



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 17, 2024 – 11:09 AM EST

PDB ID : 4RLD
Title : Crystal structure of kkf mutant of bla G 2 protein
Authors : Li, M.; Gustchina, A.; Pomes, A.; Wlodawer, A.
Deposited on : 2014-10-16
Resolution : 2.90 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

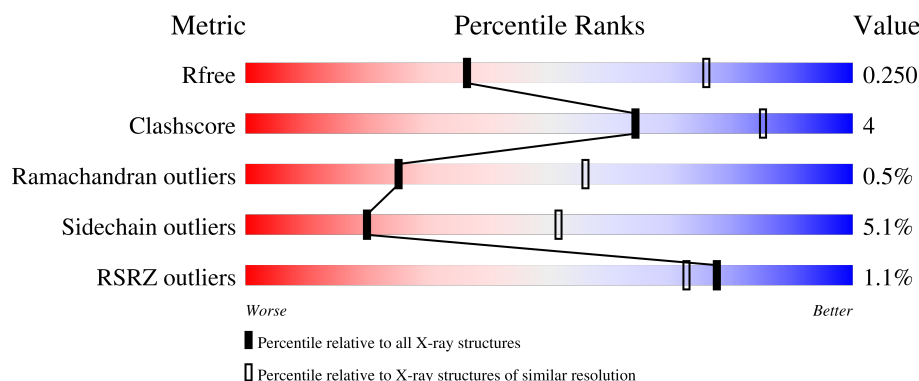
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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.



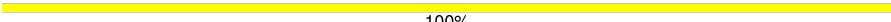
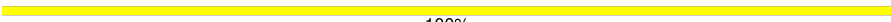

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	2335 (2.90-2.90)
Clashscore	180529	2564 (2.90-2.90)
Ramachandran outliers	177936	2514 (2.90-2.90)
Sidechain outliers	177891	2516 (2.90-2.90)
RSRZ outliers	164620	2337 (2.90-2.90)

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	330	
1	B	330	
1	C	330	
1	D	330	
2	E	2	

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

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10320 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 Aspartic protease Bla g 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	3	0
			2542	1620	418	492	12			
1	B	330	Total	C	N	O	S	0	3	0
			2542	1620	418	492	12			
1	C	330	Total	C	N	O	S	0	3	0
			2542	1620	418	492	12			
1	D	330	Total	C	N	O	S	0	3	0
			2542	1620	418	492	12			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	GLY	-	expression tag	UNP P54958
A	-7	ALA	-	expression tag	UNP P54958
A	-6	SER	-	expression tag	UNP P54958
A	-5	ILE	-	expression tag	UNP P54958
A	93	GLN	ASN	conflict	UNP P54958
A	132	ALA	LYS	engineered mutation	UNP P54958
A	162	TYR	PHE	engineered mutation	UNP P54958
A	251	ALA	LYS	engineered mutation	UNP P54958
B	-8	GLY	-	expression tag	UNP P54958
B	-7	ALA	-	expression tag	UNP P54958
B	-6	SER	-	expression tag	UNP P54958
B	-5	ILE	-	expression tag	UNP P54958
B	93	GLN	ASN	engineered mutation	UNP P54958
B	132	ALA	LYS	engineered mutation	UNP P54958
B	162	TYR	PHE	engineered mutation	UNP P54958
B	251	ALA	LYS	engineered mutation	UNP P54958
C	-8	GLY	-	expression tag	UNP P54958
C	-7	ALA	-	expression tag	UNP P54958
C	-6	SER	-	expression tag	UNP P54958
C	-5	ILE	-	expression tag	UNP P54958
C	93	GLN	ASN	conflict	UNP P54958

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Chain	Residue	Modelled	Actual	Comment	Reference
C	132	ALA	LYS	engineered mutation	UNP P54958
C	162	TYR	PHE	engineered mutation	UNP P54958
C	251	ALA	LYS	engineered mutation	UNP P54958
D	-8	GLY	-	expression tag	UNP P54958
D	-7	ALA	-	expression tag	UNP P54958
D	-6	SER	-	expression tag	UNP P54958
D	-5	ILE	-	expression tag	UNP P54958
D	93	GLN	ASN	conflict	UNP P54958
D	132	ALA	LYS	engineered mutation	UNP P54958
D	162	TYR	PHE	engineered mutation	UNP P54958
D	251	ALA	LYS	engineered mutation	UNP P54958

- 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			
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		

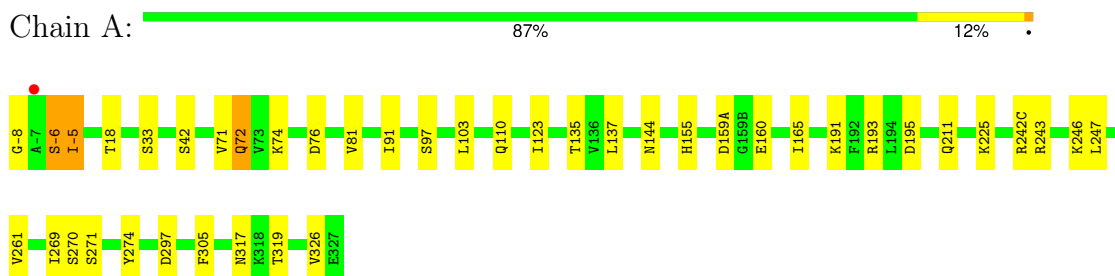
- Molecule 4 is water.

