



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 06:32 PM UTC

PDB ID : 1NWR / pdb_00001nwr
Title : Crystal structure of human cartilage gp39 (HC-gp39)
Authors : Fusetti, F.; Pijning, T.; Kalk, K.H.; Bos, E.; Dijkstra, B.W.
Deposited on : 2003-02-06
Resolution : 2.70 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	:	FAILED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

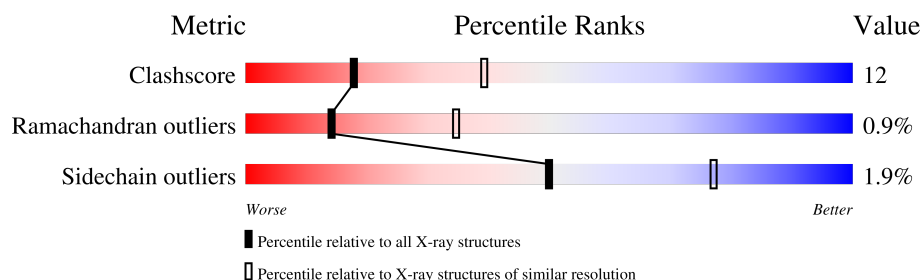
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)

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 $\leq 5\%$.

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	362	73% 25% .
1	B	362	73% 25% .
1	C	362	73% 25% .
1	D	362	70% 28% .
2	E	2	100%
2	F	2	50% 50%
2	G	2	50% 50%
2	H	2	100%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 11651 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 Chitinase-3 like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2857	1825	492	529	11			
1	B	362	Total	C	N	O	S	0	0	0
			2857	1825	492	529	11			
1	C	362	Total	C	N	O	S	0	0	0
			2857	1825	492	529	11			
1	D	362	Total	C	N	O	S	0	0	0
			2857	1825	492	529	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	311	ILE	THR	conflict	UNP P36222
B	311	ILE	THR	conflict	UNP P36222
C	311	ILE	THR	conflict	UNP P36222
D	311	ILE	THR	conflict	UNP P36222

- Molecule 2 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
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

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

- Molecule 3 is water.

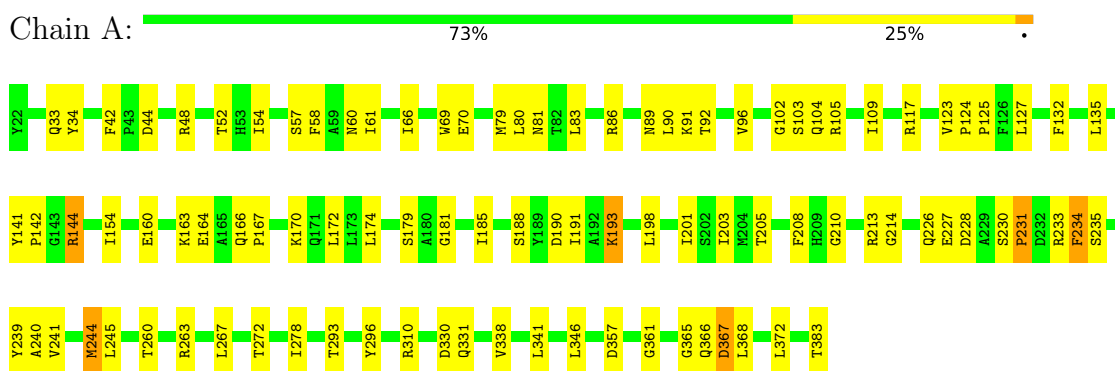
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	38	Total	O	0	0
			38	38		
3	B	16	Total	O	0	0
			16	16		
3	C	28	Total	O	0	0
			28	28		
3	D	29	Total	O	0	0
			29	29		

3 Residue-property plots

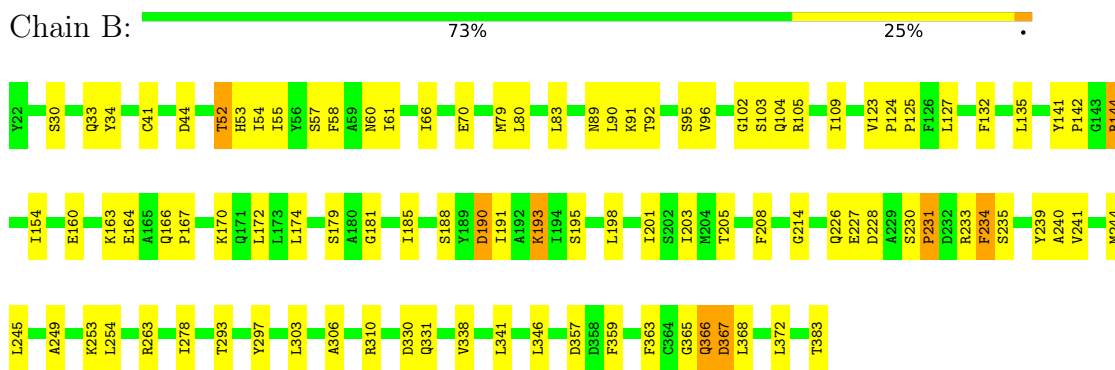
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS failed to run properly.

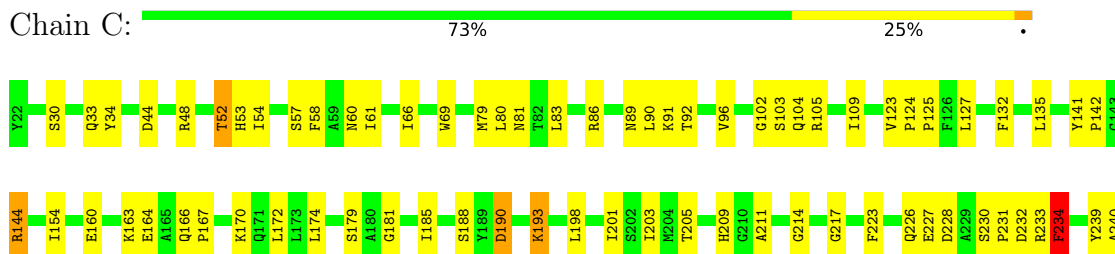
• Molecule 1: Chitinase-3 like protein 1



• Molecule 1: Chitinase-3 like protein 1



• Molecule 1: Chitinase-3 like protein 1





- Molecule 1: Chitinase-3 like protein 1

Chain D: 70% 28%



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

Chain E: 100%



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

Chain F: 50% 50%



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

Chain G: 50% 50%



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

Chain H: 100%



4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	127.30Å 127.30Å 107.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.79 – 2.70	Depositor
% Data completeness (in resolution range)	82.8 (19.79-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.212 , 0.239	Depositor
Wilson B-factor (Å ²)	25.7	Xtriage
Anisotropy	0.757	Xtriage
L-test for twinning ¹	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.054 for h,-k,-l	Xtriage
Total number of atoms	11651	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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 32.46 % 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 9.3595e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG

