



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 12:06 AM UTC

PDB ID : 5SZL / pdb_00005szl
Title : Protocadherin gamma A1 extracellular cadherin domains 1-4
Authors : Goodman, K.M.; Bahna, F.; Mannepalli, S.; Honig, B.; Shapiro, L.
Deposited on : 2016-08-14
Resolution : 4.20 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

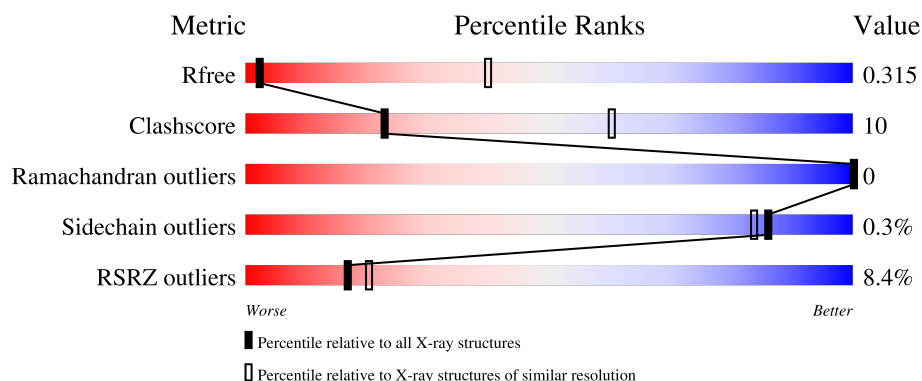
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.20 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	1004 (4.52-3.88)
Clashscore	190562	1019 (4.50-3.90)
Ramachandran outliers	187476	1020 (4.56-3.84)
Sidechain outliers	187428	1006 (4.56-3.84)
RSRZ outliers	180081	1001 (4.52-3.88)

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	427	5% 74% 22% .
1	B	427	7% 74% 22% .
1	C	427	9% 78% 19% .
1	D	427	11% 78% 19% .
2	E	3	100%

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Mol	Chain	Length	Quality of chain
3	F	2	 100%
4	G	4	 50%50%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 12505 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 PROTOCADHERIN GAMMA A1 EXTRACELLULAR CADHERIN DOMAINS 1-4, Protein Pcdhga1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	413	Total	C	N	O	S	0	0	0
			3004	1859	518	618	9			
1	B	412	Total	C	N	O	S	0	0	0
			2974	1840	512	614	8			
1	C	418	Total	C	N	O	S	0	0	0
			3119	1936	535	638	10			
1	D	416	Total	C	N	O	S	0	0	0
			3098	1928	521	640	9			

There are 32 discrepancies between the modelled and reference sequences:

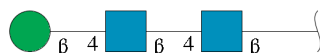
Chain	Residue	Modelled	Actual	Comment	Reference
A	420	HIS	-	expression tag	UNP A0A0A6YW27
A	421	HIS	-	expression tag	UNP A0A0A6YW27
A	422	HIS	-	expression tag	UNP A0A0A6YW27
A	423	HIS	-	expression tag	UNP A0A0A6YW27
A	424	HIS	-	expression tag	UNP A0A0A6YW27
A	425	HIS	-	expression tag	UNP A0A0A6YW27
A	426	HIS	-	expression tag	UNP A0A0A6YW27
A	427	HIS	-	expression tag	UNP A0A0A6YW27
B	420	HIS	-	expression tag	UNP A0A0A6YW27
B	421	HIS	-	expression tag	UNP A0A0A6YW27
B	422	HIS	-	expression tag	UNP A0A0A6YW27
B	423	HIS	-	expression tag	UNP A0A0A6YW27
B	424	HIS	-	expression tag	UNP A0A0A6YW27
B	425	HIS	-	expression tag	UNP A0A0A6YW27
B	426	HIS	-	expression tag	UNP A0A0A6YW27
B	427	HIS	-	expression tag	UNP A0A0A6YW27
C	420	HIS	-	expression tag	UNP A0A0A6YW27
C	421	HIS	-	expression tag	UNP A0A0A6YW27
C	422	HIS	-	expression tag	UNP A0A0A6YW27
C	423	HIS	-	expression tag	UNP A0A0A6YW27

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Chain	Residue	Modelled	Actual	Comment	Reference
C	424	HIS	-	expression tag	UNP A0A0A6YW27
C	425	HIS	-	expression tag	UNP A0A0A6YW27
C	426	HIS	-	expression tag	UNP A0A0A6YW27
C	427	HIS	-	expression tag	UNP A0A0A6YW27
D	420	HIS	-	expression tag	UNP A0A0A6YW27
D	421	HIS	-	expression tag	UNP A0A0A6YW27
D	422	HIS	-	expression tag	UNP A0A0A6YW27
D	423	HIS	-	expression tag	UNP A0A0A6YW27
D	424	HIS	-	expression tag	UNP A0A0A6YW27
D	425	HIS	-	expression tag	UNP A0A0A6YW27
D	426	HIS	-	expression tag	UNP A0A0A6YW27
D	427	HIS	-	expression tag	UNP A0A0A6YW27

- Molecule 2 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



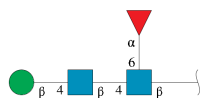
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



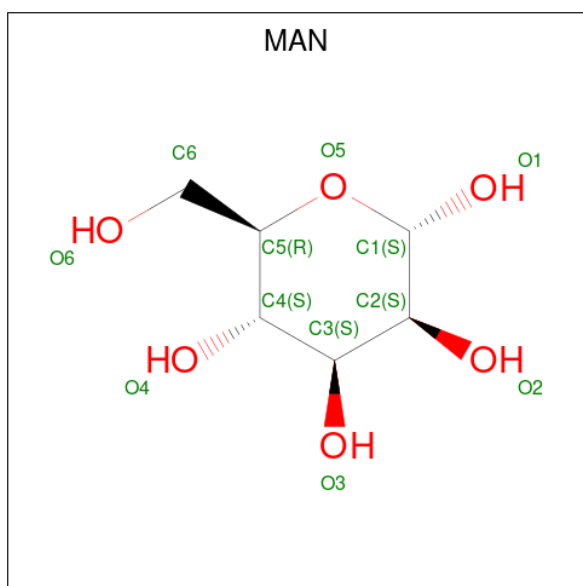
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	4	Total	C	N	O	0	0	0
			49	28	2	19			

- Molecule 5 is alpha-D-mannopyranose (CCD ID: MAN) (formula: $C_6H_{12}O_6$).

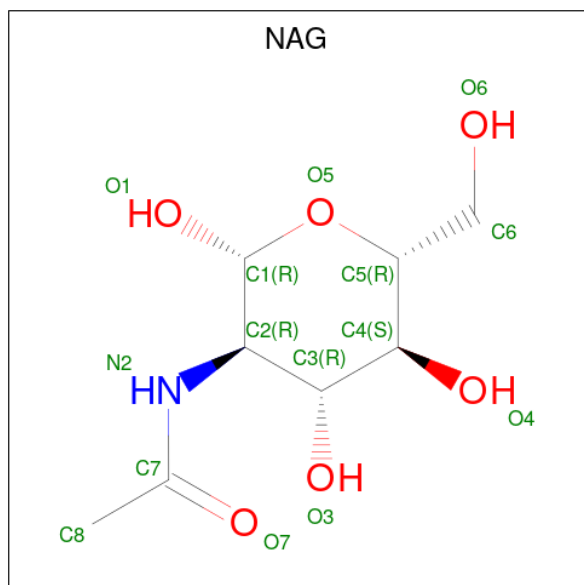


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			11	6	5		
5	A	1	Total	C	O	0	0
			11	6	5		
5	B	1	Total	C	O	0	0
			11	6	5		
5	B	1	Total	C	O	0	0
			11	6	5		
5	C	1	Total	C	O	0	0
			11	6	5		
5	C	1	Total	C	O	0	0
			11	6	5		
5	D	1	Total	C	O	0	0
			11	6	5		
5	D	1	Total	C	O	0	0
			11	6	5		

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	9	Total	Ca	0	0
			9	9		
6	B	9	Total	Ca	0	0
			9	9		
6	C	9	Total	Ca	0	0
			9	9		
6	D	9	Total	Ca	0	0
			9	9		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

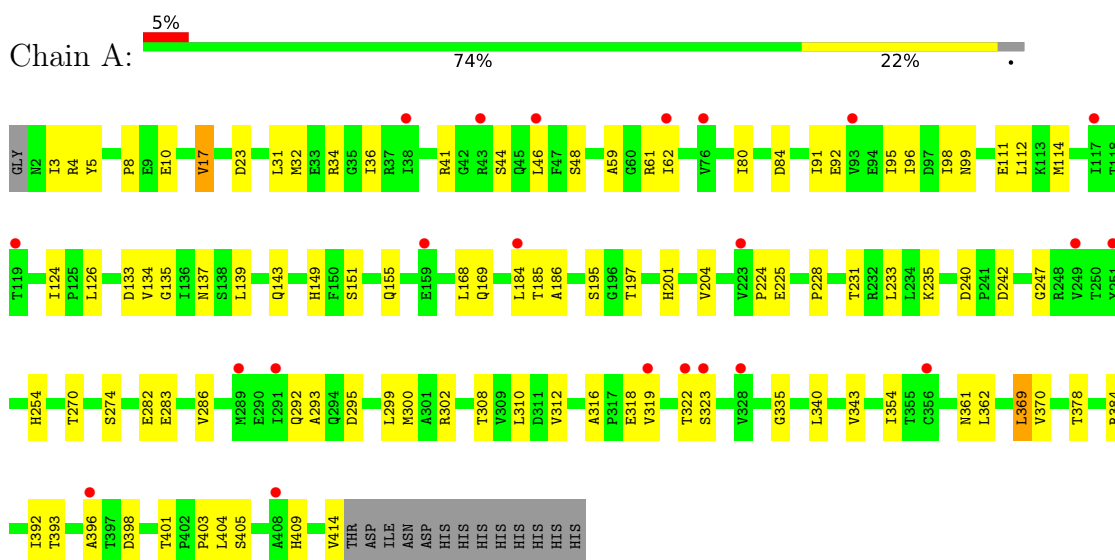


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		
7	D	1	Total	C	N	O	0	0
			14	8	1	5		

