



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 10:27 AM JST

PDB ID : 5Z0V
Title : Structural insight into the Zika virus capsid encapsulating the viral genome
Authors : Li, T.; Zhao, Q.; Yang, X.; Chen, C.; Yang, K.; Wu, C.; Zhang, T.; Duan, Y.;
Xue, X.; Mi, K.; Ji, X.; Wang, Z.; Yang, H.
Deposited on : 2017-12-21
Resolution : 2.91 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

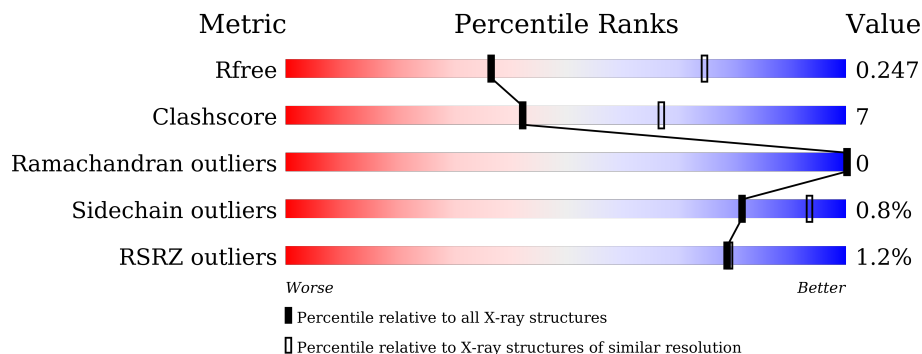
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

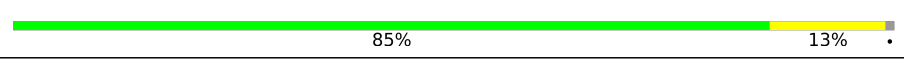
The reported resolution of this entry is 2.91 Å.

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}	130704	2307 (2.94-2.90)
Clashscore	141614	2531 (2.94-2.90)
Ramachandran outliers	138981	2462 (2.94-2.90)
Sidechain outliers	138945	2464 (2.94-2.90)
RSRZ outliers	127900	2248 (2.94-2.90)

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	445	 2% 80% 19%
1	B	445	 0% 85% 14%
1	C	445	 2% 81% 18%
1	D	445	 0% 85% 13%
2	E	2	 100%
2	F	2	 100%

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

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 13947 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 Extracellular solute-binding protein family 1,viral genome protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	444	Total	C	N	O	S	0	0	0
			3446	2229	572	636	9			
1	B	441	Total	C	N	O	S	0	0	0
			3420	2212	568	631	9			
1	C	444	Total	C	N	O	S	0	0	0
			3446	2229	572	636	9			
1	D	440	Total	C	N	O	S	0	0	0
			3411	2207	567	628	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A140NCD0
A	268	ASP	ASN	engineered mutation	UNP A0A140NCD0
B	1	MET	-	initiating methionine	UNP A0A140NCD0
B	268	ASP	ASN	engineered mutation	UNP A0A140NCD0
C	1	MET	-	initiating methionine	UNP A0A140NCD0
C	268	ASP	ASN	engineered mutation	UNP A0A140NCD0
D	1	MET	-	initiating methionine	UNP A0A140NCD0
D	268	ASP	ASN	engineered mutation	UNP A0A140NCD0

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	E	2	Total	C	O	0	0	0
			23	12	11			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	F	2	Total	C	O	0	0	0
			23	12	11			
2	G	2	Total	C	O	0	0	0
			23	12	11			
2	H	2	Total	C	O	0	0	0
			23	12	11			