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	0/2933	0.90	3/3973 (0.1%)
1	B	0.44	0/2933	0.90	4/3973 (0.1%)
1	C	0.47	0/2933	0.93	6/3973 (0.2%)
1	D	0.49	0/2933	0.93	4/3973 (0.1%)
All	All	0.46	0/11732	0.91	17/15892 (0.1%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	233	ARG	N-CA-C	-8.94	102.09	113.72
1	D	233	ARG	N-CA-C	-8.84	102.49	113.28
1	A	233	ARG	N-CA-C	-7.25	104.43	113.28
1	B	233	ARG	N-CA-C	-6.73	104.88	113.23
1	B	357	ASP	N-CA-C	-6.32	102.38	110.53
1	D	190	ASP	N-CA-C	-6.32	100.14	109.11
1	D	357	ASP	N-CA-C	-6.24	102.48	110.53
1	B	190	ASP	N-CA-C	-6.17	100.35	109.11
1	A	357	ASP	N-CA-C	-6.09	102.68	110.53
1	C	357	ASP	N-CA-C	-6.00	102.79	110.53
1	C	190	ASP	N-CA-C	-5.99	100.60	109.11
1	B	52	THR	N-CA-C	-5.62	106.55	113.41
1	C	234	PHE	N-CA-C	5.57	120.80	112.94
1	C	211	ALA	N-CA-C	-5.52	106.50	113.18
1	C	52	THR	N-CA-C	-5.33	106.91	113.41
1	D	52	THR	N-CA-C	-5.09	107.20	113.41
1	A	52	THR	N-CA-C	-5.03	107.27	113.41

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	2857	0	2775	61	0
1	B	2857	0	2775	66	0
1	C	2857	0	2775	70	0
1	D	2857	0	2775	78	0
2	E	28	0	25	5	0
2	F	28	0	25	1	0
2	G	28	0	25	2	0
2	H	28	0	25	2	0
3	A	38	0	0	1	0
3	B	16	0	0	0	0
3	C	28	0	0	4	0
3	D	29	0	0	3	0
All	All	11651	0	11200	280	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 12.