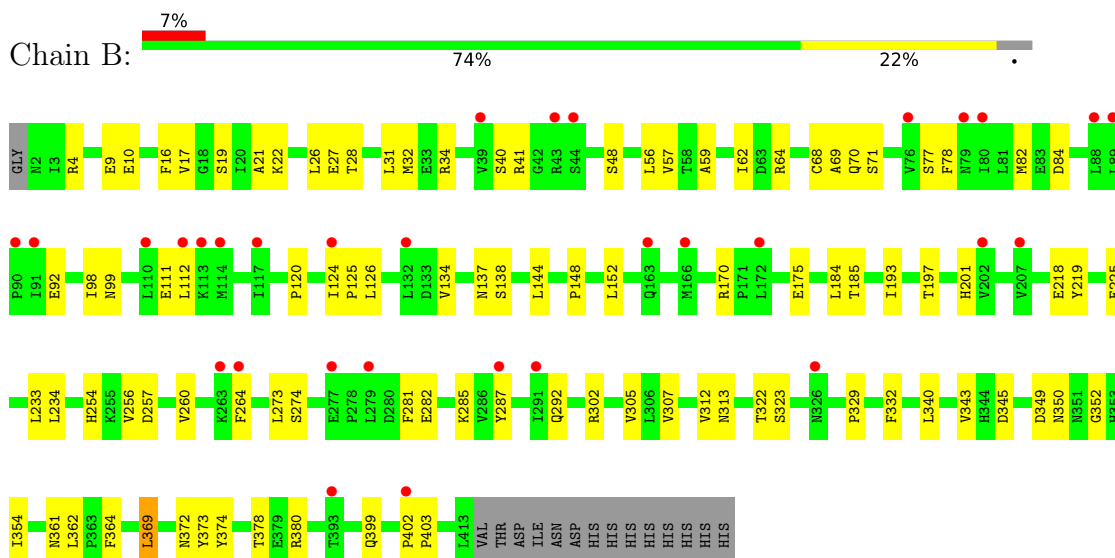
3 Residue-property plots

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.

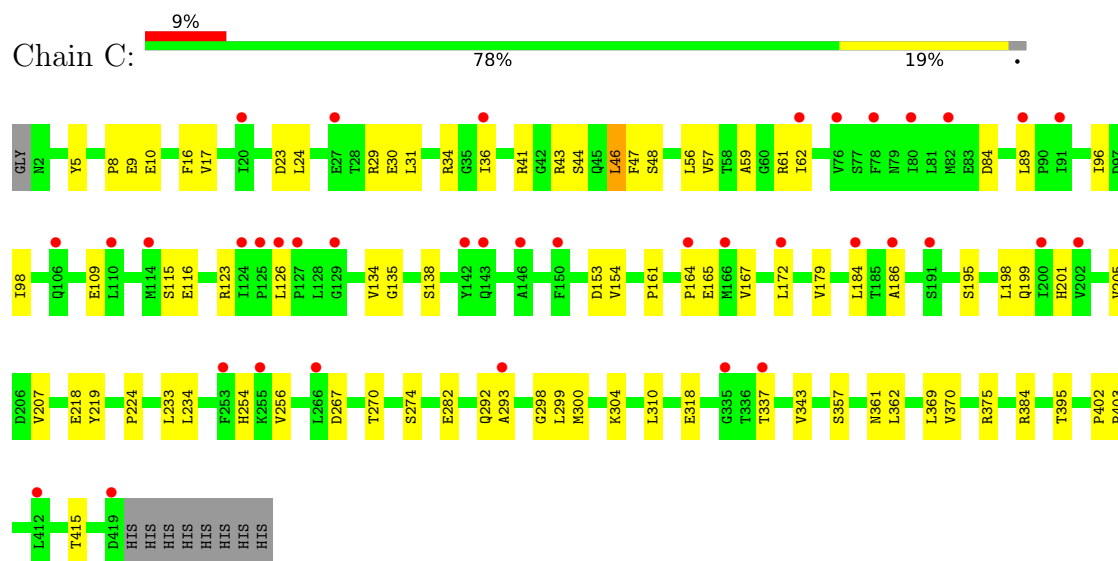
- Molecule 1: PROTOCADHERIN GAMMA A1 EXTRACELLULAR CADHERIN DOMAINS 1-4, Protein Pcdhgal



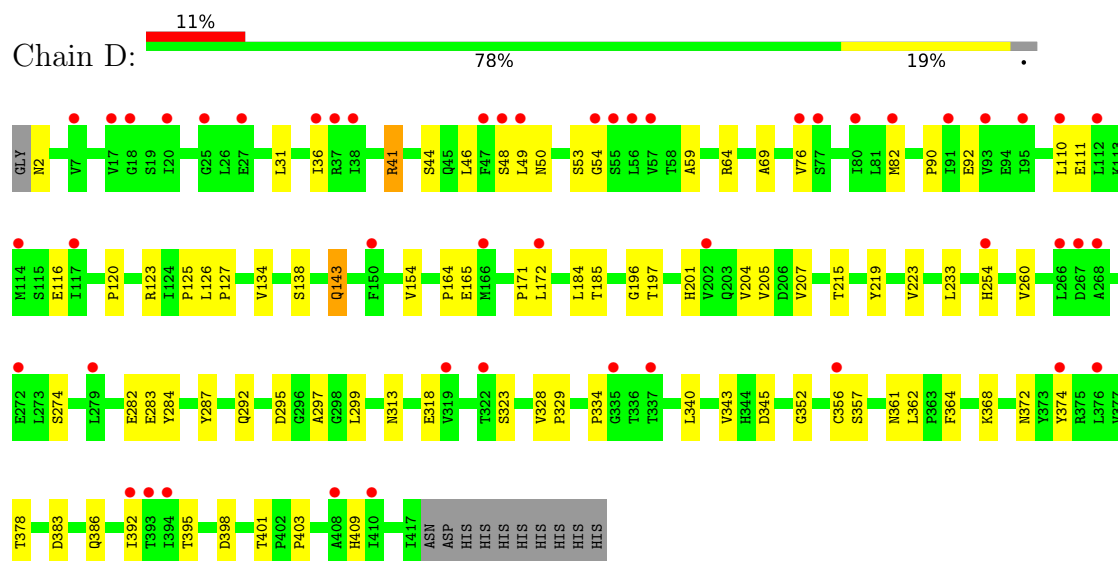
- Molecule 1: PROTOCADHERIN GAMMA A1 EXTRACELLULAR CADHERIN DOMAINS 1-4, Protein Pcdhgal



- Molecule 1: PROTOCADHERIN GAMMA A1 EXTRACELLULAR CADHERIN DOMAINS 1-4, Protein Pcdhga1



- Molecule 1: PROTOCADHERIN GAMMA A1 EXTRACELLULAR CADHERIN DOMAINS 1-4, Protein Pcdhga1



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



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



MAG1
MAG2

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

Chain G:



MAG1
MAG2
BMA3
FUC4

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	107.87Å 107.87Å 463.08Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.96 – 4.20 39.96 – 4.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (39.96-4.20) 79.8 (39.96-4.20)	Depositor EDS
R_{merge}	0.38	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.48 (at 4.13Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.287 , 0.314 0.286 , 0.315	Depositor DCC
R_{free} test set	1192 reflections (4.33%)	wwPDB-VP
Wilson B-factor (Å ²)	109.7	Xtriage
Anisotropy	0.491	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.21 , 142.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.042 for -h,-k,l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	12505	wwPDB-VP
Average B, all atoms (Å ²)	196.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, MAN, BMA, CA, FUC

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	0.21	0/3061	0.48	1/4195 (0.0%)
1	B	0.20	0/3028	0.47	2/4147 (0.0%)
1	C	0.17	0/3177	0.45	0/4348
1	D	0.23	0/3156	0.54	3/4321 (0.1%)
All	All	0.20	0/12422	0.49	6/17011 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	223	VAL	CB-CA-C	7.33	116.51	109.33
1	B	40	SER	CA-C-N	6.01	131.73	121.87
1	B	40	SER	C-N-CA	6.01	131.73	121.87
1	A	17	VAL	N-CA-C	-5.30	106.50	111.48
1	D	356	CYS	CA-CB-SG	-5.24	102.35	114.40
1	D	41	ARG	N-CA-C	-5.06	101.72	109.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3004	0	2731	68	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2974	0	2688	68	0
1	C	3119	0	2934	57	0
1	D	3098	0	2905	60	0
2	E	39	0	34	2	0
3	F	28	0	25	1	0
4	G	49	0	43	4	0
5	A	22	0	20	1	0
5	B	22	0	20	1	0
5	C	22	0	20	1	0
5	D	22	0	20	0	0
6	A	9	0	0	0	0
6	B	9	0	0	0	0
6	C	9	0	0	0	0
6	D	9	0	0	0	0
7	A	14	0	13	1	0
7	B	28	0	26	0	0
7	C	14	0	13	0	0
7	D	14	0	13	0	0
All	All	12505	0	11505	249	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 10.

