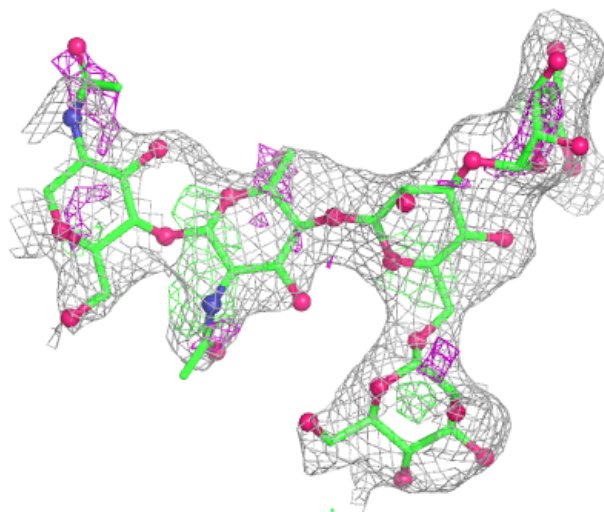
Electron density around Chain I:

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



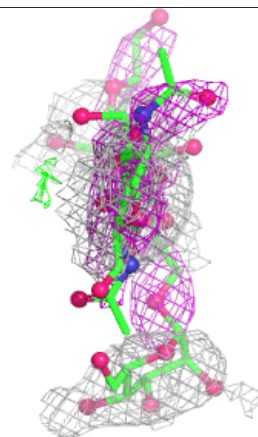
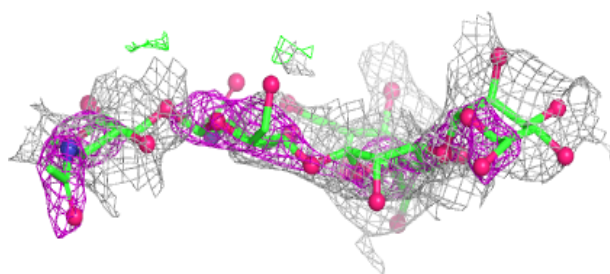
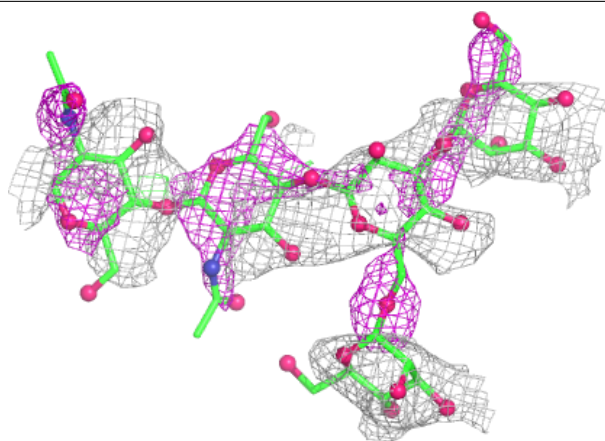
Electron density around Chain J:

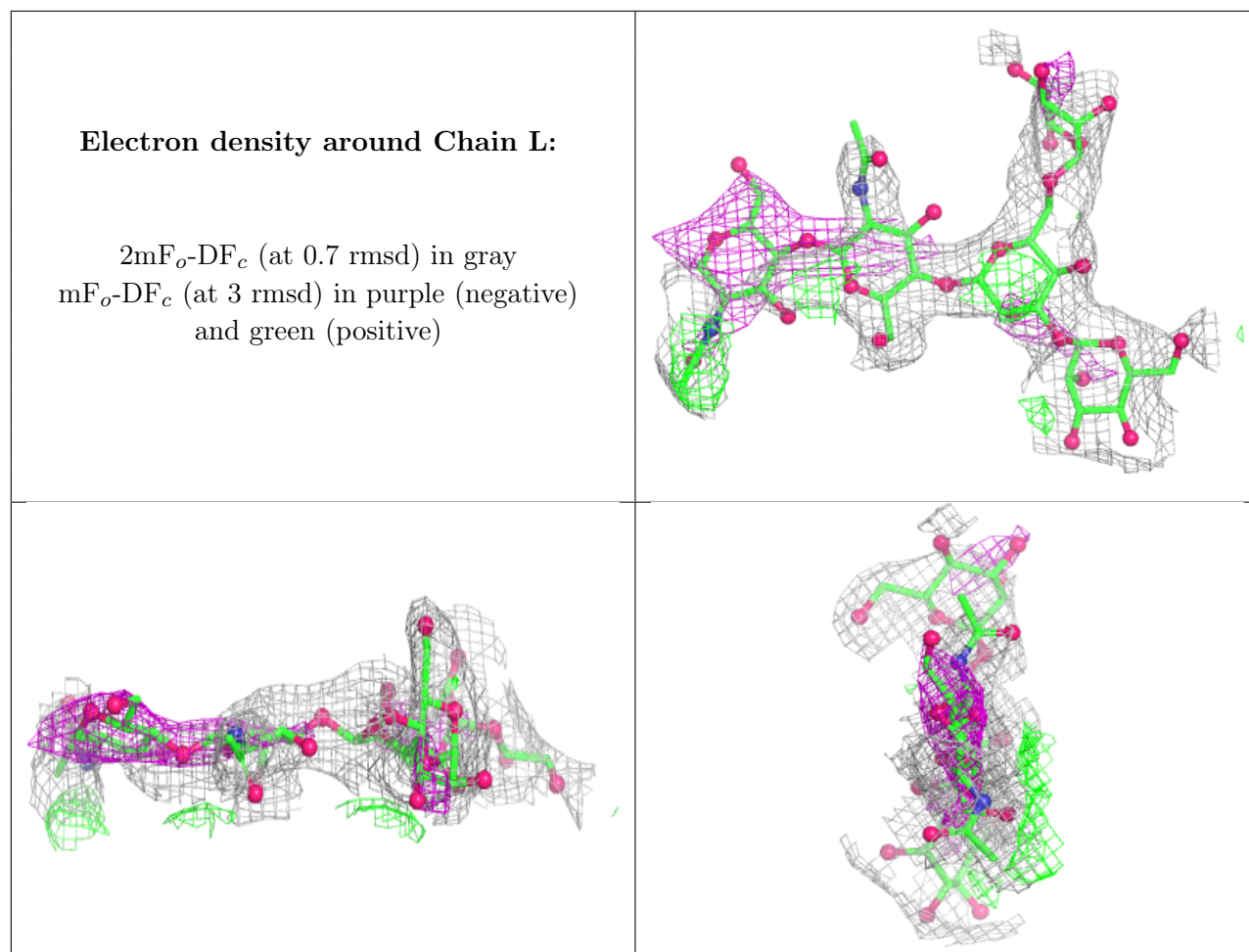
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain K:

$2mF_o-DF_c$ (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.