All (280) 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:226:GLN:HE21	1:C:228:ASP:HB2	1.42	0.82
1:B:234:PHE:HB3	1:B:239:TYR:CD2	2.15	0.81
2:E:1:NAG:H61	2:E:2:NAG:H82	1.63	0.81
1:D:234:PHE:HB3	1:D:239:TYR:CD2	2.16	0.80
1:C:234:PHE:HB3	1:C:239:TYR:CD2	2.16	0.80
1:B:92:THR:HG22	1:B:132:PHE:HD1	1.46	0.80
1:D:226:GLN:HE21	1:D:228:ASP:HB2	1.48	0.79
1:D:92:THR:HG22	1:D:132:PHE:HD1	1.49	0.77
1:A:92:THR:HG22	1:A:132:PHE:HD1	1.48	0.77
1:C:92:THR:HG22	1:C:132:PHE:HD1	1.49	0.75
1:A:226:GLN:HE21	1:A:228:ASP:HB2	1.53	0.72
1:B:92:THR:HG22	1:B:132:PHE:CD1	2.27	0.70
1:C:92:THR:HG22	1:C:132:PHE:CD1	2.28	0.68
1:D:92:THR:HG22	1:D:132:PHE:CD1	2.28	0.68
1:D:210:GLY:O	1:D:213:ARG:HG3	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:1:NAG:H61	2:E:2:NAG:C7	2.24	0.67
1:A:92:THR:HG22	1:A:132:PHE:CD1	2.28	0.66
2:E:1:NAG:H61	2:E:2:NAG:C8	2.25	0.66
1:A:234:PHE:HB3	1:A:239:TYR:CD2	2.32	0.65
1:A:203:ILE:HD11	1:A:240:ALA:HB1	1.78	0.64
2:E:1:NAG:C6	2:E:2:NAG:H82	2.26	0.64
1:C:341:LEU:HB2	1:C:346:LEU:HD12	1.81	0.63
1:B:341:LEU:HB2	1:B:346:LEU:HD12	1.80	0.63
1:D:54:ILE:HD11	3:D:455:HOH:O	2.00	0.62
1:D:341:LEU:HB2	1:D:346:LEU:HD12	1.82	0.62
1:A:341:LEU:HB2	1:A:346:LEU:HD12	1.80	0.62
2:H:1:NAG:H62	2:H:2:NAG:C7	2.30	0.61
1:C:338:VAL:O	1:C:341:LEU:HG	2.00	0.61
1:C:234:PHE:HB3	1:C:239:TYR:CE2	2.35	0.60
1:B:44:ASP:HB3	1:B:79:MET:HG2	1.85	0.59
1:A:203:ILE:HG12	1:A:205:THR:HG23	1.83	0.59
1:B:203:ILE:HG12	1:B:205:THR:HG23	1.84	0.59
1:B:234:PHE:HB3	1:B:239:TYR:CE2	2.37	0.59
1:B:338:VAL:O	1:B:341:LEU:HG	2.03	0.59
1:D:44:ASP:HB3	1:D:79:MET:HG2	1.84	0.59
1:A:44:ASP:HB3	1:A:79:MET:HG2	1.85	0.58
1:D:338:VAL:O	1:D:341:LEU:HG	2.03	0.58
1:C:310:ARG:NH2	1:C:330:ASP:OD2	2.36	0.58
1:C:86:ARG:HD3	3:C:460:HOH:O	2.02	0.58
1:C:123:VAL:HB	1:C:124:PRO:HD3	1.85	0.58
1:C:124:PRO:HB2	1:C:125:PRO:HD3	1.86	0.58
1:D:124:PRO:HB2	1:D:125:PRO:HD3	1.86	0.58
1:C:198:LEU:HD13	1:C:201:ILE:CG2	2.34	0.58
1:A:198:LEU:HD13	1:A:201:ILE:CG2	2.34	0.58
1:A:310:ARG:NH2	1:A:330:ASP:OD2	2.36	0.57
1:B:310:ARG:NH2	1:B:330:ASP:OD2	2.38	0.57
1:B:166:GLN:N	1:B:167:PRO:HD2	2.20	0.57
1:D:44:ASP:HB3	1:D:79:MET:CG	2.35	0.57
1:C:44:ASP:HB3	1:C:79:MET:HG2	1.86	0.57
1:A:166:GLN:N	1:A:167:PRO:HD2	2.20	0.56
1:A:124:PRO:HB3	1:A:164:GLU:HG3	1.87	0.56
1:B:127:LEU:HD12	1:B:172:LEU:HD13	1.87	0.56
1:C:44:ASP:HB3	1:C:79:MET:CG	2.36	0.56
1:D:166:GLN:N	1:D:167:PRO:HD2	2.21	0.56
1:D:198:LEU:HD13	1:D:201:ILE:CG2	2.36	0.56
1:B:44:ASP:HB3	1:B:79:MET:CG	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:VAL:O	1:A:341:LEU:HG	2.06	0.56
1:C:166:GLN:N	1:C:167:PRO:HD2	2.21	0.56
1:B:123:VAL:HB	1:B:124:PRO:HD3	1.88	0.56
1:B:198:LEU:HD13	1:B:201:ILE:CG2	2.36	0.56
1:D:123:VAL:HB	1:D:124:PRO:HD3	1.88	0.55
1:D:203:ILE:HG12	1:D:205:THR:HG23	1.89	0.55
1:C:127:LEU:HD12	1:C:172:LEU:HD13	1.87	0.55
1:C:203:ILE:HG12	1:C:205:THR:HG23	1.88	0.55
1:A:127:LEU:HD12	1:A:172:LEU:HD13	1.89	0.55
1:D:141:TYR:CE2	1:D:179:SER:HB2	2.42	0.55
1:D:69:TRP:CE2	2:H:1:NAG:H82	2.42	0.54
1:D:127:LEU:HD12	1:D:172:LEU:HD13	1.88	0.54
1:D:310:ARG:NH2	1:D:330:ASP:OD2	2.40	0.54
1:A:141:TYR:CE2	1:A:179:SER:HB2	2.42	0.54
1:A:124:PRO:HB2	1:A:125:PRO:HD3	1.89	0.54
1:C:241:VAL:O	1:C:245:LEU:HG	2.08	0.54
1:B:54:ILE:HG13	1:B:90:LEU:HD11	1.89	0.54
1:B:124:PRO:HB2	1:B:125:PRO:HD3	1.90	0.54
1:D:124:PRO:HB3	1:D:164:GLU:HG3	1.88	0.54
1:A:44:ASP:HB3	1:A:79:MET:CG	2.37	0.54
1:B:105:ARG:O	1:B:109:ILE:HG13	2.07	0.54
1:B:141:TYR:CE2	1:B:179:SER:HB2	2.42	0.53
1:A:123:VAL:HB	1:A:124:PRO:HD3	1.89	0.53
1:C:198:LEU:HD13	1:C:201:ILE:HG22	1.90	0.53
1:A:66:ILE:O	1:A:66:ILE:HG23	2.08	0.53
1:B:124:PRO:HB3	1:B:164:GLU:HG3	1.89	0.53
1:C:105:ARG:O	1:C:109:ILE:HG13	2.09	0.53
1:A:69:TRP:CE2	2:E:1:NAG:H82	2.43	0.53
1:D:234:PHE:HB3	1:D:239:TYR:CE2	2.43	0.53
1:D:105:ARG:O	1:D:109:ILE:HG13	2.08	0.52
1:B:198:LEU:O	1:B:253:LYS:NZ	2.32	0.52
1:B:365:GLY:C	1:B:367:ASP:H	2.17	0.52
1:C:54:ILE:HG13	1:C:90:LEU:HD11	1.91	0.52
1:C:365:GLY:C	1:C:367:ASP:H	2.18	0.52
1:D:211:ALA:HA	1:D:265:PHE:CE1	2.45	0.52
1:D:198:LEU:HD13	1:D:201:ILE:HG22	1.92	0.52
1:A:105:ARG:O	1:A:109:ILE:HG13	2.09	0.52
1:D:365:GLY:C	1:D:367:ASP:H	2.17	0.52