All (249) 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:D:154:VAL:HG12	1:D:164:PRO:HA	1.47	0.96
1:D:41:ARG:N	1:D:44:SER:OG	2.04	0.90
1:B:364:PHE:HA	1:B:378:THR:HA	1.51	0.89
1:C:46:LEU:HD11	1:C:62:ILE:HG12	1.62	0.81
1:A:369:LEU:HD23	1:A:370:VAL:H	1.47	0.79
1:B:34:ARG:HA	1:B:84:ASP:HB2	1.65	0.78
1:D:323:SER:HB3	1:D:340:LEU:HB3	1.65	0.78
1:B:21:ALA:HB2	1:B:31:LEU:HD11	1.67	0.76
1:A:48:SER:HB2	1:A:59:ALA:HB2	1.70	0.74
2:E:1:NAG:H62	2:E:2:NAG:HN2	1.52	0.74
1:B:256:VAL:HG13	1:B:257:ASP:H	1.52	0.73
1:D:334:PRO:HA	1:D:378:THR:HG23	1.69	0.73
1:C:282:GLU:OE1	1:C:282:GLU:N	2.22	0.73
1:C:41:ARG:N	1:C:44:SER:OG	2.20	0.71
1:B:282:GLU:N	1:B:282:GLU:OE1	2.25	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:282:GLU:N	1:D:282:GLU:OE1	2.25	0.70
1:B:144:LEU:H	1:B:152:LEU:HD22	1.57	0.69
1:A:282:GLU:OE1	1:A:282:GLU:N	2.24	0.69
1:A:34:ARG:NH2	1:A:84:ASP:OD2	2.26	0.69
1:C:224:PRO:HA	1:C:310:LEU:HB2	1.75	0.67
4:G:1:NAG:H61	4:G:2:NAG:C7	2.26	0.66
1:D:260:VAL:HG13	1:D:287:TYR:HD2	1.61	0.66
1:B:64:ARG:NH2	1:B:69:ALA:O	2.29	0.65
1:C:47:PHE:HB3	1:C:56:LEU:HD11	1.80	0.64
1:A:240:ASP:OD2	1:A:247:GLY:HA2	1.98	0.63
1:A:5:TYR:HB3	1:A:17:VAL:HG12	1.79	0.63
1:D:31:LEU:HD21	1:D:54:GLY:HA3	1.80	0.63
1:A:155:GLN:HE21	1:A:155:GLN:HA	1.64	0.62
1:A:343:VAL:HG23	1:A:354:ILE:HD12	1.81	0.62
1:D:185:THR:HG22	1:D:197:THR:HG22	1.82	0.62
1:C:9:GLU:HG3	1:C:98:ILE:HG12	1.80	0.62
4:G:1:NAG:H3	4:G:1:NAG:H83	1.81	0.62
1:C:126:LEU:HB2	1:C:164:PRO:HG2	1.82	0.61
1:B:148:PRO:HB2	1:B:170:ARG:HD3	1.83	0.61
1:C:154:VAL:HG11	1:C:161:PRO:HB3	1.83	0.61
1:C:43:ARG:HH12	1:C:61:ARG:HG3	1.66	0.60
1:B:28:THR:O	1:B:32:MET:HG2	2.01	0.59
1:B:41:ARG:NH2	1:B:78:PHE:HA	2.17	0.59
1:D:392:ILE:O	1:D:409:HIS:HA	2.02	0.59
1:A:323:SER:HB2	1:A:340:LEU:HB3	1.86	0.58
1:D:215:THR:HG1	1:D:219:TYR:HH	1.47	0.58
7:A:515:NAG:O7	7:A:515:NAG:O3	2.22	0.58
1:C:298:GLY:HA3	1:D:205:VAL:HG21	1.85	0.58
1:C:123:ARG:HG2	1:C:167:VAL:HG22	1.86	0.58
1:D:172:LEU:HD13	1:D:204:VAL:HG22	1.86	0.57
1:D:46:LEU:O	1:D:59:ALA:N	2.33	0.57
1:B:126:LEU:HD13	1:B:184:LEU:HD22	1.85	0.57
1:B:193:ILE:O	5:B:502:MAN:O6	2.19	0.57
1:B:16:PHE:HA	1:B:57:VAL:HG12	1.86	0.57
1:A:5:TYR:HB3	1:A:17:VAL:CG1	2.36	0.56
1:B:369:LEU:HB3	1:B:373:TYR:O	2.05	0.56
1:A:151:SER:OG	1:A:169:GLN:NE2	2.31	0.56
1:D:233:LEU:HG	1:D:274:SER:HA	1.87	0.56
1:B:345:ASP:N	1:B:354:ILE:HD11	2.21	0.56
1:A:98:ILE:HG22	1:A:99:ASN:H	1.71	0.56
1:B:287:TYR:HB2	1:B:307:VAL:HB	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:43:ARG:HD3	1:C:46:LEU:HD12	1.89	0.55
1:D:254:HIS:CD2	1:D:292:GLN:HB2	2.41	0.55
1:D:343:VAL:O	1:D:372:ASN:HB3	2.07	0.55
1:A:392:ILE:O	1:A:409:HIS:HA	2.07	0.55
1:A:396:ALA:O	1:A:405:SER:HA	2.07	0.55
1:D:126:LEU:HD13	1:D:184:LEU:HD22	1.90	0.54
1:B:134:VAL:HA	1:B:138:SER:HB3	1.90	0.54
1:B:345:ASP:OD2	1:B:352:GLY:HA2	2.07	0.54
1:C:46:LEU:O	1:C:59:ALA:N	2.33	0.54
1:B:345:ASP:HB2	1:B:354:ILE:HD11	1.90	0.54
1:D:64:ARG:NH2	1:D:69:ALA:O	2.41	0.54
1:D:36:ILE:HG22	1:D:82:MET:HG2	1.90	0.54
1:A:322:THR:HA	1:B:41:ARG:HD3	1.89	0.53
1:D:2:ASN:HA	1:D:90:PRO:O	2.07	0.53
1:A:4:ARG:NH2	1:A:92:GLU:OE1	2.34	0.53
1:C:337:THR:HG22	1:C:375:ARG:HD2	1.89	0.53
1:C:153:ASP:O	1:C:165:GLU:N	2.36	0.53
1:D:41:ARG:H	1:D:44:SER:HG	1.51	0.53
1:A:41:ARG:N	1:A:44:SER:OG	2.41	0.53
1:A:111:GLU:HG2	1:A:201:HIS:HB3	1.90	0.53
1:B:264:PHE:HB3	1:B:273:LEU:HD11	1.92	0.52
1:D:2:ASN:OD1	1:D:90:PRO:HB2	2.09	0.52
1:A:335:GLY:N	1:A:378:THR:O	2.42	0.52
1:A:286:VAL:HG12	1:A:308:THR:HG23	1.91	0.52
1:C:116:GLU:HA	1:C:172:LEU:HB2	1.91	0.52
1:B:120:PRO:HG2	1:D:120:PRO:HG2	1.91	0.51
1:D:31:LEU:HD21	1:D:54:GLY:CA	2.41	0.51
1:B:4:ARG:HA	1:B:92:GLU:O	2.10	0.51
1:D:50:ASN:ND2	1:D:53:SER:OG	2.44	0.51
1:A:398:ASP:OD1	1:A:404:LEU:N	2.31	0.51
1:C:207:VAL:HG13	1:D:207:VAL:HG13	1.91	0.51
1:B:323:SER:HB2	1:B:340:LEU:H	1.75	0.51
1:D:368:LYS:HG2	1:D:374:TYR:HE1	1.74	0.51
1:A:240:ASP:OD2	1:A:242:ASP:HB2	2.11	0.51
1:A:316:ALA:HB2	1:A:404:LEU:HB3	1.92	0.51
1:C:115:SER:HB2	1:D:297:ALA:O	2.11	0.51
1:C:17:VAL:HB	1:C:56:LEU:O	2.10	0.50
1:B:68:CYS:HB3	1:B:71:SER:OG	2.11	0.50
1:B:343:VAL:HG12	1:B:354:ILE:HD13	1.92	0.50
1:C:48:SER:HB2	1:C:59:ALA:HB2	1.93	0.50
1:D:111:GLU:HG2	1:D:201:HIS:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:313:ASN:ND2	1:D:398:ASP:OD2	2.44	0.50
1:D:368:LYS:HG2	1:D:374:TYR:CE1	2.47	0.50
1:A:185:THR:HG22	1:A:197:THR:HG22	1.94	0.50
3:F:1:NAG:H61	3:F:2:NAG:N2	2.27	0.50
1:D:361:ASN:OD1	1:D:362:LEU:N	2.44	0.49
1:C:43:ARG:NH1	1:C:61:ARG:HG3	2.27	0.49
1:A:62:ILE:HG21	1:A:95:ILE:HD12	1.94	0.49
1:D:383:ASP:HB2	1:D:386:GLN:HB3	1.94	0.49
1:C:361:ASN:OD1	1:C:362:LEU:N	2.45	0.49
1:A:126:LEU:HD13	1:A:184:LEU:HD22	1.93	0.49
1:B:340:LEU:HD12	1:B:374:TYR:O	2.12	0.49
1:A:319:VAL:HA	1:A:343:VAL:HG12	1.94	0.49
1:A:393:THR:HG22	1:A:409:HIS:ND1	2.28	0.49
1:A:41:ARG:HD3	1:B:322:THR:HB	1.95	0.49
1:C:134:VAL:HG12	1:C:135:GLY:H	1.76	0.49
1:C:293:ALA:O	1:C:300:MET:HA	2.13	0.49
1:C:384:ARG:HH21	1:C:415:THR:HA	1.78	0.48
1:D:357:SER:HB2	1:D:395:THR:OG1	2.12	0.48
1:C:29:ARG:HG3	1:C:30:GLU:H	1.77	0.48
1:C:9:GLU:CG	1:C:98:ILE:HG12	2.43	0.48
1:C:43:ARG:HG2	1:C:46:LEU:HB2	1.94	0.48
1:D:215:THR:HG22	4:G:4:FUC:H61	1.95	0.48
1:B:10:GLU:N	1:B:62:ILE:O	2.36	0.48
1:A:361:ASN:OD1	1:A:362:LEU:N	2.46	0.48
1:C:267:ASP:OD1	1:C:270:THR:N	2.34	0.48
1:A:293:ALA:O	1:A:300:MET:HA	2.14	0.48
1:C:233:LEU:HG	1:C:274:SER:HA	1.93	0.48
1:B:313:ASN:ND2	1:B:402:PRO:O	2.47	0.47
1:D:36:ILE:HD12	1:D:49:LEU:HD21	1.95	0.47
1:B:48:SER:HB2	1:B:59:ALA:HB2	1.95	0.47
1:B:41:ARG:HH22	1:B:78:PHE:HA	1.77	0.47
1:B:343:VAL:HB	1:B:372:ASN:O	2.13	0.47
1:B:19:SER:HB3	1:B:22:LYS:HG2	1.97	0.47