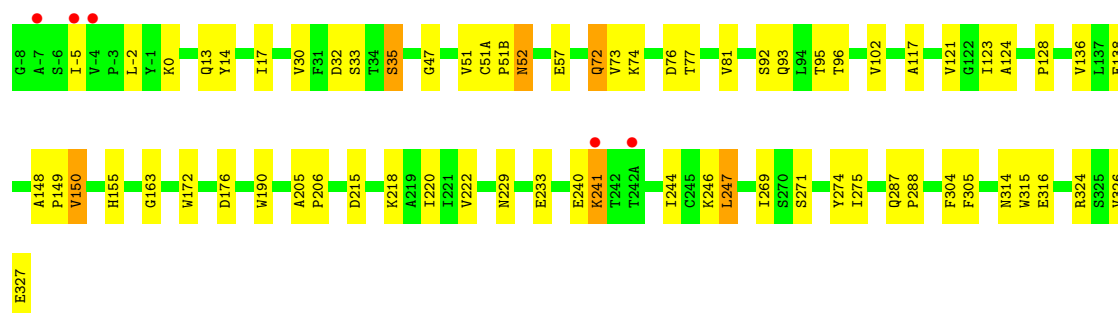
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total 12	O 12	0	0
4	B	13	Total 13	O 13	0	0
4	C	7	Total 7	O 7	0	0
4	D	4	Total 4	O 4	0	0

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: Aspartic protease Bla g 2





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

Chain E:  100%

NAG1
NAG2

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

Chain F:  100%

NAG1
NAG2

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

Chain G:  100%

NAG1
NAG2

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

Chain H:  50%  50%

NAG1
NAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.82Å 75.41Å 339.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.90 50.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	88.0 (50.00-2.90) 88.0 (50.00-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.196 , 0.257 0.191 , 0.250	Depositor DCC
R_{free} test set	1074 reflections (3.11%)	wwPDB-VP
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10320	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.48	1/2618 (0.0%)	0.58	0/3559
1	B	0.46	0/2618	0.58	0/3559
1	C	0.46	1/2618 (0.0%)	0.57	0/3559
1	D	0.52	2/2618 (0.1%)	0.58	0/3559
All	All	0.48	4/10472 (0.0%)	0.58	0/14236

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	138	GLU	CB-CG	6.40	1.64	1.52
1	D	172	TRP	CD2-CE3	6.20	1.49	1.40
1	C	-6	SER	CB-OG	5.12	1.49	1.42
1	A	-8	GLY	N-CA	5.10	1.53	1.46

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	2542	0	2454	13	0
1	B	2542	0	2454	20	0
1	C	2542	0	2454	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2542	0	2454	36	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	12	0	0	1	0
4	B	13	0	0	2	0
4	C	7	0	0	0	0
4	D	4	0	0	0	0
All	All	10320	0	9916	76	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 4.

All (76) 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:150:VAL:HG12	1:D:314:ASN:HA	1.53	0.89
1:B:29:THR:HG21	1:B:121:VAL:HG23	1.62	0.81
1:D:229:ASN:O	1:D:233:GLU:HG2	1.88	0.73
1:D:149:PRO:HG2	1:D:316:GLU:HB2	1.71	0.72
1:D:72:GLN:HG2	1:D:81:VAL:HG22	1.73	0.71
1:B:29:THR:HG23	1:B:119:VAL:HG12	1.73	0.70
1:A:-5:ILE:HD13	1:A:144:ASN:HD21	1.58	0.68
1:A:269:ILE:HD11	1:A:305:PHE:HD1	1.61	0.64
1:B:29:THR:HG21	1:B:121:VAL:CG2	2.27	0.62
1:B:242:THR:HG23	4:B:511:HOH:O	2.00	0.61
1:D:326:VAL:O	1:D:327:GLU:HB2	1.99	0.61
1:D:269:ILE:HD11	1:D:305:PHE:HD1	1.64	0.61
1:A:317:ASN:O	1:A:319:THR:HG23	2.01	0.60
1:B:269:ILE:HD11	1:B:305:PHE:HD1	1.67	0.59
1:A:195:ASP:HB2	1:A:261:VAL:HG12	1.86	0.58
1:B:191:LYS:HD2	1:B:211:GLN:HE21	1.69	0.57
1:A:193:ARG:HD2	4:A:601:HOH:O	2.04	0.56
1:D:51:VAL:O	1:D:51(B):PRO:HD3	2.05	0.56
1:D:215:ASP:OD2	1:D:218:LYS:HE2	2.07	0.55
1:C:222:VAL:HG22	1:C:287:GLN:HB3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:4:VAL:HB	1:C:165:ILE:HG23	1.89	0.55
1:B:233:GLU:HA	1:B:233:GLU:OE1	2.07	0.54
1:C:33:SER:HA	1:C:123:ILE:HG13	1.90	0.53
1:D:74:LYS:HE2	1:D:76:ASP:O	2.09	0.53
1:D:73:VAL:HG11	1:D:102:VAL:HG21	1.90	0.53
1:D:121:VAL:HG13	1:D:136:VAL:HG21	1.91	0.53
1:B:29:THR:CG2	1:B:121:VAL:HG23	2.36	0.52
1:C:72:GLN:HG2	1:C:81:VAL:HG22	1.92	0.52
1:A:159(A):ASP:OD2	1:A:160:GLU:OE1	2.29	0.52
1:D:33:SER:HA	1:D:123:ILE:HG13	1.94	0.50
1:D:13:GLN:O	1:D:14:TYR:HB2	2.13	0.49
1:D:32:ASP:OD2	1:D:35:SER:HB3	2.13	0.48
1:D:72:GLN:CG	1:D:81:VAL:HG22	2.42	0.48
1:D:92[B]:SER:HB3	1:D:93:GLN:H	1.52	0.48
1:D:220:ILE:HA	1:D:304:PHE:CE1	2.49	0.47
1:C:29:THR:HG23	1:C:119:VAL:HG12	1.96	0.47
1:B:-5:ILE:HD12	1:B:144:ASN:HD21	1.79	0.47
1:A:271:SER:HA	1:A:274:TYR:CE2	2.50	0.46
1:B:237:CYS:HB3	1:B:247:LEU:HB3	1.98	0.46
1:C:18:THR:HG22	1:C:91:ILE:HG12	1.97	0.46
1:B:146:ILE:HD13	1:B:167:PHE:HB3	1.98	0.45
1:A:42:SER:HB3	1:A:103:LEU:HG	1.99	0.45
1:B:15:ALA:HA	1:B:29:THR:O	2.17	0.45
1:D:176:ASP:HB3	1:D:324:ARG:HB2	1.98	0.45
1:A:191:LYS:HD2	1:A:211:GLN:HE21	1.82	0.44
1:B:313:PHE:CD1	1:B:313:PHE:N	2.86	0.44
1:B:221:ILE:HG12	1:B:301:ILE:HA	1.99	0.44
1:C:53:LEU:HD12	1:C:118:ASP:HA	2.00	0.44
1:D:47:GLY:HA2	1:D:51:VAL:CG2	2.48	0.44
1:D:271:SER:HA	1:D:274:TYR:CE2	2.53	0.44
1:B:92[A]:SER:OG	1:B:93:GLN:NE2	2.51	0.43
1:D:155:HIS:HA	1:D:163:GLY:HA3	2.01	0.43
1:D:148:ALA:O	1:D:150:VAL:N	2.45	0.42
1:B:214:ILE:HD13	1:B:301:ILE:HB	2.01	0.42
1:D:218:LYS:HD3	1:D:222:VAL:HG23	2.00	0.42
1:B:243:ARG:HA	4:B:511:HOH:O	2.19	0.42
1:D:51(A):CYS:HA	1:D:51(B):PRO:HD2	1.92	0.42
1:D:30:VAL:HG23	1:D:117:ALA:CB	2.49	0.42
1:A:165:ILE:O	1:A:165:ILE:HG23	2.18	0.42
1:C:25:GLN:HB3	1:C:27:PHE:CE1	2.54	0.42
1:D:52:ASN:HD22	1:D:52:ASN:HA	1.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:269:ILE:HD11	1:D:305:PHE:CD1	2.51	0.41
1:A:33:SER:HA	1:A:123:ILE:HG13	2.02	0.41
1:B:110:GLN:O	1:B:114:ILE:HG12	2.20	0.41
1:A:72:GLN:HG2	1:A:81:VAL:HG22	2.03	0.41
1:D:149:PRO:O	1:D:315:TRP:HB3	2.21	0.41
1:D:287:GLN:HA	1:D:288:PRO:HD3	1.92	0.41
1:D:124:ALA:HA	1:D:190:TRP:CD1	2.57	0.40
1:D:205:ALA:HA	1:D:206:PRO:HD3	1.90	0.40
1:D:247:LEU:HD21	1:D:275:ILE:HD12	2.03	0.40
1:A:18:THR:HG22	1:A:91:ILE:HG12	2.02	0.40
1:B:238:VAL:HG22	1:D:93:GLN:HG2	2.04	0.40
1:D:47:GLY:HA2	1:D:51:VAL:HG22	2.03	0.40
1:B:195:ASP:HB2	1:B:261:VAL:HG12	2.04	0.40
1:D:176:ASP:HB2	1:D:326:VAL:HG23	2.03	0.40
1:D:241:LYS:HG3	1:D:241:LYS:O	2.21	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	331/330 (100%)	310 (94%)	20 (6%)	1 (0%)	37	66
1	B	331/330 (100%)	315 (95%)	14 (4%)	2 (1%)	22	52
1	C	331/330 (100%)	312 (94%)	16 (5%)	3 (1%)	14	43
1	D	331/330 (100%)	312 (94%)	18 (5%)	1 (0%)	37	66
All	All	1324/1320 (100%)	1249 (94%)	68 (5%)	7 (0%)	25	56