1:A:365:GLY:C	1:A:367:ASP:H	2.17	0.52
1:A:198:LEU:HD13	1:A:201:ILE:HG22	1.91	0.52
1:D:209:HIS:CE1	1:D:213:ARG:HD3	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:LYS:O	1:A:170:LYS:HG3	2.11	0.51
1:D:66:ILE:HG23	1:D:66:ILE:O	2.10	0.51
1:D:54:ILE:HG13	1:D:90:LEU:HD11	1.93	0.51
1:B:80:LEU:HD12	1:B:83:LEU:HD12	1.92	0.51
1:B:198:LEU:HD13	1:B:201:ILE:HG22	1.92	0.51
1:B:66:ILE:HG23	1:B:66:ILE:O	2.10	0.51
1:D:210:GLY:HA3	1:D:212:TRP:CD1	2.46	0.50
1:B:154:ILE:HD13	1:B:198:LEU:HD11	1.94	0.50
1:B:191:ILE:HD11	1:B:244:MET:HE2	1.93	0.50
1:C:141:TYR:CE2	1:C:179:SER:HB2	2.46	0.50
1:C:230:SER:HB3	3:C:441:HOH:O	2.10	0.50
1:A:203:ILE:HD12	1:A:244:MET:SD	2.52	0.50
1:B:41:CYS:HB2	1:B:359:PHE:CZ	2.46	0.50
1:B:226:GLN:HE21	1:B:228:ASP:HB2	1.75	0.50
1:B:170:LYS:HG3	1:B:170:LYS:O	2.12	0.50
1:B:208:PHE:HB3	1:B:235:SER:HB3	1.94	0.50
1:C:170:LYS:O	1:C:170:LYS:HG3	2.11	0.50
1:B:33:GLN:HG3	1:B:34:TYR:CD1	2.47	0.50
1:C:124:PRO:HB3	1:C:164:GLU:HG3	1.92	0.50
1:D:203:ILE:HD11	1:D:240:ALA:HB1	1.94	0.49
1:A:263:ARG:NH1	1:A:293:THR:OG1	2.45	0.49
1:C:69:TRP:CE2	2:G:1:NAG:H82	2.47	0.49
1:A:54:ILE:HG13	1:A:90:LEU:HD11	1.95	0.49
1:C:80:LEU:HD12	1:C:83:LEU:HD12	1.94	0.49
1:C:86:ARG:NH1	3:C:460:HOH:O	2.45	0.48
1:A:154:ILE:HD13	1:A:198:LEU:HD11	1.95	0.48
1:A:208:PHE:HB3	1:A:235:SER:HB3	1.94	0.48
1:A:234:PHE:HB3	1:A:239:TYR:CE2	2.47	0.48
1:D:154:ILE:HD13	1:D:198:LEU:HD11	1.94	0.48
1:A:33:GLN:HG3	1:A:34:TYR:CD1	2.48	0.48
1:B:142:PRO:HB2	1:B:188:SER:HB3	1.96	0.48
1:C:154:ILE:HD13	1:C:198:LEU:HD11	1.95	0.48
1:C:209:HIS:CD2	1:C:217:GLY:HA3	2.49	0.48
1:C:263:ARG:NH1	1:C:293:THR:OG1	2.46	0.48
1:A:214:GLY:C	1:A:278:ILE:HG12	2.39	0.48
1:B:366:GLN:C	1:B:368:LEU:H	2.22	0.48
1:A:80:LEU:HD12	1:A:83:LEU:HD12	1.96	0.48
1:C:174:LEU:HG	1:C:198:LEU:HD23	1.96	0.48
1:A:190:ASP:OD1	1:A:193:LYS:HE3	2.14	0.48
1:C:33:GLN:HG3	1:C:34:TYR:CD1	2.48	0.48
1:C:203:ILE:HD11	1:C:240:ALA:HB1	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:GLY:O	1:A:213:ARG:HG3	2.14	0.47
1:B:191:ILE:CG2	1:B:249:ALA:HB2	2.44	0.47
1:C:66:ILE:HG23	1:C:66:ILE:O	2.13	0.47
1:A:241:VAL:O	1:A:245:LEU:HG	2.14	0.47
1:B:105:ARG:NH1	2:F:1:NAG:H62	2.29	0.47
1:C:366:GLN:C	1:C:368:LEU:H	2.22	0.47
1:D:84:LYS:NZ	1:D:133:ASP:OD1	2.47	0.47
1:D:144:ARG:NH1	1:D:144:ARG:HB3	2.29	0.47
1:D:170:LYS:HG3	1:D:170:LYS:O	2.13	0.47
1:C:366:GLN:C	1:C:368:LEU:N	2.72	0.47
1:D:142:PRO:HB2	1:D:188:SER:HB3	1.97	0.47
1:D:211:ALA:HB2	3:D:452:HOH:O	2.14	0.47
1:D:365:GLY:C	1:D:367:ASP:N	2.73	0.47
1:D:33:GLN:HG3	1:D:34:TYR:CD1	2.50	0.47
1:B:181:GLY:O	1:B:185:ILE:HG13	2.14	0.47
1:B:263:ARG:NH1	1:B:293:THR:OG1	2.48	0.47
1:C:144:ARG:HB3	1:C:144:ARG:NH1	2.30	0.47
1:D:80:LEU:HD12	1:D:83:LEU:HD12	1.96	0.47
1:D:211:ALA:HA	1:D:265:PHE:HE1	1.80	0.46
1:B:190:ASP:OD1	1:B:193:LYS:HE3	2.16	0.46
1:C:142:PRO:HB2	1:C:188:SER:HB3	1.97	0.46
1:B:174:LEU:HG	1:B:198:LEU:HD23	1.96	0.46
1:B:89:ASN:O	1:B:91:LYS:HD3	2.16	0.46
1:D:366:GLN:C	1:D:368:LEU:N	2.74	0.46
1:C:190:ASP:OD1	1:C:193:LYS:HE3	2.16	0.46
1:B:365:GLY:C	1:B:367:ASP:N	2.74	0.45
1:C:57:SER:HA	1:C:58:PHE:HA	1.65	0.45
1:A:174:LEU:HG	1:A:198:LEU:HD23	1.98	0.45
1:D:263:ARG:NH1	1:D:293:THR:OG1	2.49	0.45
1:A:365:GLY:C	1:A:367:ASP:N	2.74	0.45
1:A:366:GLN:C	1:A:368:LEU:H	2.24	0.45
1:C:181:GLY:O	1:C:185:ILE:HG13	2.16	0.45
1:A:366:GLN:C	1:A:368:LEU:N	2.74	0.45
1:A:383:THR:HG22	1:A:383:THR:O	2.17	0.45
1:D:57:SER:HA	1:D:58:PHE:HA	1.67	0.45
1:B:144:ARG:NH1	1:B:144:ARG:HB3	2.32	0.45
1:D:96:VAL:HG12	1:D:135:LEU:HD11	1.97	0.45
1:A:60:ASN:CG	1:A:61:ILE:H	2.25	0.45
1:B:366:GLN:C	1:B:368:LEU:N	2.72	0.45
1:C:214:GLY:C	1:C:278:ILE:HG12	2.42	0.45
1:C:365:GLY:C	1:C:367:ASP:N	2.74	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:366:GLN:C	1:D:368:LEU:H	2.23	0.45
1:C:232:ASP:HB3	1:C:234:PHE:H	1.80	0.45
1:A:181:GLY:O	1:A:185:ILE:HG13	2.17	0.45
1:B:383:THR:HG22	1:B:383:THR:O	2.17	0.44
1:C:96:VAL:HG12	1:C:135:LEU:HD11	1.99	0.44
1:D:190:ASP:OD1	1:D:193:LYS:HE3	2.16	0.44
1:D:174:LEU:HG	1:D:198:LEU:HD23	1.98	0.44
1:D:303:LEU:HA	1:D:306:ALA:HB3	1.99	0.44
1:B:61:ILE:HG21	1:B:109:ILE:HD13	1.99	0.44
1:C:81:ASN:OD1	1:C:92:THR:HG21	2.17	0.44
1:B:201:ILE:HG13	1:B:254:LEU:HD12	1.99	0.44