1:B:254:HIS:CD2	1:B:292:GLN:HB2	2.49	0.47
1:C:8:PRO:HA	1:C:96:ILE:HB	1.97	0.47
1:B:260:VAL:HG13	1:B:287:TYR:HD2	1.78	0.47
1:D:328:VAL:HG22	1:D:329:PRO:HD2	1.96	0.47
1:A:134:VAL:HG12	1:A:135:GLY:N	2.30	0.47
1:B:233:LEU:HG	1:B:274:SER:HA	1.97	0.47
2:E:1:NAG:H62	2:E:2:NAG:N2	2.26	0.46
1:C:31:LEU:HG	1:C:36:ILE:HG12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:402:PRO:HA	1:C:403:PRO:HD3	1.85	0.46
1:D:343:VAL:CG1	1:D:374:TYR:HE2	2.27	0.46
1:C:195:SER:HB2	5:C:501:MAN:H2	1.48	0.46
1:A:31:LEU:HB2	1:A:36:ILE:HD12	1.97	0.46
1:B:19:SER:HB3	1:B:22:LYS:HE3	1.97	0.46
1:B:329:PRO:HG2	1:B:332:PHE:CB	2.46	0.46
1:D:120:PRO:HG3	1:D:171:PRO:HG3	1.97	0.46
1:A:4:ARG:HA	1:A:92:GLU:O	2.16	0.46
1:D:134:VAL:HA	1:D:138:SER:HB3	1.97	0.46
1:B:185:THR:HG22	1:B:197:THR:HG22	1.99	0.45
1:B:218:GLU:N	1:B:218:GLU:OE1	2.49	0.45
1:C:29:ARG:HG3	1:C:30:GLU:N	2.32	0.45
1:D:41:ARG:O	1:D:44:SER:OG	2.28	0.45
1:D:295:ASP:OD2	1:D:299:LEU:HB2	2.17	0.45
1:B:402:PRO:HA	1:B:403:PRO:HD3	1.87	0.45
1:A:36:ILE:HG23	1:A:80:ILE:HG23	1.98	0.45
1:A:224:PRO:HA	1:A:310:LEU:HB2	1.98	0.45
1:A:233:LEU:HG	1:A:274:SER:HA	1.99	0.45
1:C:134:VAL:HA	1:C:138:SER:HB3	1.98	0.45
1:C:254:HIS:CD2	1:C:292:GLN:HB2	2.51	0.45
1:C:218:GLU:HA	1:C:304:LYS:O	2.16	0.45
1:B:99:ASN:HB2	1:B:137:ASN:CG	2.41	0.45
1:C:369:LEU:HG	1:C:370:VAL:H	1.82	0.45
1:A:155:GLN:HA	1:A:155:GLN:NE2	2.31	0.44
1:B:219:TYR:HB3	1:B:234:LEU:HD21	1.99	0.44
1:D:143:GLN:NE2	1:D:185:THR:O	2.50	0.44
1:A:398:ASP:O	1:A:403:PRO:HB3	2.17	0.44
1:B:219:TYR:HB3	1:B:234:LEU:CD2	2.47	0.44
1:C:179:VAL:HG13	1:C:201:HIS:CE1	2.51	0.44
1:D:343:VAL:HG11	1:D:374:TYR:HE2	1.83	0.44
1:A:343:VAL:CG2	1:A:354:ILE:HD12	2.47	0.44
1:C:184:LEU:HD23	1:C:198:LEU:HD23	1.99	0.44
1:D:123:ARG:NE	1:D:165:GLU:OE1	2.43	0.44
1:B:361:ASN:OD1	1:B:362:LEU:N	2.50	0.44
1:B:175:GLU:OE1	1:B:175:GLU:N	2.45	0.44
1:C:134:VAL:HG12	1:C:135:GLY:N	2.32	0.44
1:A:228:PRO:HG2	1:A:231:THR:OG1	2.17	0.44
1:D:318:GLU:O	1:D:343:VAL:HA	2.18	0.44
1:D:328:VAL:CG2	1:D:329:PRO:HD2	2.48	0.44
1:D:76:VAL:O	1:D:92:GLU:HA	2.18	0.44
1:A:384:ARG:HG3	1:A:414:VAL:HB	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:TYR:HB2	1:B:305:VAL:HG22	2.00	0.44
1:D:345:ASP:OD2	1:D:352:GLY:HA2	2.18	0.44
1:A:139:LEU:HD11	1:A:186:ALA:HB1	2.00	0.43
1:A:3:ILE:HG21	1:A:23:ASP:OD2	2.18	0.43
1:C:186:ALA:O	1:C:195:SER:HA	2.18	0.43
1:D:46:LEU:HD23	1:D:46:LEU:HA	1.67	0.43
1:B:225:GLU:OE1	1:B:312:VAL:O	2.35	0.43
1:C:10:GLU:N	1:C:62:ILE:O	2.49	0.43
1:A:112:LEU:HD13	1:A:124:ILE:HD13	2.00	0.43
1:A:369:LEU:CD2	1:A:370:VAL:H	2.23	0.43
1:B:27:GLU:OE2	1:B:27:GLU:N	2.51	0.43
1:B:112:LEU:HD13	1:B:124:ILE:HD13	2.00	0.43
1:A:235:LYS:NZ	1:A:270:THR:O	2.52	0.43
1:A:149:HIS:O	1:A:168:LEU:HD12	2.19	0.43
1:C:126:LEU:HD13	1:C:184:LEU:HD22	2.00	0.43
1:B:17:VAL:HB	1:B:56:LEU:O	2.19	0.43
1:B:111:GLU:HG2	1:B:201:HIS:HB3	2.01	0.43
1:A:354:ILE:O	1:A:354:ILE:HG13	2.19	0.42
1:C:34:ARG:HA	1:C:84:ASP:HB2	2.01	0.42
1:D:185:THR:HA	1:D:196:GLY:O	2.19	0.42
1:A:8:PRO:HA	1:A:96:ILE:HB	2.01	0.42
1:C:5:TYR:OH	1:C:23:ASP:OD2	2.31	0.42
1:B:378:THR:HG22	1:B:380:ARG:H	1.84	0.42
1:A:10:GLU:OE1	1:A:61:ARG:NE	2.48	0.42
1:B:26:LEU:HD13	1:B:34:ARG:NE	2.34	0.42
1:D:283:GLU:HG2	1:D:284:TYR:CD2	2.54	0.42
1:B:56:LEU:HD21	1:B:78:PHE:HZ	1.83	0.42
1:A:282:GLU:O	1:A:283:GLU:HB3	2.20	0.42
1:B:281:PHE:CE1	1:B:285:LYS:HG3	2.55	0.42
1:C:109:GLU:HA	1:C:199:GLN:O	2.19	0.42
1:C:16:PHE:HA	1:C:57:VAL:HG12	2.00	0.42
1:A:99:ASN:HB2	1:A:137:ASN:CG	2.45	0.42
1:A:133:ASP:HB2	1:A:137:ASN:O	2.19	0.42
1:A:254:HIS:CD2	1:A:292:GLN:HB2	2.55	0.42
1:A:302:ARG:NH1	1:B:125:PRO:HD3	2.35	0.42
1:B:345:ASP:CB	1:B:354:ILE:HD11	2.49	0.42
1:C:219:TYR:HB3	1:C:234:LEU:HD23	2.01	0.42
1:C:357:SER:HB2	1:C:395:THR:OG1	2.20	0.42
1:B:77:SER:OG	1:B:92:GLU:HG2	2.19	0.41
1:D:116:GLU:HA	1:D:172:LEU:HB2	2.01	0.41
1:A:41:ARG:CD	1:B:322:THR:HB	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:ARG:HD3	1:B:82:MET:SD	2.61	0.41
1:A:143:GLN:HB2	1:A:185:THR:HG1	1.84	0.41
1:A:322:THR:O	1:B:41:ARG:NH1	2.50	0.41
1:C:205:VAL:HG11	1:D:297:ALA:HA	2.02	0.41
1:C:299:LEU:HD23	1:C:299:LEU:HA	1.85	0.41
1:A:3:ILE:HB	1:A:91:ILE:CD1	2.50	0.41
1:A:32:MET:HE3	1:A:32:MET:HA	2.02	0.41
1:D:401:THR:HA	1:D:403:PRO:HD3	2.02	0.41
1:B:9:GLU:HG3	1:B:98:ILE:HG12	2.02	0.41
1:B:70:GLN:HA	1:B:134:VAL:HG11	2.03	0.41
1:B:302:ARG:HE	1:B:302:ARG:HB3	1.69	0.41
1:B:349:ASP:CG	1:B:350:ASN:H	2.29	0.41
1:C:224:PRO:HG3	1:C:310:LEU:HD12	2.01	0.41
1:C:256:VAL:CG2	1:D:125:PRO:HB2	2.50	0.41
1:A:295:ASP:OD2	1:A:299:LEU:HB2	2.20	0.41
1:D:110:LEU:HD22	1:D:127:PRO:HD3	2.02	0.41
1:A:114:MET:O	1:A:204:VAL:HA	2.21	0.41
1:A:318:GLU:O	1:A:343:VAL:HA	2.21	0.41
1:D:299:LEU:HD23	1:D:299:LEU:HA	1.95	0.41
1:C:318:GLU:O	1:C:343:VAL:HA	2.21	0.41
1:A:195:SER:HB2	5:A:501:MAN:H2	1.73	0.40
1:A:225:GLU:OE1	1:A:312:VAL:O	2.38	0.40
1:A:401:THR:HA	1:A:403:PRO:HD3	2.04	0.40
1:B:399:GLN:O	1:B:399:GLN:HG2	2.21	0.40
1:C:24:LEU:HD11	1:C:89:LEU:HD22	2.03	0.40
1:D:48:SER:HB2	1:D:59:ALA:HB2	2.04	0.40
1:A:46:LEU:C	1:A:59:ALA:HB3	2.47	0.40
4:G:2:NAG:O3	4:G:3:BMA:O5	2.34	0.40
1:D:362:LEU:HD13	1:D:364:PHE:CZ	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	411/427 (96%)	388 (94%)	23 (6%)	0	100	100
1	B	410/427 (96%)	388 (95%)	22 (5%)	0	100	100
1	C	416/427 (97%)	393 (94%)	23 (6%)	0	100	100
1	D	414/427 (97%)	392 (95%)	22 (5%)	0	100	100
All	All	1651/1708 (97%)	1561 (94%)	90 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	312/380 (82%)	311 (100%)	1 (0%)	86	84
1	B	303/380 (80%)	302 (100%)	1 (0%)	86	84
1	C	340/380 (90%)	339 (100%)	1 (0%)	86	84
1	D	338/380 (89%)	337 (100%)	1 (0%)	86	84
All	All	1293/1520 (85%)	1289 (100%)	4 (0%)	86	84