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	-6	SER
1	C	76	ASP
1	C	242(B)	THR
1	C	225	LYS
1	B	76	ASP
1	B	126	PRO
1	D	128	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	283/280 (101%)	264 (93%)	19 (7%)	13	39
1	B	283/280 (101%)	275 (97%)	8 (3%)	38	73
1	C	283/280 (101%)	270 (95%)	13 (5%)	23	55
1	D	283/280 (101%)	266 (94%)	17 (6%)	16	44
All	All	1132/1120 (101%)	1075 (95%)	57 (5%)	20	52

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-6	SER
1	A	-5	ILE
1	A	71	VAL
1	A	72	GLN
1	A	74	LYS
1	A	76	ASP
1	A	97	SER
1	A	110	GLN
1	A	135	THR
1	A	137	LEU
1	A	155	HIS
1	A	225	LYS
1	A	242(C)	ARG
1	A	243	ARG
1	A	246	LYS

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Mol	Chain	Res	Type
1	A	247	LEU
1	A	270	SER
1	A	297	ASP
1	A	326	VAL
1	B	-5	ILE
1	B	96	THR
1	B	179	PHE
1	B	242(C)	ARG
1	B	243	ARG
1	B	247	LEU
1	B	270	SER
1	B	287	GLN
1	C	51	VAL
1	C	57	GLU
1	C	134	LYS
1	C	137	LEU
1	C	155	HIS
1	C	244	ILE
1	C	246	LYS
1	C	248	ASP
1	C	269	ILE
1	C	270	SER
1	C	292	SER
1	C	297	ASP
1	C	316	GLU
1	D	-5	ILE
1	D	-2	LEU
1	D	0	LYS
1	D	17	ILE
1	D	35	SER
1	D	52	ASN
1	D	57	GLU
1	D	72	GLN
1	D	77	THR
1	D	95	THR
1	D	96	THR
1	D	150	VAL
1	D	240	GLU
1	D	241	LYS
1	D	244	ILE
1	D	246	LYS
1	D	247	LEU