1:C:69:TRP:CD2	2:G:1:NAG:H82	2.53	0.44
1:C:201:ILE:HG13	1:C:254:LEU:HD12	1.99	0.44
1:C:366:GLN:O	1:C:368:LEU:N	2.51	0.44
1:D:34:TYR:HE2	1:D:70:GLU:OE2	2.01	0.44
1:A:96:VAL:HG12	1:A:135:LEU:HD11	1.98	0.44
1:B:366:GLN:O	1:B:368:LEU:N	2.51	0.44
1:C:60:ASN:CG	1:C:61:ILE:H	2.26	0.44
1:A:102:GLY:C	1:A:104:GLN:N	2.76	0.43
1:D:208:PHE:HB3	1:D:235:SER:HB3	2.00	0.43
1:A:160:GLU:OE1	1:A:163:LYS:HD3	2.18	0.43
1:C:102:GLY:C	1:C:104:GLN:N	2.76	0.43
1:C:383:THR:O	1:C:383:THR:HG22	2.17	0.43
1:A:57:SER:HA	1:A:58:PHE:HA	1.67	0.43
1:A:117:ARG:HB2	3:A:416:HOH:O	2.18	0.43
1:A:144:ARG:NH1	1:A:144:ARG:HB3	2.32	0.43
1:B:365:GLY:C	1:B:366:GLN:HG2	2.44	0.43
1:D:383:THR:HG22	1:D:383:THR:O	2.18	0.43
1:A:260:THR:O	1:A:296:TYR:HB2	2.19	0.43
1:B:96:VAL:HG12	1:B:135:LEU:HD11	1.99	0.43
1:A:89:ASN:O	1:A:91:LYS:HD3	2.19	0.43
1:B:60:ASN:CG	1:B:61:ILE:H	2.26	0.43
1:B:241:VAL:O	1:B:245:LEU:HG	2.18	0.43
1:D:102:GLY:C	1:D:104:GLN:N	2.76	0.43
1:D:211:ALA:O	1:D:278:ILE:HD11	2.19	0.43
1:C:160:GLU:OE1	1:C:163:LYS:HD3	2.18	0.43
1:D:330:ASP:HA	1:D:372:LEU:HD21	2.00	0.43
1:D:366:GLN:O	1:D:368:LEU:N	2.52	0.43
1:D:267:LEU:HD13	1:D:272:THR:HG22	2.01	0.43
1:A:330:ASP:HA	1:A:372:LEU:HD21	2.00	0.42
1:D:55:ILE:CG2	1:D:95:SER:HB2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:214:GLY:C	1:D:278:ILE:HG12	2.44	0.42
1:C:89:ASN:O	1:C:91:LYS:HD3	2.19	0.42
1:D:89:ASN:O	1:D:91:LYS:HD3	2.19	0.42
1:D:181:GLY:O	1:D:185:ILE:HG13	2.19	0.42
1:D:361:GLY:O	1:D:367:ASP:HA	2.20	0.42
1:D:160:GLU:OE1	1:D:163:LYS:HD3	2.20	0.42
1:D:209:HIS:ND1	1:D:213:ARG:HD3	2.34	0.42
1:A:34:TYR:HE2	1:A:70:GLU:OE2	2.02	0.42
1:D:60:ASN:CG	1:D:61:ILE:H	2.26	0.42
1:D:81:ASN:OD1	1:D:92:THR:HG21	2.20	0.42
1:C:52:THR:OG1	1:C:53:HIS:ND1	2.39	0.42
1:C:303:LEU:HA	1:C:306:ALA:HB3	2.02	0.42
1:D:48:ARG:CZ	1:D:86:ARG:HD2	2.50	0.42
1:B:203:ILE:HD11	1:B:240:ALA:HB1	2.01	0.42
1:D:48:ARG:NE	1:D:86:ARG:HB2	2.35	0.42
1:D:365:GLY:C	1:D:366:GLN:HG2	2.45	0.42
1:A:81:ASN:OD1	1:A:92:THR:HG21	2.20	0.42
1:A:230:SER:HA	1:A:231:PRO:HD3	1.84	0.42
1:B:330:ASP:HA	1:B:372:LEU:HD21	2.02	0.42
1:A:142:PRO:HB2	1:A:188:SER:HB3	2.02	0.41
1:B:57:SER:HA	1:B:58:PHE:HA	1.65	0.41
1:C:267:LEU:HD13	1:C:272:THR:HG22	2.02	0.41
1:B:297:TYR:HB2	1:B:363:PHE:CG	2.55	0.41
1:D:116:ARG:NH2	3:D:467:HOH:O	2.52	0.41
1:B:303:LEU:HA	1:B:306:ALA:HB3	2.00	0.41
1:D:297:TYR:HB2	1:D:363:PHE:CG	2.55	0.41
1:B:195:SER:O	1:B:253:LYS:HD3	2.21	0.41
1:B:214:GLY:C	1:B:278:ILE:HG12	2.45	0.41
1:B:263:ARG:HD3	1:B:263:ARG:HA	1.94	0.41
1:D:232:ASP:HB3	1:D:234:PHE:H	1.85	0.41
1:A:48:ARG:NE	1:A:86:ARG:HB2	2.35	0.41
1:B:230:SER:HA	1:B:231:PRO:HD3	1.85	0.41
1:A:33:GLN:HB2	1:A:42:PHE:CZ	2.55	0.41
1:B:160:GLU:OE1	1:B:163:LYS:HD3	2.20	0.41
1:B:102:GLY:C	1:B:104:GLN:N	2.76	0.41
1:C:30:SER:O	1:C:33:GLN:HG2	2.20	0.41
1:C:365:GLY:C	1:C:366:GLN:HG2	2.45	0.41
1:A:361:GLY:O	1:A:367:ASP:HA	2.21	0.41
1:C:61:ILE:HG21	1:C:109:ILE:HD13	2.02	0.41
1:C:86:ARG:NH1	1:C:86:ARG:HG2	2.36	0.41
1:C:297:TYR:HB2	1:C:363:PHE:CG	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:330:ASP:HA	1:C:372:LEU:HD21	2.03	0.41
1:D:86:ARG:NH1	1:D:86:ARG:HG2	2.36	0.41
1:D:140:LEU:HA	1:D:141:TYR:HA	1.80	0.41
1:A:267:LEU:HD13	1:A:272:THR:HG22	2.03	0.41
1:C:331:GLN:NE2	3:C:454:HOH:O	2.46	0.41
1:C:341:LEU:CB	1:C:346:LEU:HD12	2.51	0.41
1:D:34:TYR:CE2	1:D:70:GLU:OE2	2.74	0.41
1:A:366:GLN:O	1:A:368:LEU:N	2.54	0.40
1:B:30:SER:O	1:B:33:GLN:HG2	2.21	0.40
1:B:34:TYR:HE2	1:B:70:GLU:OE2	2.03	0.40
1:D:41:CYS:HB2	1:D:359:PHE:CZ	2.57	0.40
1:B:55:ILE:CG2	1:B:95:SER:HB2	2.51	0.40
1:C:263:ARG:HA	1:C:263:ARG:HD3	1.93	0.40
1:C:48:ARG:CZ	1:C:86:ARG:HD2	2.52	0.40
1:C:223:PHE:HB2	1:C:313:GLY:O	2.21	0.40
1:D:139:TRP:O	1:D:141:TYR:HA	2.21	0.40
1:B:52:THR:OG1	1:B:53:HIS:ND1	2.38	0.40
1:D:62:SER:O	1:D:63:ASN:C	2.63	0.40
1:D:309:HIS:HB2	1:D:318:TYR:CZ	2.57	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	360/362 (99%)	339 (94%)	17 (5%)	4 (1%)	11	29
1	B	360/362 (99%)	338 (94%)	19 (5%)	3 (1%)	16	37
1	C	360/362 (99%)	339 (94%)	18 (5%)	3 (1%)	16	37
1	D	360/362 (99%)	341 (95%)	16 (4%)	3 (1%)	16	37
All	All	1440/1448 (99%)	1357 (94%)	70 (5%)	13 (1%)	14	35