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	369	LEU
1	B	369	LEU
1	C	46	LEU
1	D	143	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	147	ASN
1	A	155	GLN

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Mol	Chain	Res	Type
1	A	199	GLN
1	A	210	ASN
1	B	79	ASN
1	B	246	ASN
1	B	254	HIS
1	B	294	GLN
1	C	180	HIS
1	C	210	ASN
1	D	106	GLN
1	D	137	ASN
1	D	210	ASN
1	D	254	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](#)

9 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	E	1	1,2	14,14,15	0.35	0	17,19,21	0.48	0
2	NAG	E	2	2	14,14,15	0.22	0	17,19,21	0.44	0
2	BMA	E	3	2	11,11,12	0.89	0	15,15,17	0.88	1 (6%)
3	NAG	F	1	1,3	14,14,15	0.51	0	17,19,21	0.73	0
3	NAG	F	2	3	14,14,15	0.28	0	17,19,21	0.42	0
4	NAG	G	1	1,4	14,14,15	0.86	1 (7%)	17,19,21	1.42	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	G	2	4	14,14,15	0.19	0	17,19,21	0.77	0
4	BMA	G	3	4	11,11,12	0.96	1 (9%)	15,15,17	1.30	2 (13%)
4	FUC	G	4	4	10,10,11	0.80	0	14,14,16	0.83	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	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	BMA	E	3	2	-	1/2/19/22	0/1/1/1
3	NAG	F	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
4	NAG	G	1	1,4	-	4/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
4	BMA	G	3	4	-	0/2/19/22	0/1/1/1
4	FUC	G	4	4	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	3	BMA	C1-C2	2.79	1.58	1.52
4	G	1	NAG	O5-C1	-2.66	1.39	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	1	NAG	C2-N2-C7	4.52	128.96	122.90
4	G	3	BMA	C1-C2-C3	2.89	113.86	109.64
4	G	3	BMA	C1-O5-C5	2.41	115.42	112.19
4	G	1	NAG	C1-O5-C5	2.28	115.24	112.19
2	E	3	BMA	C2-C3-C4	2.08	114.51	110.86

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	2	NAG	O5-C5-C6-O6

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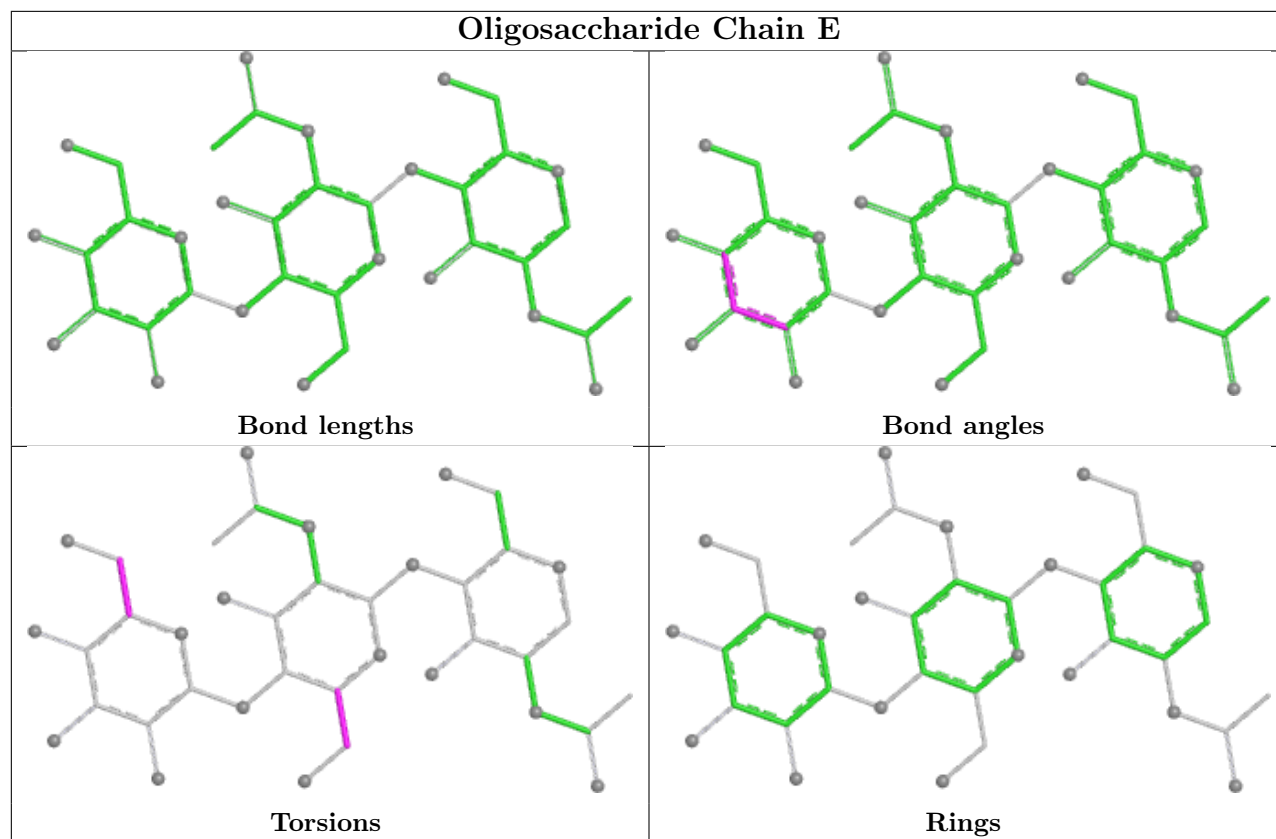
Mol	Chain	Res	Type	Atoms
4	G	1	NAG	C8-C7-N2-C2
4	G	1	NAG	O7-C7-N2-C2
3	F	2	NAG	C4-C5-C6-O6
2	E	3	BMA	O5-C5-C6-O6
4	G	2	NAG	C3-C2-N2-C7
2	E	2	NAG	C4-C5-C6-O6
3	F	1	NAG	C3-C2-N2-C7
4	G	1	NAG	C3-C2-N2-C7
2	E	2	NAG	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
3	F	1	NAG	C1-C2-N2-C7
4	G	1	NAG	C1-C2-N2-C7