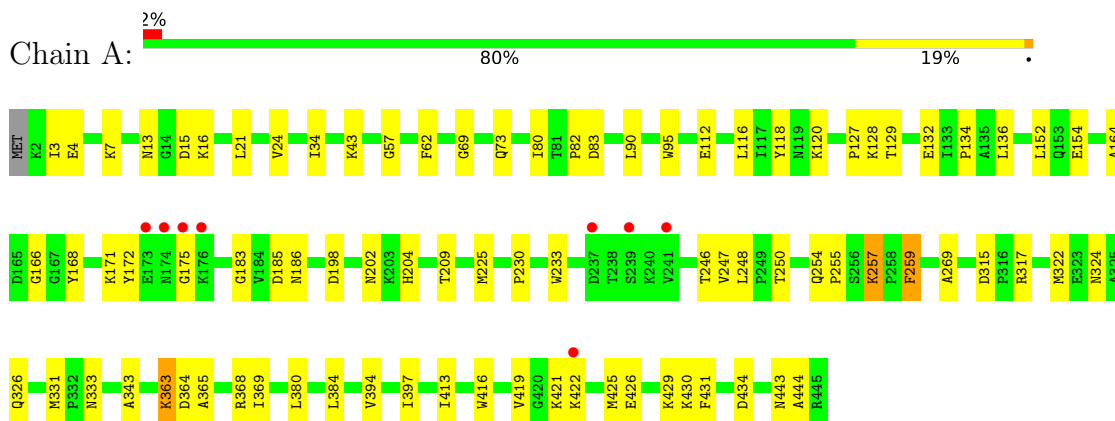
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	29	Total	O	0	0
			29	29		
3	B	26	Total	O	0	0
			26	26		
3	C	26	Total	O	0	0
			26	26		
3	D	51	Total	O	0	0
			51	51		

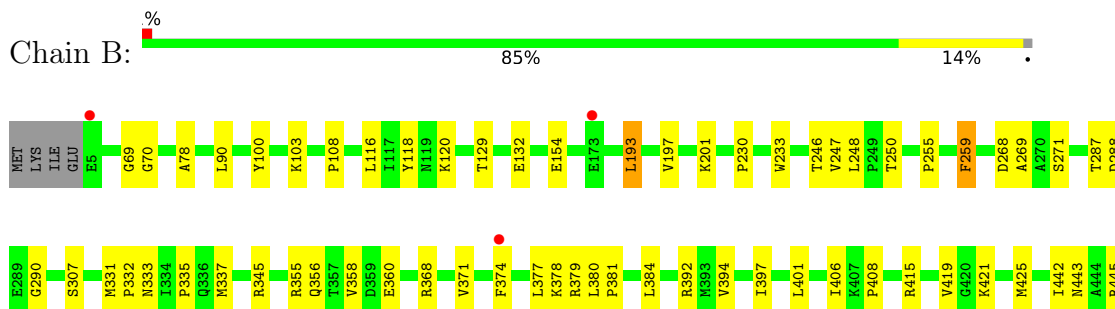
3 Residue-property plots [i](#)

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.

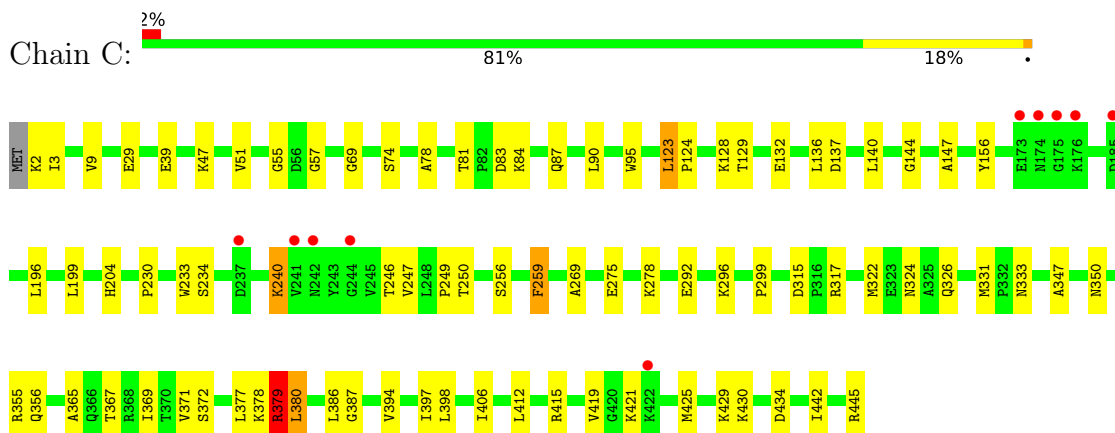
- Molecule 1: Extracellular solute-binding protein family 1,viral genome protein