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

Mol	Chain	Res	Type
1	A	26	ASN
1	A	52	ASN
1	A	98	GLN
1	A	144	ASN
1	A	211	GLN
1	B	26	ASN
1	B	52	ASN
1	B	93	GLN
1	B	98	GLN
1	B	211	GLN
1	B	287	GLN
1	C	26	ASN
1	C	52	ASN
1	C	93	GLN
1	C	98	GLN
1	C	159	GLN
1	C	211	GLN
1	C	272	GLN
1	C	278	ASN
1	D	52	ASN
1	D	110	GLN
1	D	159	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	0.83	1 (7%)	17,19,21	1.03	1 (5%)
2	NAG	E	2	2	14,14,15	0.47	0	17,19,21	1.05	1 (5%)
2	NAG	F	1	1,2	14,14,15	0.59	0	17,19,21	1.48	1 (5%)
2	NAG	F	2	2	14,14,15	0.71	0	17,19,21	1.28	2 (11%)
2	NAG	G	1	1,2	14,14,15	0.55	0	17,19,21	1.34	2 (11%)
2	NAG	G	2	2	14,14,15	0.62	0	17,19,21	1.02	1 (5%)
2	NAG	H	1	1,2	14,14,15	0.60	0	17,19,21	1.20	1 (5%)
2	NAG	H	2	2	14,14,15	0.58	0	17,19,21	1.05	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	NAG	F	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
2	NAG	H	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1	NAG	C1-C2	2.07	1.55	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	NAG	C1-O5-C5	4.97	118.84	112.19
2	G	1	NAG	C1-O5-C5	3.96	117.49	112.19
2	H	1	NAG	C1-O5-C5	3.67	117.11	112.19
2	F	2	NAG	C4-C3-C2	3.58	116.27	111.02
2	G	2	NAG	C4-C3-C2	2.95	115.34	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	1	NAG	C4-C3-C2	2.76	115.07	111.02
2	E	1	NAG	C4-C3-C2	2.40	114.53	111.02
2	E	2	NAG	C1-O5-C5	2.09	114.99	112.19
2	F	2	NAG	C1-C2-N2	-2.05	107.19	110.43

There are no chirality outliers.

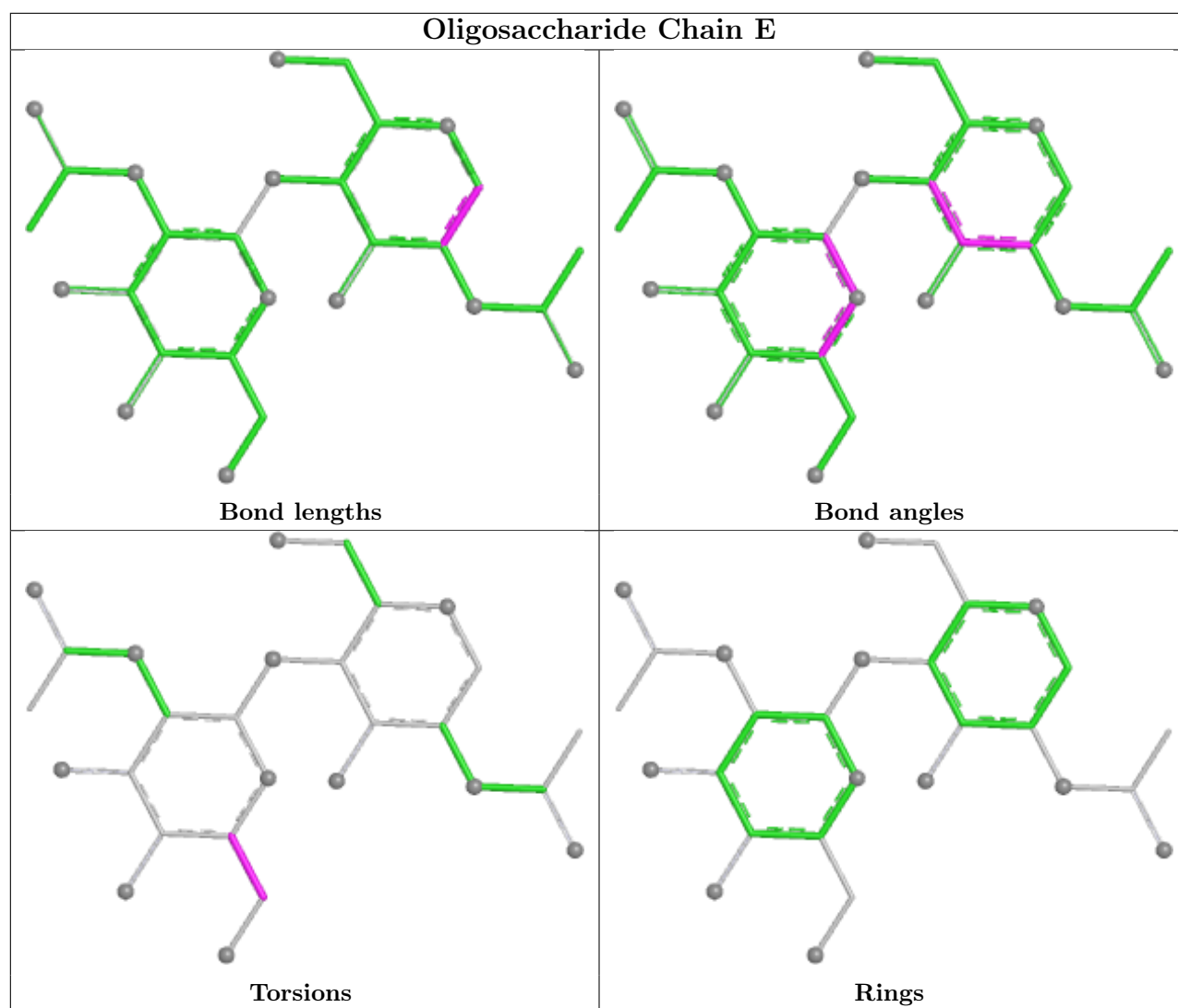
All (8) torsion outliers are listed below:

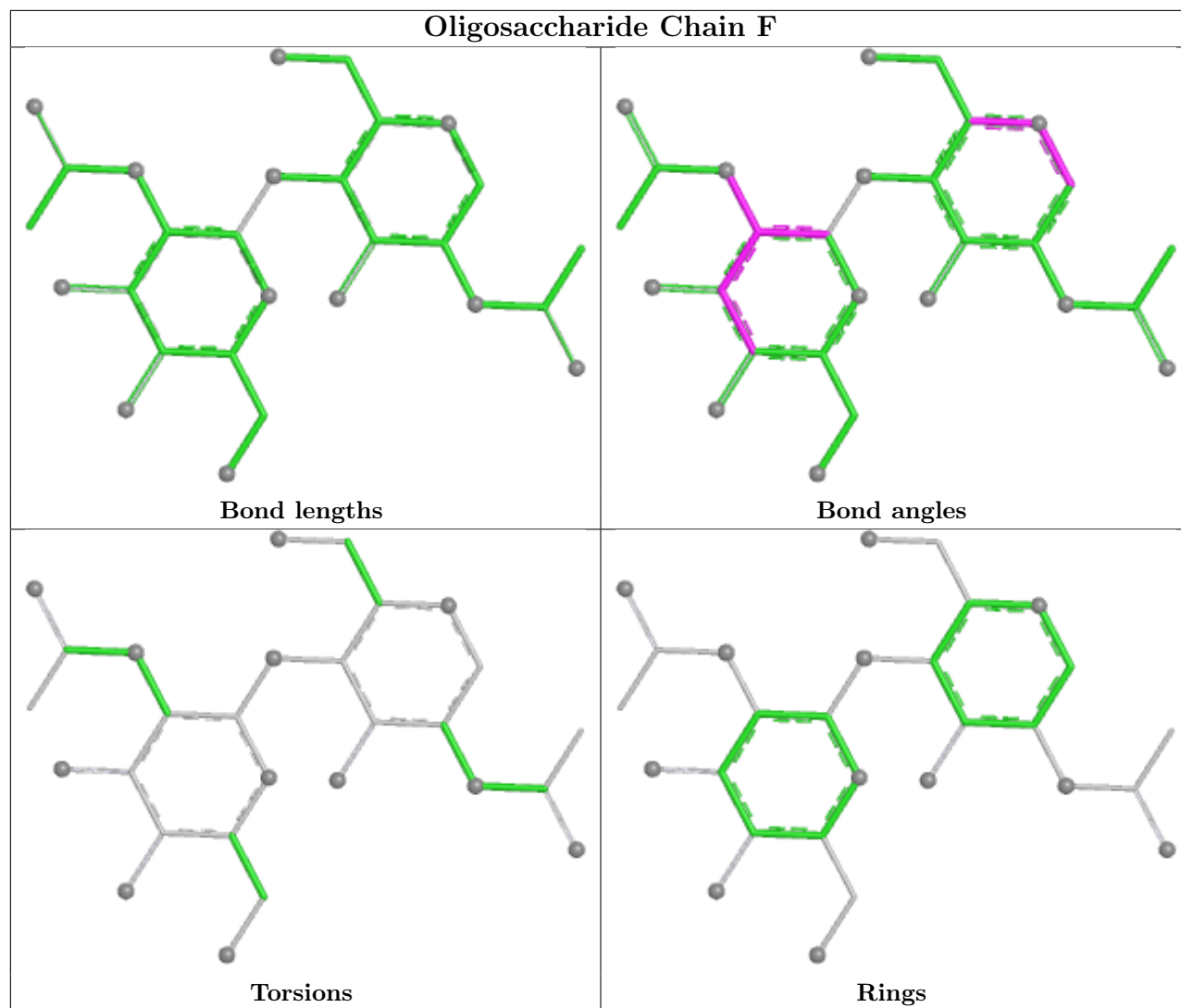
Mol	Chain	Res	Type	Atoms
2	G	2	NAG	O5-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	G	1	NAG	C4-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6

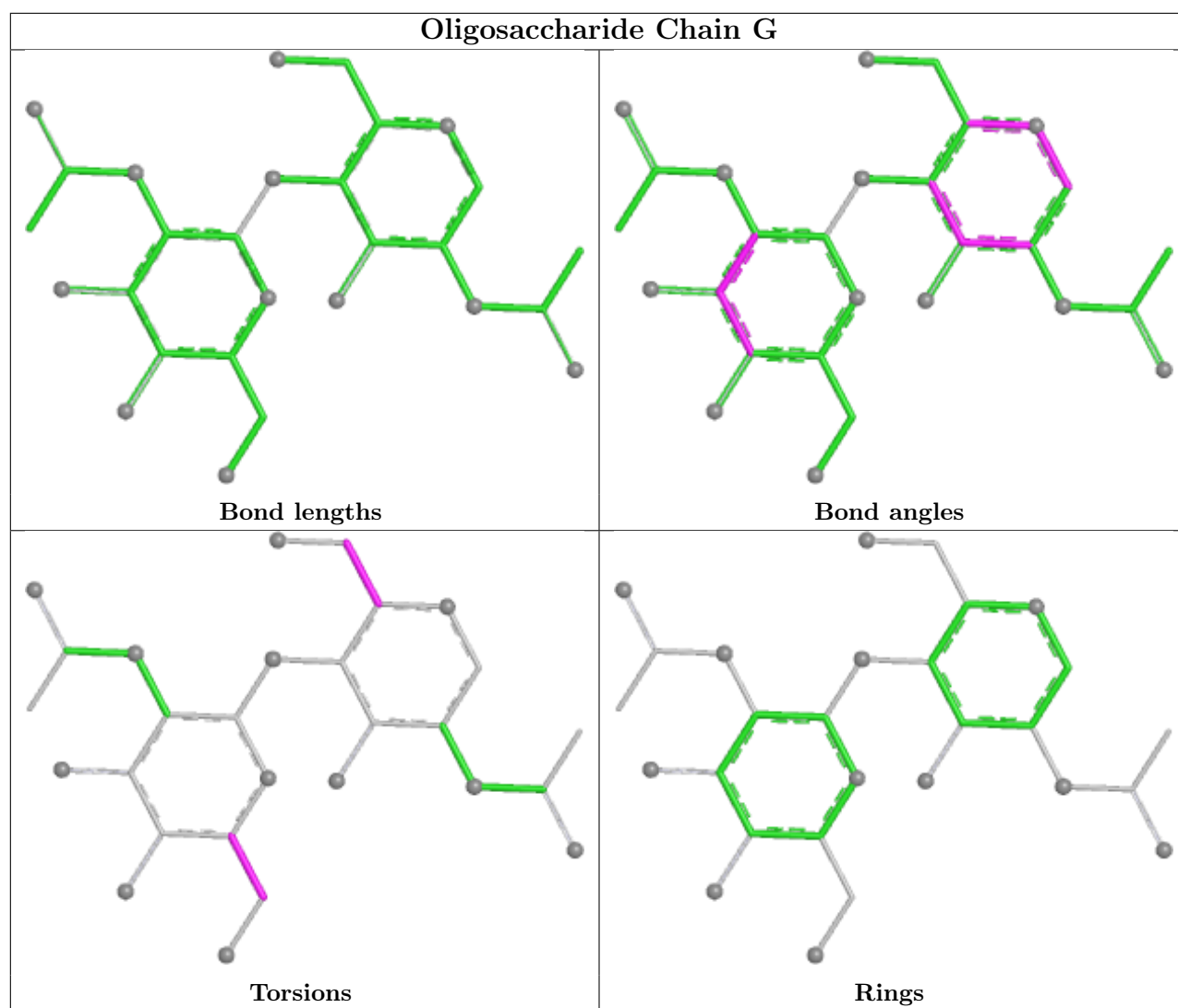
There are no ring outliers.