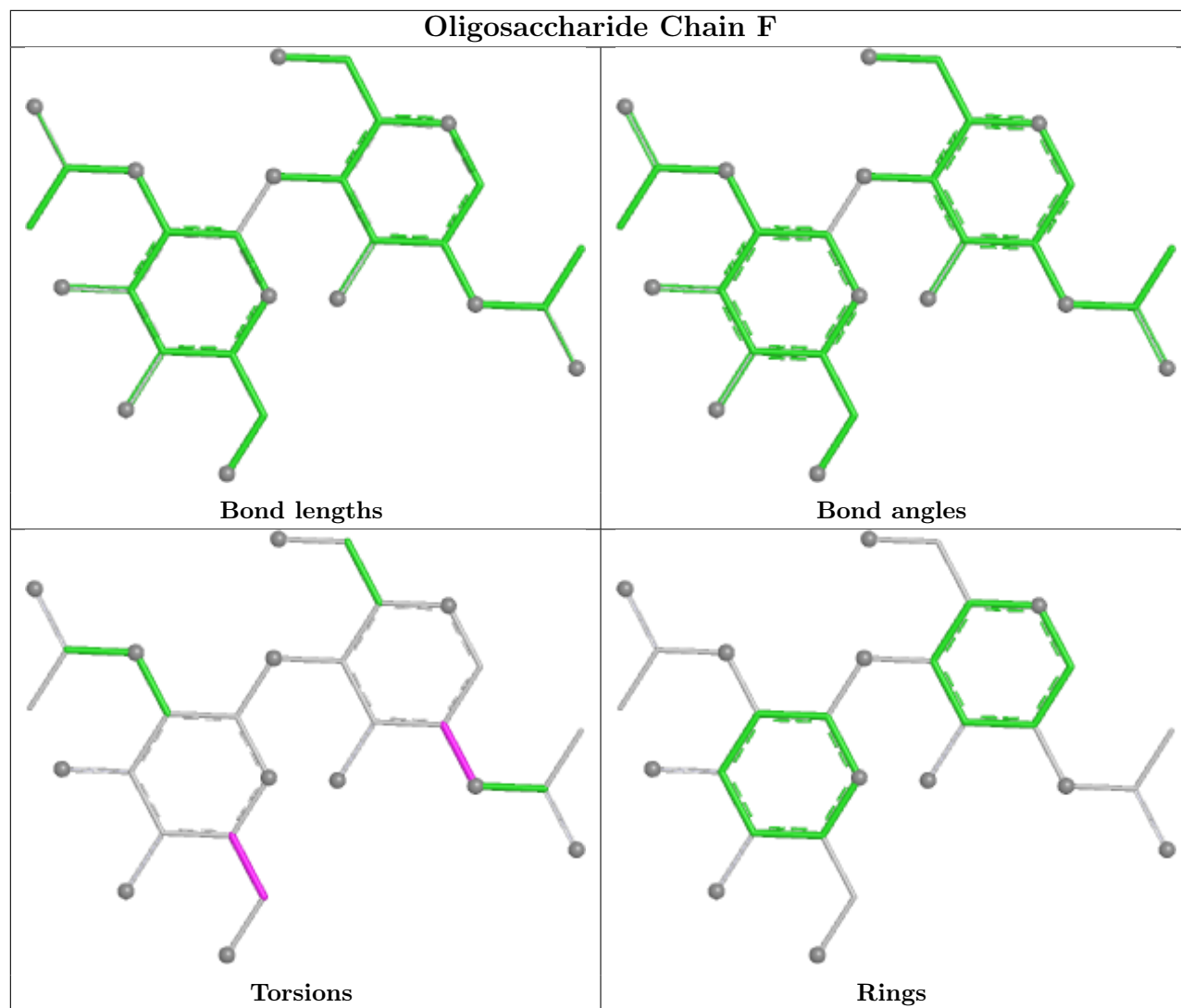
There are no ring outliers.

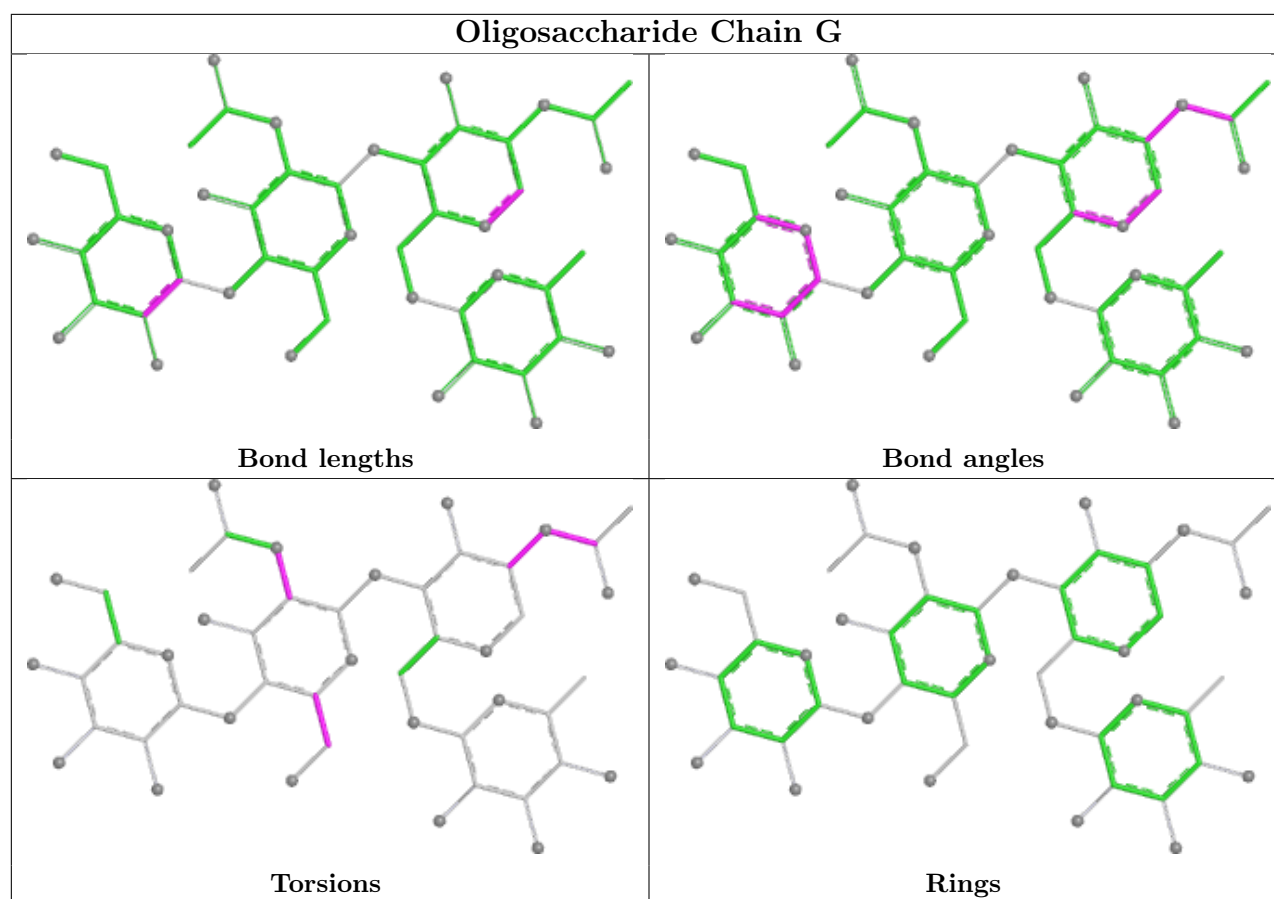
8 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	4	FUC	1	0
3	F	2	NAG	1	0
4	G	2	NAG	2	0
2	E	2	NAG	2	0
4	G	1	NAG	2	0
4	G	3	BMA	1	0
3	F	1	NAG	1	0
2	E	1	NAG	2	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](#)

Of 49 ligands modelled in this entry, 36 are monoatomic - leaving 13 for Mogul analysis.

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	$\# Z > 2$
7	NAG	D	516	1	14,14,15	0.31	0	17,19,21	0.45	0
5	MAN	C	502	1	11,11,12	0.86	1 (9%)	15,15,17	0.85	1 (6%)
7	NAG	C	514	1	14,14,15	0.32	0	17,19,21	0.45	0
5	MAN	D	501	1	11,11,12	0.69	0	15,15,17	0.94	1 (6%)
5	MAN	D	502	1	11,11,12	0.81	0	15,15,17	0.86	1 (6%)
7	NAG	B	513	1	14,14,15	0.57	0	17,19,21	0.48	0
5	MAN	C	501	1	11,11,12	0.81	1 (9%)	15,15,17	0.92	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MAN	A	501	1	11,11,12	0.78	0	15,15,17	0.92	2 (13%)
7	NAG	B	512	1	14,14,15	0.27	0	17,19,21	0.44	0
5	MAN	A	502	1	11,11,12	0.94	1 (9%)	15,15,17	0.84	1 (6%)
5	MAN	B	501	1	11,11,12	0.68	0	15,15,17	0.95	2 (13%)
5	MAN	B	502	1	11,11,12	0.85	0	15,15,17	0.78	0
7	NAG	A	515	1	14,14,15	0.30	0	17,19,21	0.51	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
7	NAG	D	516	1	-	2/6/23/26	0/1/1/1
5	MAN	C	502	1	-	0/2/19/22	0/1/1/1
7	NAG	C	514	1	-	2/6/23/26	0/1/1/1
5	MAN	D	501	1	-	0/2/19/22	0/1/1/1
5	MAN	D	502	1	-	0/2/19/22	0/1/1/1
7	NAG	B	513	1	-	3/6/23/26	0/1/1/1
5	MAN	C	501	1	-	0/2/19/22	0/1/1/1
5	MAN	A	501	1	-	0/2/19/22	0/1/1/1
7	NAG	B	512	1	-	2/6/23/26	0/1/1/1
5	MAN	A	502	1	-	0/2/19/22	0/1/1/1
5	MAN	B	501	1	-	0/2/19/22	0/1/1/1
5	MAN	B	502	1	-	0/2/19/22	0/1/1/1
7	NAG	A	515	1	-	4/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	502	MAN	O5-C1	-2.45	1.39	1.43
5	C	501	MAN	O5-C1	-2.21	1.40	1.43
5	C	502	MAN	O5-C1	-2.16	1.40	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	501	MAN	O2-C2-C3	-2.43	105.13	110.15
5	C	502	MAN	O2-C2-C3	-2.27	105.45	110.15
5	C	501	MAN	O2-C2-C3	-2.23	105.53	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	501	MAN	C1-O5-C5	2.18	115.11	112.19
5	D	502	MAN	O2-C2-C3	-2.18	105.63	110.15
5	A	501	MAN	O2-C2-C3	-2.15	105.70	110.15
5	B	501	MAN	O2-C2-C3	-2.14	105.72	110.15
5	A	502	MAN	O2-C2-C3	-2.08	105.84	110.15
5	A	501	MAN	C1-O5-C5	2.00	114.87	112.19