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- Molecule 1: Extracellular solute-binding protein family 1,viral genome protein



- Molecule 1: Extracellular solute-binding protein family 1,viral genome protein

Chain D:  85% 13%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain E:  100%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain F:  100%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain G:  50% 50%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain H:  50% 50%



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.10Å 223.08Å 85.17Å 90.00° 109.99° 90.00°	Depositor
Resolution (Å)	46.36 – 2.91 46.36 – 2.91	Depositor EDS
% Data completeness (in resolution range)	95.5 (46.36-2.91) 95.5 (46.36-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.91Å)	Xtrriage
Refinement program	PHENIX (1.11.1_2575)	Depositor
R, R_{free}	0.204 , 0.248 0.204 , 0.247	Depositor DCC
R_{free} test set	2340 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	50.9	Xtrriage
Anisotropy	0.316	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 33.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.049 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13947	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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: GLC

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	0.29	0/3526	0.49	0/4776
1	B	0.30	0/3500	0.50	0/4742
1	C	0.30	0/3526	0.51	3/4776 (0.1%)
1	D	0.30	1/3491 (0.0%)	0.49	0/4730
All	All	0.30	1/14043 (0.0%)	0.50	3/19024 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	171	LYS	CE-NZ	-5.15	1.36	1.49

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	123	LEU	CA-CB-CG	6.33	129.85	115.30
1	C	379	ARG	C-N-CA	-5.98	106.75	121.70
1	C	380	LEU	CA-CB-CG	5.48	127.90	115.30

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	3446	0	3492	59	0
1	B	3420	0	3462	39	0
1	C	3446	0	3492	61	0
1	D	3411	0	3456	46	0
2	E	23	0	21	0	0
2	F	23	0	21	0	0
2	G	23	0	21	0	0
2	H	23	0	21	1	0
3	A	29	0	0	4	0
3	B	26	0	0	1	0
3	C	26	0	0	7	0
3	D	51	0	0	7	0
All	All	13947	0	13986	193	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 7.