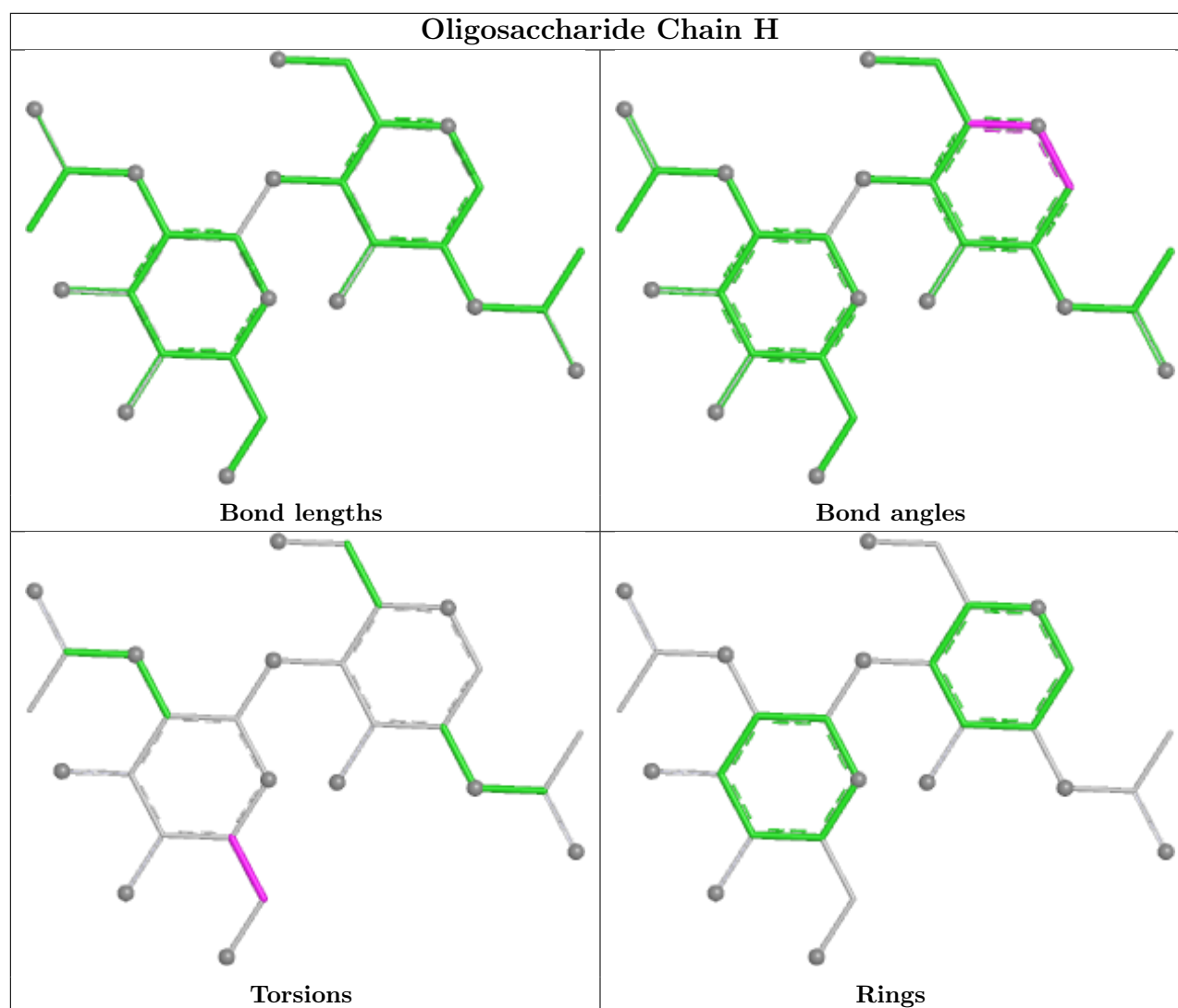
No monomer is involved in short contacts.

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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	330/330 (100%)	-0.61	1 (0%) 90 88	21, 39, 60, 109	2 (0%)
1	B	330/330 (100%)	-0.57	3 (0%) 81 76	24, 42, 63, 112	2 (0%)
1	C	330/330 (100%)	-0.29	6 (1%) 67 61	32, 51, 75, 130	2 (0%)
1	D	330/330 (100%)	-0.05	5 (1%) 71 66	31, 63, 94, 185	2 (0%)
All	All	1320/1320 (100%)	-0.38	15 (1%) 77 72	21, 47, 82, 185	8 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	-5	ILE	4.0
1	A	-7	ALA	3.8
1	C	-5	ILE	3.7
1	D	-7	ALA	3.4
1	D	242(A)	THR	3.3
1	C	-7	ALA	3.1
1	B	-7	ALA	3.0
1	D	241	LYS	2.9
1	C	76	ASP	2.8
1	B	-5	ILE	2.6
1	C	-8	GLY	2.5
1	B	-8	GLY	2.4
1	D	-4	VAL	2.4
1	C	297	ASP	2.4
1	C	233	GLU	2.2