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	515	NAG	C4-C5-C6-O6
7	B	513	NAG	C4-C5-C6-O6
7	C	514	NAG	O5-C5-C6-O6
7	B	512	NAG	O5-C5-C6-O6
7	A	515	NAG	O5-C5-C6-O6
7	B	513	NAG	O5-C5-C6-O6
7	C	514	NAG	C4-C5-C6-O6
7	B	512	NAG	C4-C5-C6-O6
7	D	516	NAG	C4-C5-C6-O6
7	A	515	NAG	C1-C2-N2-C7
7	A	515	NAG	C3-C2-N2-C7
7	B	513	NAG	C3-C2-N2-C7
7	D	516	NAG	O5-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	501	MAN	1	0
5	A	501	MAN	1	0
5	B	502	MAN	1	0
7	A	515	NAG	1	0

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 ⓘ

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	413/427 (96%)	0.43	22 (5%)	32	29	106, 201, 312, 366	0
1	B	412/427 (96%)	0.50	31 (7%)	20	22	112, 210, 310, 350	0
1	C	418/427 (97%)	0.52	38 (9%)	15	18	112, 163, 234, 279	0
1	D	416/427 (97%)	0.69	49 (11%)	9	12	119, 204, 263, 314	0
All	All	1659/1708 (97%)	0.54	140 (8%)	17	20	106, 188, 296, 366	0

All (140) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	263	LYS	6.1
1	B	89	LEU	6.0
1	D	76	VAL	5.2
1	C	166	MET	5.2
1	C	80	ILE	4.8
1	B	113	LYS	4.7
1	D	394	ILE	4.6
1	D	57	VAL	4.5
1	D	48	SER	4.4
1	D	49	LEU	4.4
1	C	127	PRO	4.2
1	D	93	VAL	4.2
1	D	80	ILE	4.1
1	D	91	ILE	3.8
1	B	279	LEU	3.8
1	B	287	TYR	3.8
1	D	56	LEU	3.8
1	B	39	VAL	3.8
1	C	124	ILE	3.8
1	C	293	ALA	3.7
1	C	202	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	47	PHE	3.6
1	D	393	THR	3.6
1	A	184	LEU	3.5
1	D	254	HIS	3.5
1	D	172	LEU	3.5
1	D	20	ILE	3.5
1	D	202	VAL	3.5
1	B	79	ASN	3.5
1	B	43	ARG	3.4
1	C	266	LEU	3.4
1	B	277	GLU	3.3
1	A	43	ARG	3.3
1	A	396	ALA	3.3
1	D	337	THR	3.3
1	A	249	VAL	3.3
1	C	125	PRO	3.2
1	C	20	ILE	3.2
1	B	264	PHE	3.2
1	C	114	MET	3.2
1	B	80	ILE	3.2
1	D	410	ILE	3.2
1	D	36	ILE	3.2
1	A	323	SER	3.1
1	B	117	ILE	3.1
1	C	78	PHE	3.1
1	C	253	PHE	3.1
1	B	112	LEU	3.1
1	A	251	TYR	3.0
1	C	184	LEU	3.0
1	D	114	MET	3.0
1	C	129	GLY	3.0
1	D	268	ALA	3.0
1	D	356	CYS	3.0
1	D	38	ILE	2.9
1	C	150	PHE	2.9
1	A	159	GLU	2.9
1	D	392	ILE	2.9
1	C	126	LEU	2.8
1	A	119	THR	2.8
1	B	326	ASN	2.8
1	A	319	VAL	2.8
1	D	54	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	408	ALA	2.8
1	B	114	MET	2.7
1	A	223	VAL	2.7
1	C	89	LEU	2.7
1	B	91	ILE	2.7
1	C	36	ILE	2.7
1	B	132	LEU	2.6
1	D	55	SER	2.6
1	A	93	VAL	2.6
1	B	172	LEU	2.6
1	D	7	VAL	2.6
1	D	266	LEU	2.6
1	C	27	GLU	2.6
1	D	27	GLU	2.6
1	A	76	VAL	2.6
1	D	408	ALA	2.6
1	D	110	LEU	2.5
1	D	25	GLY	2.5
1	B	110	LEU	2.5
1	C	337	THR	2.5
1	D	319	VAL	2.5
1	D	376	LEU	2.5
1	B	124	ILE	2.5
1	C	62	ILE	2.5
1	C	200	ILE	2.5
1	D	95	ILE	2.5
1	D	267	ASP	2.5
1	D	112	LEU	2.4
1	D	279	LEU	2.4
1	D	150	PHE	2.4
1	D	322	THR	2.4
1	C	91	ILE	2.4
1	D	166	MET	2.4
1	B	90	PRO	2.4
1	A	38	ILE	2.3
1	A	117	ILE	2.3
1	C	419	ASP	2.3
1	A	328	VAL	2.3
1	B	76	VAL	2.3
1	D	335	GLY	2.3
1	C	255	LYS	2.3
1	A	46	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	186	ALA	2.3
1	A	289	MET	2.3
1	B	163	GLN	2.3
1	B	88	LEU	2.2
1	C	335	GLY	2.2
1	B	402	PRO	2.2
1	D	37	ARG	2.2
1	A	62	ILE	2.2
1	B	291	ILE	2.2
1	D	117	ILE	2.2
1	B	166	MET	2.2
1	B	44	SER	2.2
1	A	291	ILE	2.2
1	B	202	VAL	2.1
1	A	356	CYS	2.1
1	D	374	TYR	2.1
1	C	142	TYR	2.1
1	C	143	GLN	2.1
1	C	106	GLN	2.1
1	C	164	PRO	2.1
1	C	82	MET	2.1
1	A	322	THR	2.1
1	C	412	LEU	2.1
1	D	18	GLY	2.1
1	D	17	VAL	2.1
1	C	110	LEU	2.0
1	B	207	VAL	2.0
1	C	191	SER	2.0
1	C	146	ALA	2.0
1	B	393	THR	2.0
1	D	82	MET	2.0
1	C	172	LEU	2.0
1	C	76	VAL	2.0
1	D	272	GLU	2.0
1	D	77	SER	2.0

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

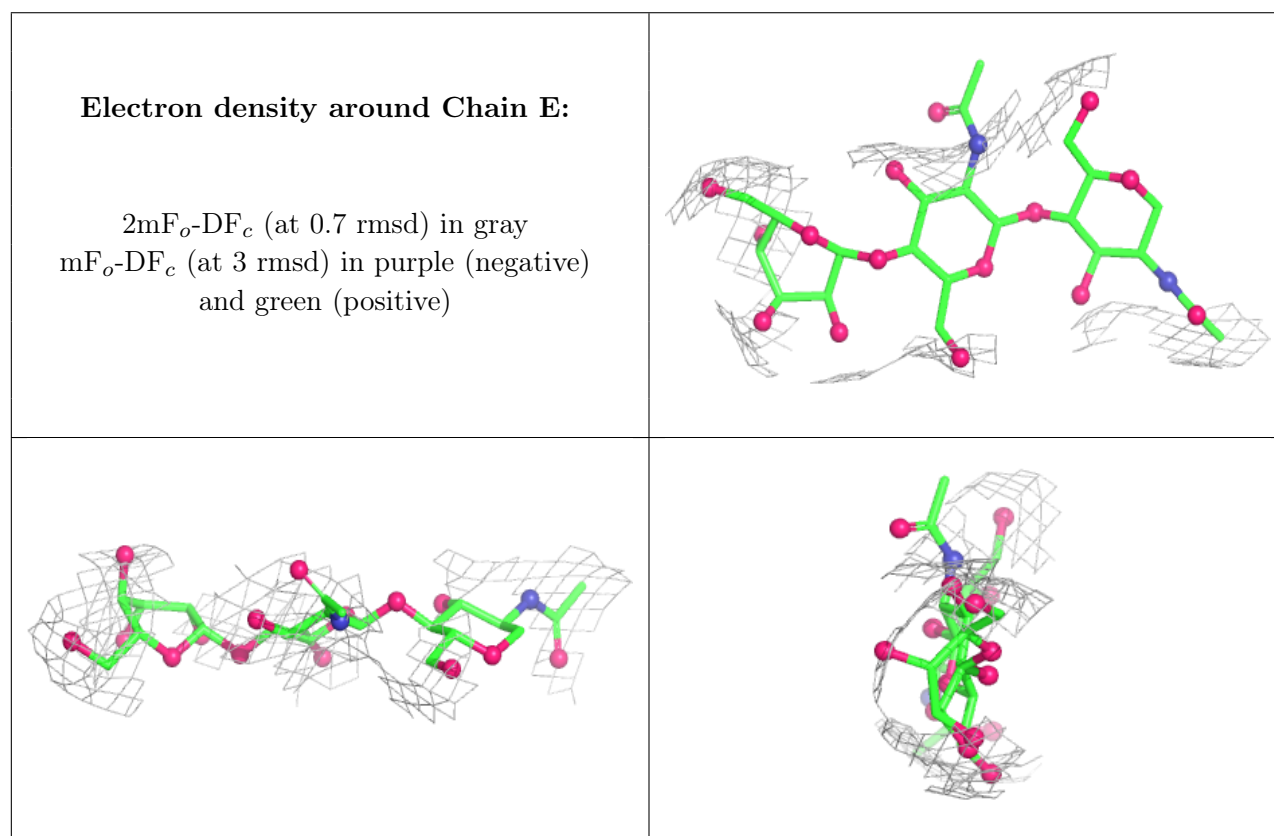
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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