All (193) 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:286:LEU:O	3:D:601:HOH:O	1.82	0.98
1:D:223:THR:O	3:D:602:HOH:O	1.84	0.93
1:C:156:TYR:O	3:C:602:HOH:O	1.94	0.86
1:C:140:LEU:O	3:C:601:HOH:O	1.92	0.85
1:D:128:LYS:NZ	3:D:603:HOH:O	1.96	0.85
1:C:129:THR:HG22	1:C:132:GLU:HG3	1.61	0.82
1:C:29:GLU:OE1	3:C:604:HOH:O	1.98	0.81
1:A:80:ILE:HG22	1:A:82:PRO:HD3	1.62	0.81
1:A:185:ASP:OD2	1:A:363:LYS:HE3	1.83	0.79
1:A:4:GLU:HB2	1:A:7:LYS:HE2	1.65	0.77
1:D:36:VAL:O	3:D:604:HOH:O	2.02	0.77
1:B:380:LEU:HG	1:C:380:LEU:HD21	1.66	0.77
1:C:129:THR:CG2	1:C:132:GLU:HG3	2.15	0.76
1:A:154:GLU:OE2	3:A:601:HOH:O	2.01	0.76
1:A:83:ASP:OD2	3:A:603:HOH:O	2.03	0.76
1:A:73:GLN:OE1	3:A:604:HOH:O	2.04	0.76
1:A:62:PHE:O	3:A:602:HOH:O	2.03	0.75
1:A:419:VAL:O	1:D:445:ARG:NH2	2.20	0.75
1:C:324:ASN:OD1	3:C:605:HOH:O	2.04	0.74
1:D:96:ASP:OD1	1:D:99:ARG:NH2	2.20	0.74
1:B:380:LEU:HD22	1:B:384:LEU:HD23	1.69	0.74
1:A:430:LYS:NZ	1:A:434:ASP:OD2	2.23	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:144:GLY:N	3:C:601:HOH:O	2.22	0.71
1:B:154:GLU:OE2	1:B:345:ARG:NH1	2.24	0.71
1:A:128:LYS:HD2	1:A:128:LYS:H	1.56	0.71
1:B:377:LEU:HD12	1:C:406:ILE:HD13	1.72	0.70
1:A:380:LEU:H	1:D:379:ARG:HH22	1.39	0.69
1:A:343:ALA:O	1:A:368:ARG:NH2	2.25	0.69
1:B:100:TYR:O	1:B:103:LYS:HG2	1.93	0.69
1:D:377:LEU:O	3:D:605:HOH:O	2.10	0.69
1:A:443:ASN:HA	1:D:421:LYS:HE3	1.76	0.68
1:D:83:ASP:OD1	3:D:606:HOH:O	2.11	0.68
1:A:315:ASP:OD2	1:A:317:ARG:NH2	2.21	0.68
1:B:374:PHE:HE2	1:B:378:LYS:HA	1.58	0.68
1:D:181:ASP:OD1	3:D:607:HOH:O	2.13	0.66
1:D:171:LYS:HD3	1:D:181:ASP:OD2	1.95	0.66
1:C:379:ARG:HG3	1:C:380:LEU:O	1.96	0.66
1:A:422:LYS:HE3	1:A:426:GLU:HG2	1.78	0.65
1:B:443:ASN:HA	1:C:421:LYS:HE3	1.79	0.64
1:D:53:ALA:O	1:D:440:ARG:NH1	2.30	0.64
1:B:374:PHE:CE2	1:B:378:LYS:HA	2.32	0.64
1:B:287:THR:HG23	1:B:290:GLY:H	1.62	0.63
1:D:322:MET:O	1:D:326:GLN:HG3	1.98	0.63
1:D:83:ASP:OD1	1:D:84:LYS:N	2.33	0.62
1:A:3:ILE:HG13	1:A:57:GLY:O	2.01	0.60
1:B:90:LEU:HD23	1:B:108:PRO:HG2	1.83	0.60
1:C:2:LYS:HE3	1:C:9:VAL:HG11	1.84	0.60
1:C:259:PHE:CD1	1:C:331:MET:HG2	2.36	0.60
1:A:128:LYS:HD2	1:A:128:LYS:N	2.17	0.60
1:B:193:LEU:HD23	1:B:358:VAL:HG22	1.85	0.59
1:A:255:PRO:HG2	1:A:257:LYS:HE2	1.85	0.59
1:C:398:LEU:HD21	1:C:412:LEU:HB3	1.86	0.58