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

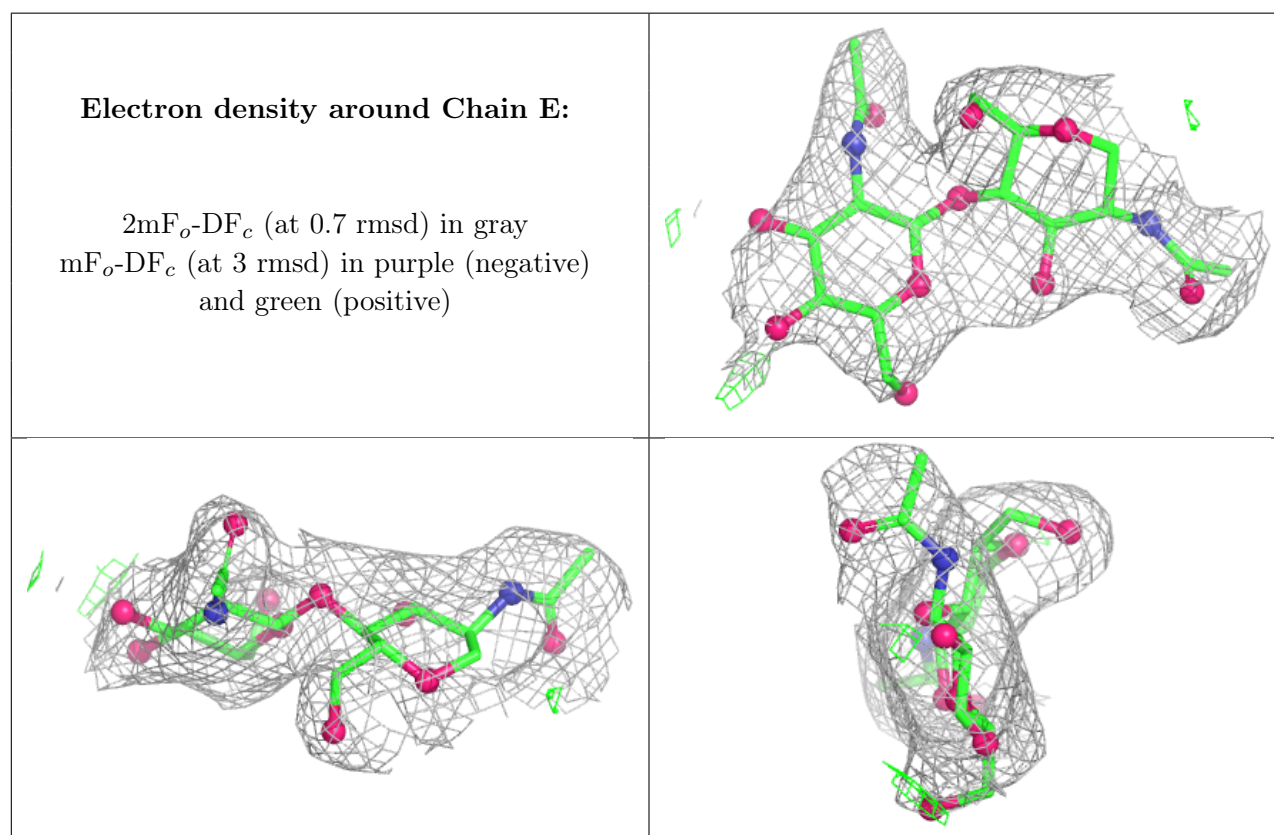
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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

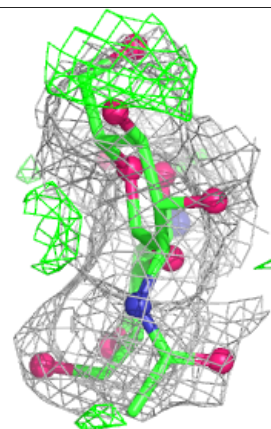
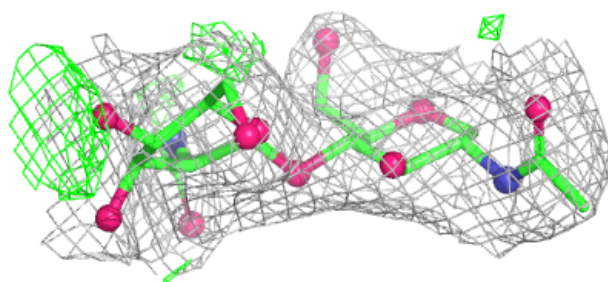
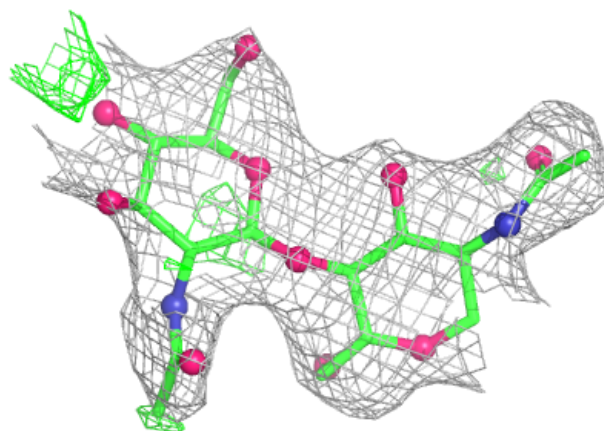
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	F	2	14/15	0.66	0.15	74,77,78,78	0
2	NAG	H	2	14/15	0.73	0.13	77,80,82,83	0
2	NAG	G	2	14/15	0.75	0.13	77,81,83,85	0
2	NAG	E	2	14/15	0.80	0.12	67,71,72,73	0
2	NAG	F	1	14/15	0.91	0.09	51,57,64,70	0
2	NAG	H	1	14/15	0.93	0.08	57,64,67,73	0
2	NAG	G	1	14/15	0.93	0.08	54,61,65,71	0
2	NAG	E	1	14/15	0.95	0.07	45,50,55,61	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.