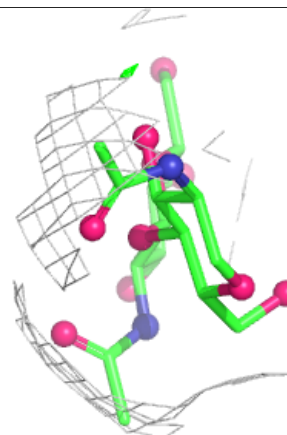
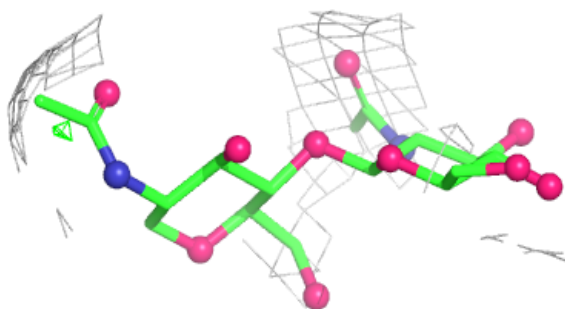
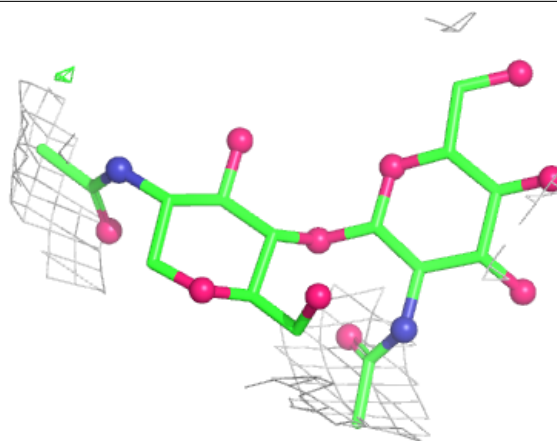
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	BMA	G	3	11/12	0.46	0.11	222,240,264,275	0
2	NAG	E	1	14/15	0.53	0.11	196,226,239,245	0
2	BMA	E	3	11/12	0.65	0.08	192,208,216,238	0
4	NAG	G	2	14/15	0.66	0.13	231,248,270,272	0
2	NAG	E	2	14/15	0.72	0.09	222,234,245,257	0
4	FUC	G	4	10/11	0.72	0.09	208,233,240,240	0
3	NAG	F	2	14/15	0.73	0.11	174,191,201,212	0
4	NAG	G	1	14/15	0.74	0.13	182,201,261,265	0
3	NAG	F	1	14/15	0.80	0.11	162,186,198,219	0

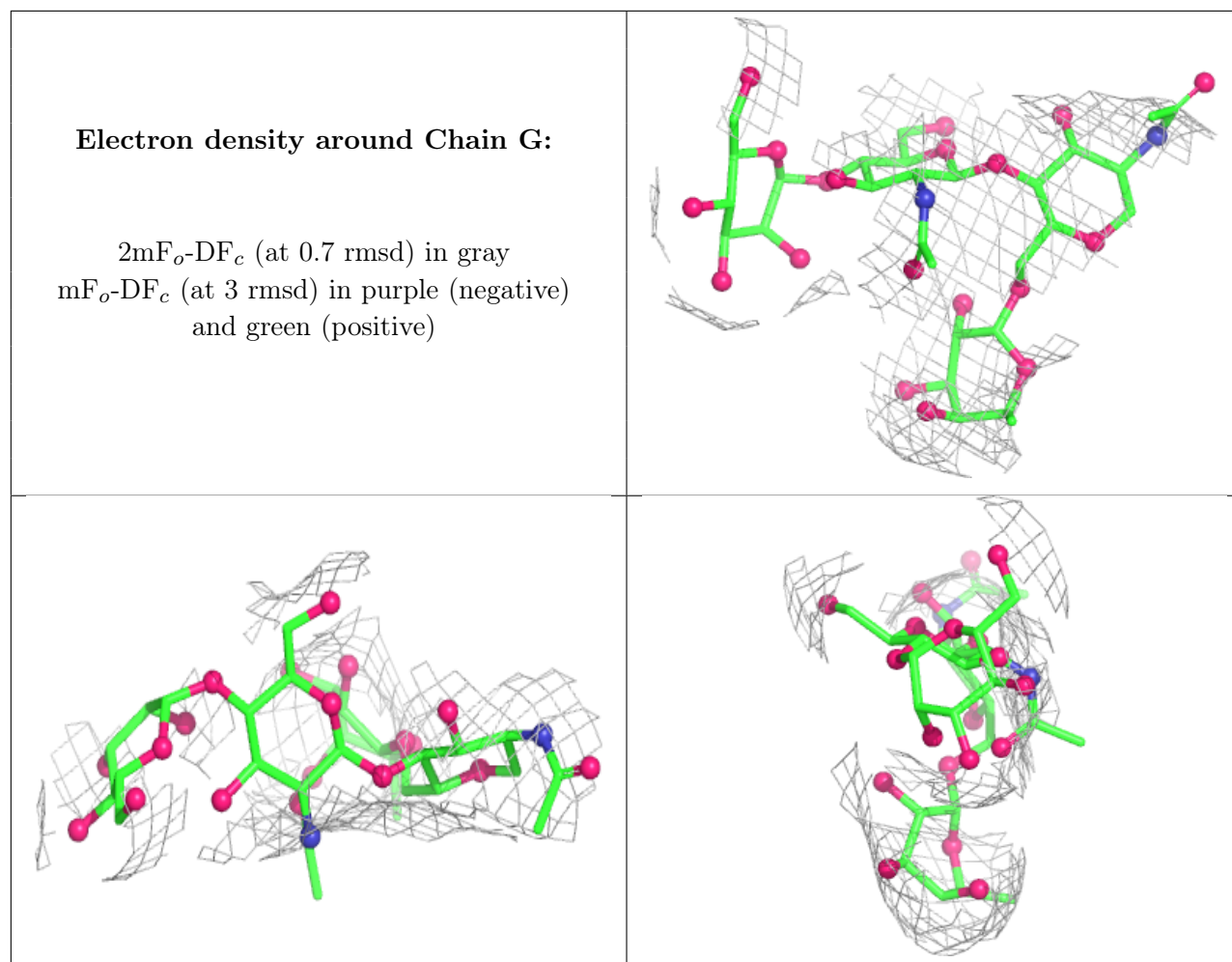
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around Chain F:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MAN	A	502	11/12	0.17	0.15	172,194,211,212	0
7	NAG	B	512	14/15	0.40	0.14	180,229,240,249	0
5	MAN	B	502	11/12	0.45	0.17	207,214,246,252	0
7	NAG	D	516	14/15	0.50	0.13	173,216,229,236	0
7	NAG	A	515	14/15	0.55	0.14	102,152,169,174	0
5	MAN	A	501	11/12	0.56	0.13	157,187,239,260	0
5	MAN	D	501	11/12	0.61	0.12	173,195,225,225	0
5	MAN	D	502	11/12	0.63	0.12	160,177,206,234	0
5	MAN	C	502	11/12	0.64	0.11	166,170,230,234	0
7	NAG	B	513	14/15	0.65	0.13	137,151,173,215	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MAN	C	501	11/12	0.75	0.12	144,161,174,234	0
5	MAN	B	501	11/12	0.75	0.09	155,167,186,195	0
6	CA	B	505	1/1	0.78	0.12	193,193,193,193	0
7	NAG	C	514	14/15	0.83	0.09	127,174,179,183	0
6	CA	D	509	1/1	0.84	0.12	192,192,192,192	0
6	CA	B	509	1/1	0.87	0.14	199,199,199,199	0
6	CA	C	508	1/1	0.88	0.10	171,171,171,171	0
6	CA	D	510	1/1	0.90	0.21	148,148,148,148	0
6	CA	C	505	1/1	0.90	0.10	97,97,97,97	0
6	CA	A	510	1/1	0.91	0.12	187,187,187,187	0
6	CA	A	505	1/1	0.91	0.14	95,95,95,95	0
6	CA	B	511	1/1	0.93	0.07	235,235,235,235	0
6	CA	A	508	1/1	0.94	0.04	239,239,239,239	0
6	CA	B	508	1/1	0.95	0.14	104,104,104,104	0
6	CA	B	507	1/1	0.95	0.08	154,154,154,154	0
6	CA	C	503	1/1	0.96	0.04	120,120,120,120	0
6	CA	B	506	1/1	0.96	0.04	96,96,96,96	0
6	CA	A	503	1/1	0.97	0.04	119,119,119,119	0
6	CA	A	506	1/1	0.97	0.05	198,198,198,198	0
6	CA	D	511	1/1	0.97	0.04	190,190,190,190	0
6	CA	A	511	1/1	0.97	0.11	150,150,150,150	0
6	CA	A	507	1/1	0.97	0.07	185,185,185,185	0
6	CA	D	504	1/1	0.97	0.04	175,175,175,175	0
6	CA	D	505	1/1	0.97	0.05	169,169,169,169	0
6	CA	D	508	1/1	0.97	0.09	100,100,100,100	0
6	CA	C	511	1/1	0.98	0.03	132,132,132,132	0
6	CA	B	510	1/1	0.98	0.06	145,145,145,145	0
6	CA	C	504	1/1	0.98	0.03	127,127,127,127	0
6	CA	C	507	1/1	0.99	0.07	95,95,95,95	0
6	CA	B	503	1/1	0.99	0.08	184,184,184,184	0
6	CA	C	509	1/1	0.99	0.07	158,158,158,158	0
6	CA	C	510	1/1	0.99	0.09	124,124,124,124	0
6	CA	B	504	1/1	0.99	0.05	209,209,209,209	0
6	CA	D	503	1/1	0.99	0.08	146,146,146,146	0
6	CA	A	504	1/1	0.99	0.02	130,130,130,130	0
6	CA	C	506	1/1	0.99	0.04	151,151,151,151	0
6	CA	D	506	1/1	0.99	0.04	129,129,129,129	0
6	CA	D	507	1/1	1.00	0.04	103,103,103,103	0
6	CA	A	509	1/1	1.00	0.03	110,110,110,110	0

6.5 Other polymers [i](#)

There are no such residues in this entry.