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	231	PRO
1	D	231	PRO
1	A	231	PRO
1	B	367	ASP
1	C	367	ASP
1	D	367	ASP
1	A	367	ASP
1	B	231	PRO
1	C	103	SER
1	D	103	SER
1	A	103	SER
1	B	103	SER
1	A	191	ILE

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	301/302 (100%)	295 (98%)	6 (2%)	48	76
1	B	301/302 (100%)	295 (98%)	6 (2%)	48	76
1	C	301/302 (100%)	295 (98%)	6 (2%)	48	76
1	D	301/302 (100%)	296 (98%)	5 (2%)	53	79
All	All	1204/1208 (100%)	1181 (98%)	23 (2%)	50	77

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	144	ARG
1	A	193	LYS
1	A	227	GLU
1	A	234	PHE
1	A	244	MET
1	A	331	GLN
1	B	144	ARG

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Mol	Chain	Res	Type
1	B	193	LYS
1	B	227	GLU
1	B	234	PHE
1	B	331	GLN
1	B	366	GLN
1	C	144	ARG
1	C	193	LYS
1	C	227	GLU
1	C	234	PHE
1	C	331	GLN
1	C	366	GLN
1	D	144	ARG
1	D	193	LYS
1	D	234	PHE
1	D	331	GLN
1	D	366	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	ASN
1	A	148	GLN
1	A	226	GLN
1	A	331	GLN
1	A	345	GLN
1	A	366	GLN
1	B	100	ASN
1	B	148	GLN
1	B	226	GLN
1	B	331	GLN
1	B	345	GLN
1	B	366	GLN
1	C	100	ASN
1	C	148	GLN
1	C	226	GLN
1	C	331	GLN
1	C	345	GLN
1	C	366	GLN
1	D	100	ASN
1	D	148	GLN
1	D	226	GLN
1	D	345	GLN

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Mol	Chain	Res	Type
1	D	366	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 ⓘ

8 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	0.70	0	17,19,21	0.86	1 (5%)
2	NAG	E	2	2	14,14,15	0.80	0	17,19,21	0.89	1 (5%)
2	NAG	F	1	1,2	14,14,15	0.58	0	17,19,21	0.66	0
2	NAG	F	2	2	14,14,15	0.70	0	17,19,21	0.76	0
2	NAG	G	1	1,2	14,14,15	0.55	0	17,19,21	0.73	1 (5%)
2	NAG	G	2	2	14,14,15	0.58	0	17,19,21	0.66	0
2	NAG	H	1	1,2	14,14,15	0.71	0	17,19,21	0.70	0
2	NAG	H	2	2	14,14,15	0.79	0	17,19,21	0.67	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	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	G	2	2	-	3/6/23/26	0/1/1/1
2	NAG	H	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	H	2	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	NAG	C2-N2-C7	-2.88	119.05	122.90
2	E	2	NAG	C2-N2-C7	-2.38	119.72	122.90
2	G	1	NAG	C2-N2-C7	-2.32	119.80	122.90

There are no chirality outliers.

All (17) torsion outliers are listed below:

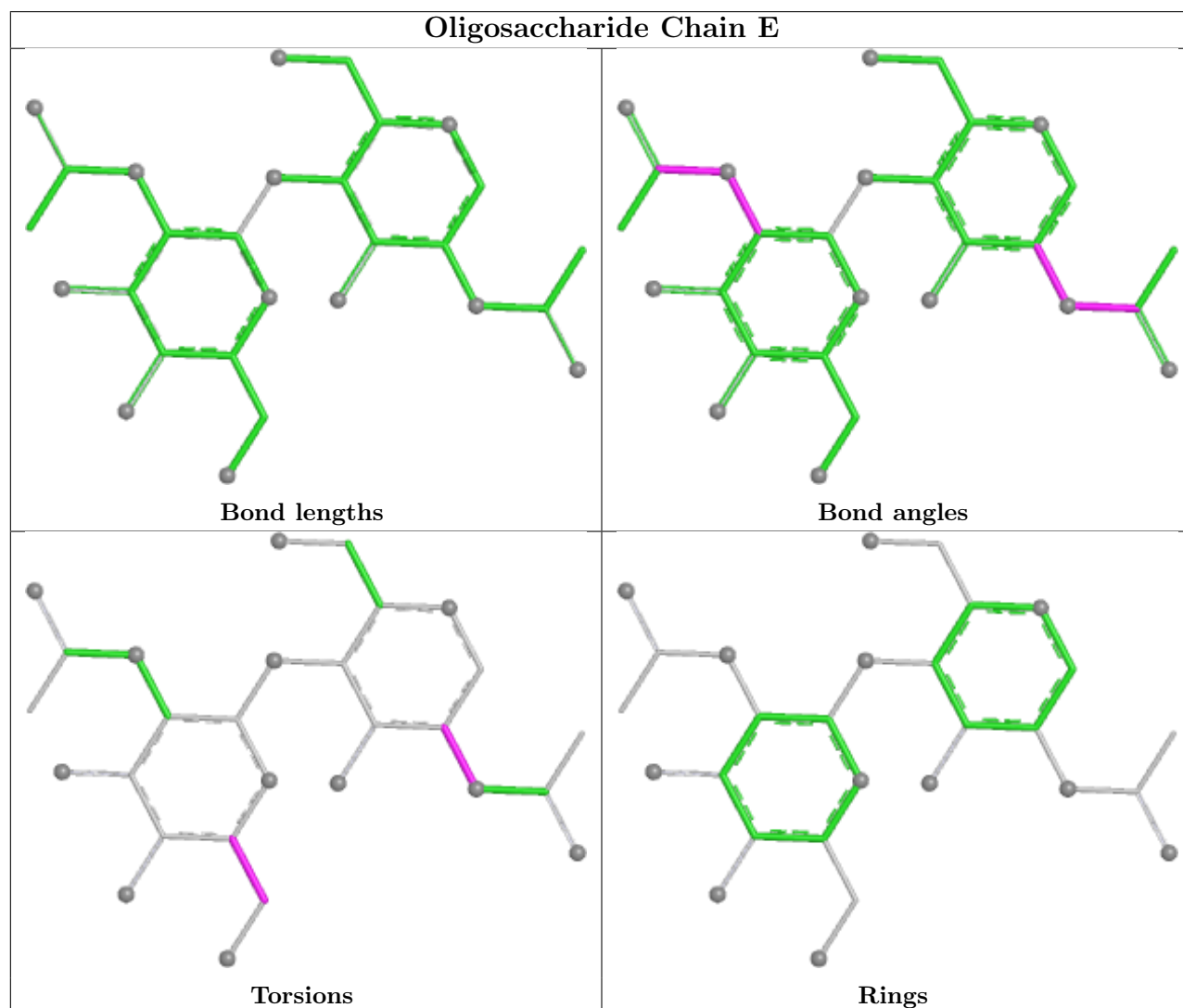
Mol	Chain	Res	Type	Atoms
2	E	2	NAG	O5-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
2	H	1	NAG	O5-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	H	1	NAG	C4-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
2	G	1	NAG	C4-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
2	E	1	NAG	C1-C2-N2-C7
2	G	2	NAG	C1-C2-N2-C7
2	H	1	NAG	O7-C7-N2-C2