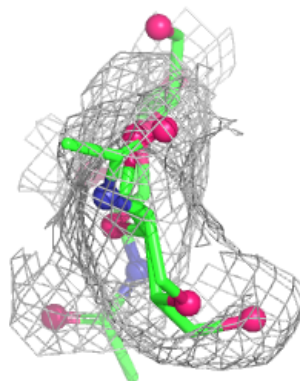
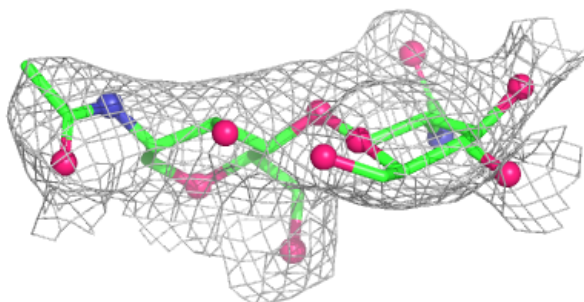
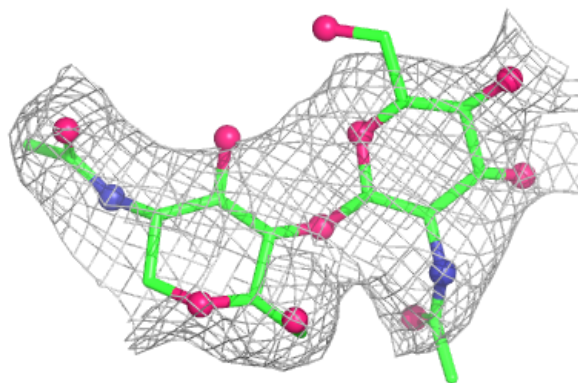


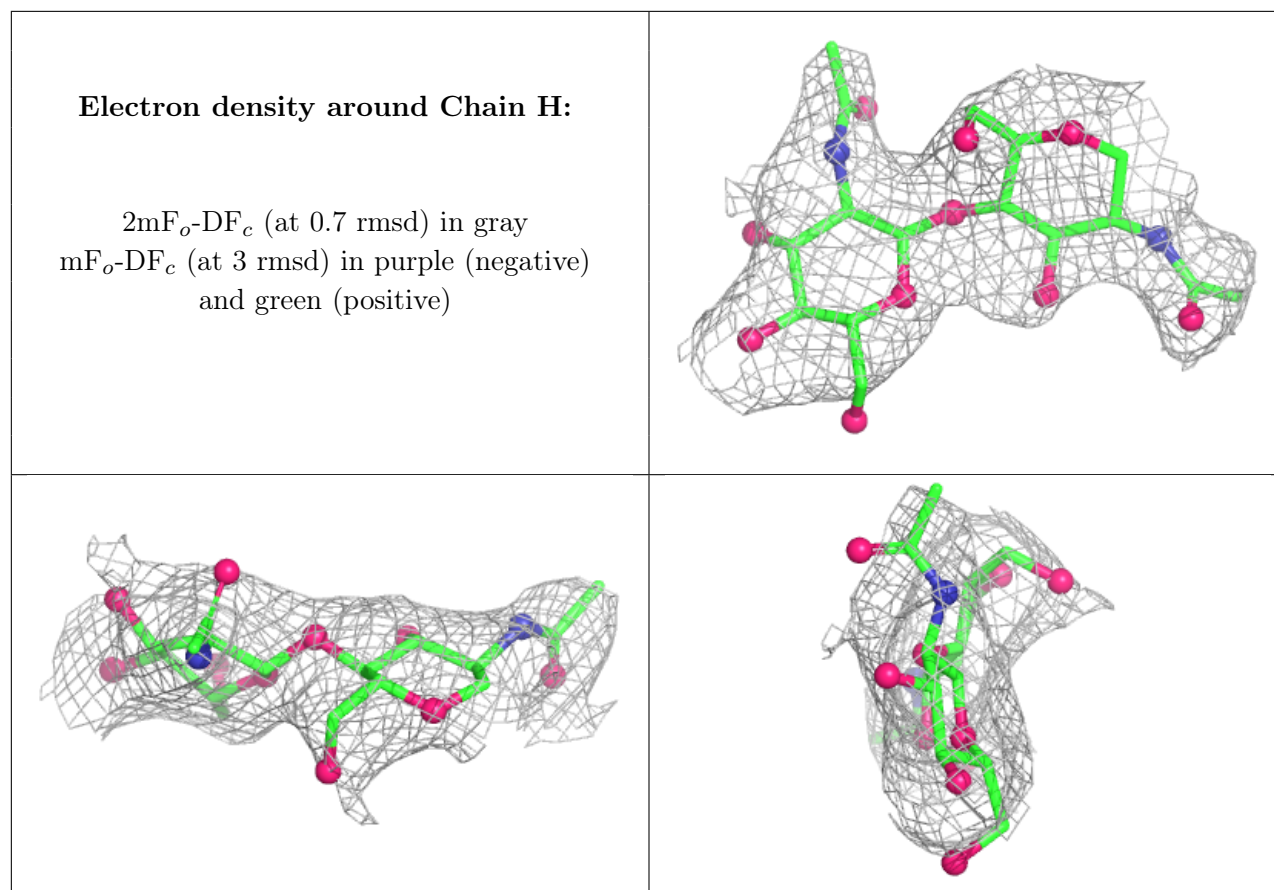
Electron density around Chain F:

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

**Electron density around Chain G:**

$2mF_o - DF_c$ (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(Å ²)	Q<0.9
3	ZN	D	403	1/1	0.99	0.02	51,51,51,51	0
3	ZN	B	403	1/1	1.00	0.01	35,35,35,35	0
3	ZN	C	403	1/1	1.00	0.01	45,45,45,45	0
3	ZN	A	503	1/1	1.00	0.03	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.