1:A:322:MET:O	1:A:326:GLN:HG3	2.04	0.58
1:C:83:ASP:OD1	1:C:84:LYS:N	2.35	0.58
1:C:90:LEU:HD12	1:C:95:TRP:CZ2	2.39	0.57
1:B:250:THR:HG22	1:B:255:PRO:HA	1.87	0.57
1:A:152:LEU:HD12	1:A:209:THR:HB	1.86	0.56
1:D:286:LEU:HD23	1:D:291:LEU:HD21	1.87	0.56
1:C:69:GLY:HA3	1:C:333:ASN:O	2.05	0.55
1:A:246:THR:OG1	1:A:247:VAL:N	2.39	0.55
1:B:421:LYS:O	1:B:425:MET:HG3	2.07	0.55
1:A:90:LEU:HD12	1:A:95:TRP:CZ2	2.42	0.54
1:A:116:LEU:HD11	1:A:225:MET:HE3	1.87	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:442:ILE:HD12	1:C:445:ARG:HH21	1.72	0.54
1:D:174:ASN:O	1:D:174:ASN:ND2	2.40	0.54
1:A:380:LEU:N	1:D:379:ARG:HH22	2.03	0.54
1:C:137:ASP:OD2	1:C:204:HIS:ND1	2.41	0.54
1:C:415:ARG:O	1:C:419:VAL:HG13	2.09	0.53
1:C:249:PRO:O	1:C:256:SER:OG	2.26	0.53
1:D:13:ASN:ND2	1:D:15:ASP:OD1	2.42	0.53
1:D:421:LYS:O	1:D:425:MET:HG3	2.08	0.53
1:A:172:TYR:CZ	1:A:175:GLY:HA2	2.44	0.53
1:C:322:MET:O	1:C:326:GLN:HG3	2.09	0.53
1:C:128:LYS:O	1:C:250:THR:HG22	2.09	0.52
1:A:129:THR:OG1	1:A:132:GLU:HG3	2.09	0.52
1:A:422:LYS:HD2	1:A:425:MET:HB2	1.92	0.52
1:C:2:LYS:NZ	1:C:39:GLU:OE1	2.30	0.52
1:D:129:THR:OG1	1:D:132:GLU:HG3	2.10	0.52
1:C:123:LEU:HD12	1:C:123:LEU:O	2.09	0.51
1:B:288:ASP:OD1	1:B:307:SER:OG	2.28	0.51
1:C:74:SER:OG	3:C:603:HOH:O	1.94	0.51
1:B:377:LEU:HB2	1:C:406:ILE:HG21	1.93	0.50
1:C:246:THR:OG1	1:C:247:VAL:N	2.44	0.50
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.47	0.50
1:C:230:PRO:HA	1:C:233:TRP:CE2	2.46	0.50
1:A:247:VAL:HA	1:A:324:ASN:HD21	1.75	0.50
1:B:118:TYR:CE2	1:B:120:LYS:HG2	2.47	0.50
1:B:406:ILE:HG21	1:C:377:LEU:HB2	1.93	0.50
1:C:367:THR:O	1:C:371:VAL:HG23	2.11	0.50
1:A:127:PRO:HG3	1:A:136:LEU:CD1	2.42	0.50
1:B:230:PRO:HA	1:B:233:TRP:CE2	2.47	0.49
1:A:421:LYS:O	1:A:425:MET:HG3	2.12	0.49
1:B:69:GLY:HA3	1:B:333:ASN:O	2.12	0.49
1:B:129:THR:OG1	1:B:132:GLU:HG3	2.13	0.49
1:A:425:MET:O	1:A:429:LYS:HG3	2.13	0.49
1:C:275:GLU:H	1:C:275:GLU:CD	2.16	0.49
1:A:3:ILE:CG1	1:A:57:GLY:O	2.61	0.49
1:B:355:ARG:NH1	3:B:607:HOH:O	2.45	0.49
1:D:230:PRO:HA	1:D:233:TRP:CE2	2.48	0.49
1:B:442:ILE:O	1:B:445:ARG:HG3	2.12	0.48
1:C:47:LYS:O	1:C:51:VAL:HG22	2.12	0.48
1:A:269:ALA:O	1:A:444:ALA:HB1	2.13	0.48
1:B:415:ARG:O	1:B:419:VAL:HG13	2.13	0.48