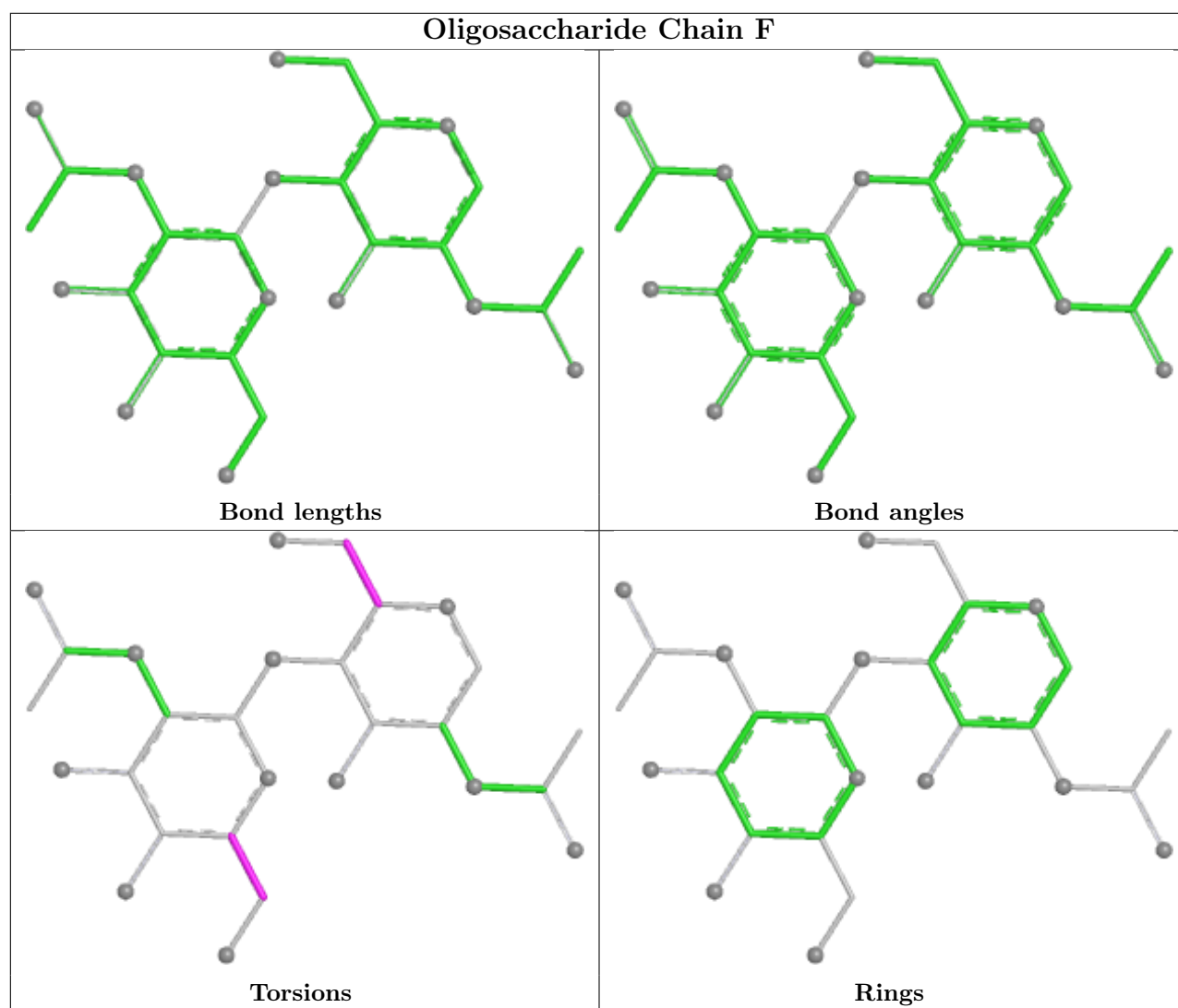
There are no ring outliers.

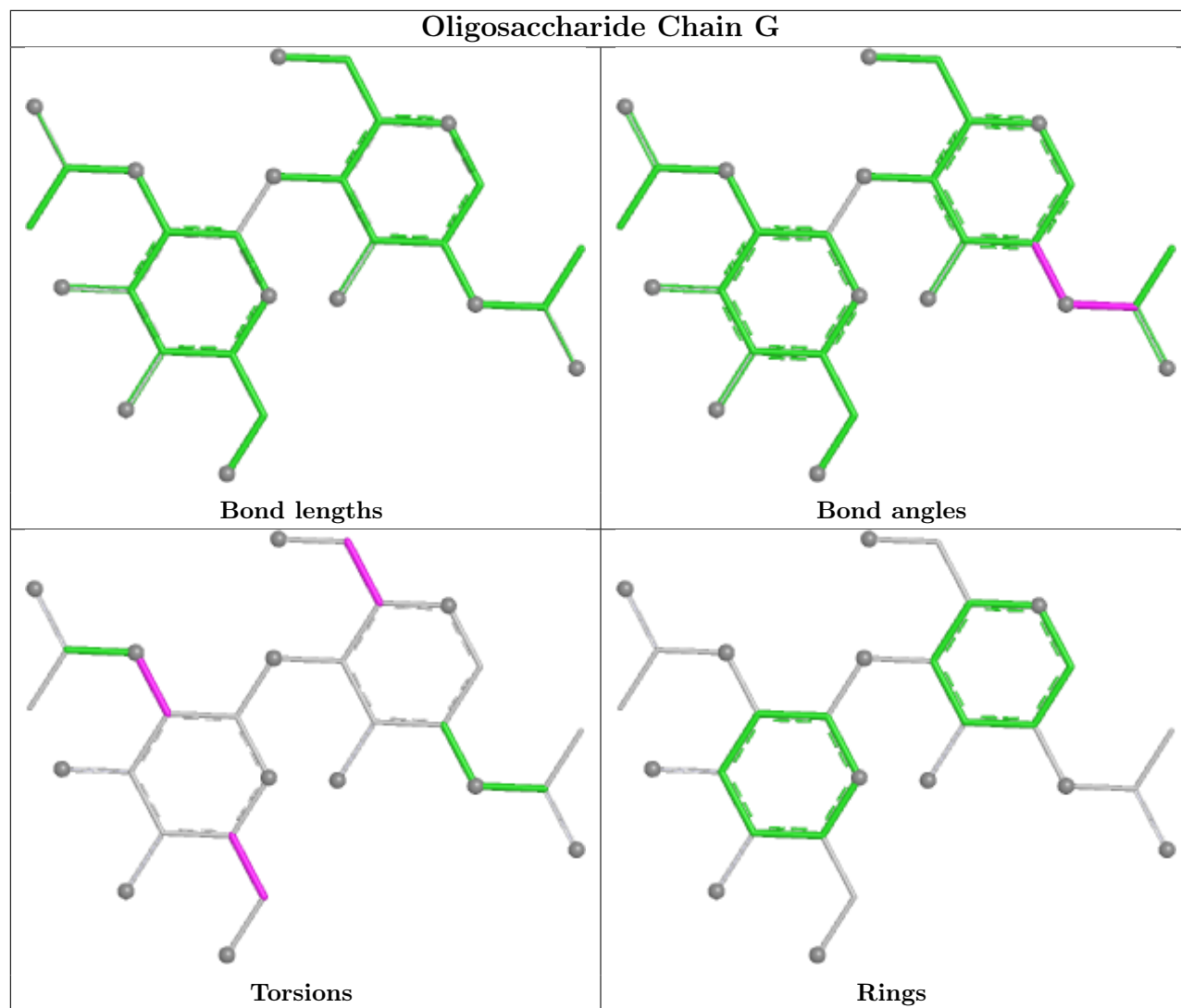
6 monomers are involved in 10 short contacts:

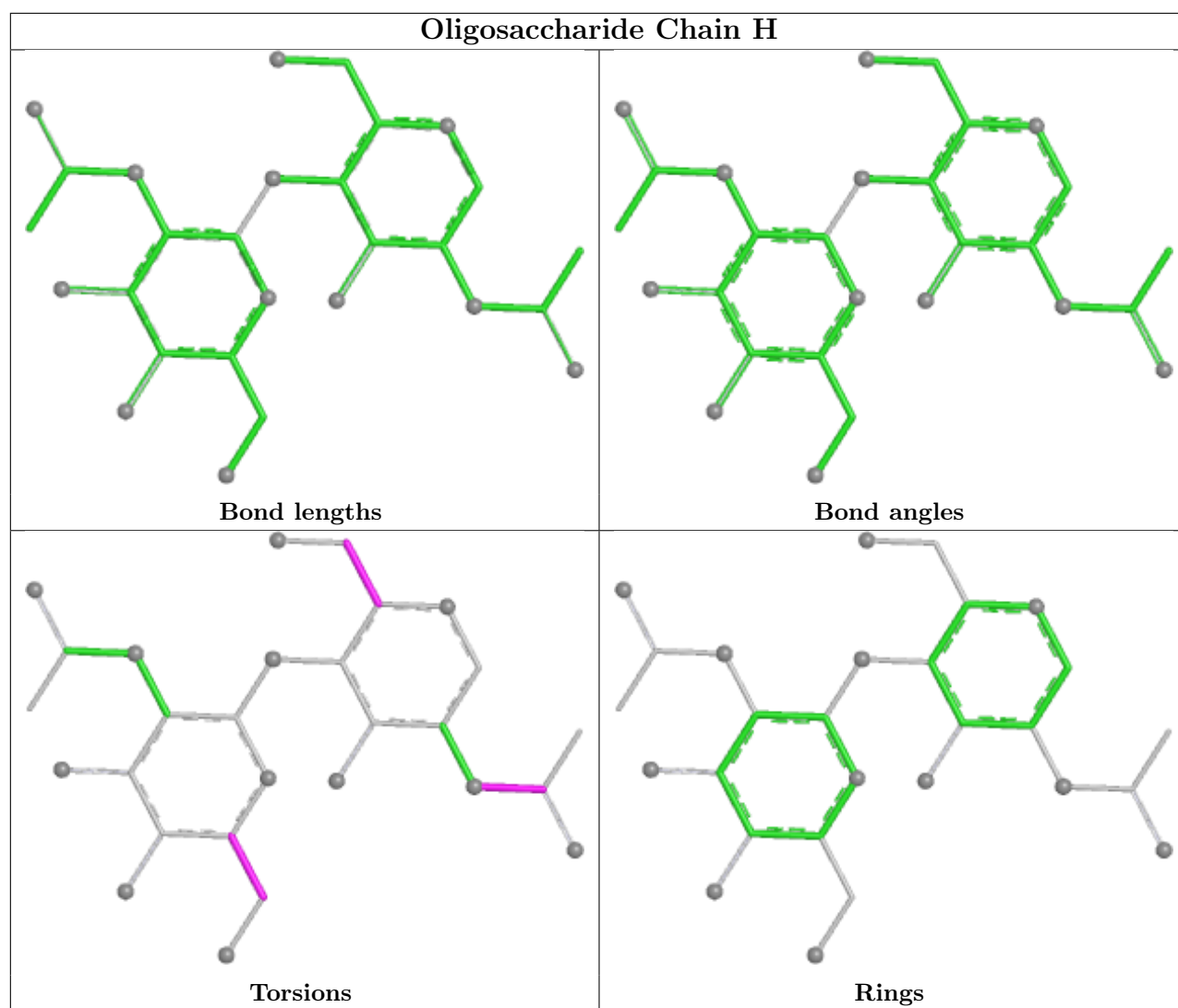
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1	NAG	1	0
2	E	2	NAG	4	0
2	H	2	NAG	1	0
2	H	1	NAG	2	0
2	G	1	NAG	2	0
2	E	1	NAG	5	0

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

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

6.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

6.4 Ligands

EDS failed to run properly - this section is therefore empty.

6.5 Other polymers

EDS failed to run properly - this section is therefore empty.