1:A:69:GLY:HA3	1:A:333:ASN:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:VAL:HG12	1:B:201:LYS:HE2	1.96	0.48
1:D:425:MET:O	1:D:429:LYS:HG3	2.13	0.48
1:B:116:LEU:HB2	1:B:248:LEU:HD23	1.95	0.47
1:D:98:VAL:HG21	1:D:108:PRO:HD3	1.96	0.47
1:A:413:ILE:O	1:A:416:TRP:HB3	2.14	0.47
1:C:430:LYS:HE3	1:C:434:ASP:OD2	2.15	0.47
1:A:198:ASP:O	1:A:202:ASN:ND2	2.47	0.47
1:A:118:TYR:CE2	1:A:120:LYS:HG2	2.49	0.47
1:A:7:LYS:HA	1:A:34:ILE:HG23	1.97	0.47
1:B:246:THR:OG1	1:B:247:VAL:N	2.47	0.47
1:B:356:GLN:HB3	1:B:360:GLU:HB2	1.95	0.47
1:A:394:VAL:O	1:A:397:ILE:HG22	2.15	0.46
1:C:369:ILE:O	1:C:372:SER:OG	2.28	0.46
1:D:21:LEU:O	1:D:24:VAL:HG22	2.15	0.46
1:C:315:ASP:OD2	1:C:317:ARG:NH2	2.32	0.46
1:A:116:LEU:HB2	1:A:248:LEU:HD23	1.98	0.46
1:A:43:LYS:HE3	1:A:43:LYS:HB2	1.84	0.46
1:D:246:THR:OG1	1:D:247:VAL:N	2.49	0.46
1:C:394:VAL:O	1:C:397:ILE:HG22	2.16	0.46
1:D:63:TRP:CD1	1:D:67:ARG:HG3	2.51	0.46
1:C:84:LYS:HA	1:C:87:GLN:HB2	1.98	0.45
1:D:288:ASP:OD1	1:D:307:SER:OG	2.34	0.45
1:A:171:LYS:HG2	1:A:172:TYR:N	2.31	0.45
1:B:401:LEU:HD13	1:B:408:PRO:HB3	1.97	0.45
1:C:81:THR:O	1:C:278:LYS:NZ	2.50	0.45
1:B:397:ILE:HD12	1:B:397:ILE:HA	1.87	0.45
1:B:380:LEU:HD23	1:B:381:PRO:HD2	1.99	0.45
1:C:3:ILE:CG1	1:C:57:GLY:O	2.65	0.45
1:D:303:VAL:HG22	1:D:312:LEU:HD12	1.98	0.45
1:A:166:GLY:O	1:A:186:ASN:ND2	2.44	0.45
1:B:368:ARG:HA	1:B:371:VAL:HG22	1.99	0.45
1:C:123:LEU:HD22	1:C:136:LEU:HD21	1.98	0.45
1:B:78:ALA:HB2	1:B:269:ALA:HA	1.99	0.45
1:A:343:ALA:CA	1:A:368:ARG:HH21	2.30	0.44
1:D:247:VAL:HA	1:D:324:ASN:HD21	1.82	0.44
1:C:355:ARG:NH1	1:D:96:ASP:OD2	2.50	0.44
1:D:78:ALA:HB2	1:D:269:ALA:HA	1.99	0.44
1:C:137:ASP:HA	1:C:147:ALA:HB2	1.99	0.44
1:A:164:ALA:HA	1:A:257:LYS:HD3	2.00	0.44
1:C:240:LYS:HE3	1:C:240:LYS:HB3	1.67	0.44
1:C:350:ASN:HB3	1:C:356:GLN:HG2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:94:THR:HB	1:D:108:PRO:HB2	2.00	0.43
1:C:425:MET:O	1:C:429:LYS:HG3	2.18	0.43
1:D:109:ILE:HD13	1:D:286:LEU:HD11	2.00	0.43
1:B:70:GLY:HA2	1:B:335:PRO:HB3	1.99	0.43
1:A:168:TYR:CE2	1:A:183:GLY:HA3	2.54	0.43
1:A:134:PRO:HB3	1:A:204:HIS:CE1	2.54	0.43
1:A:365:ALA:O	1:A:369:ILE:HG13	2.19	0.43
1:C:2:LYS:HB3	1:C:55:GLY:O	2.18	0.43
1:B:374:PHE:CD1	1:C:386:LEU:HD21	2.54	0.43
1:C:378:LYS:N	1:C:378:LYS:HD3	2.34	0.43
1:D:69:GLY:HA3	1:D:333:ASN:O	2.19	0.43
1:D:110:ALA:HA	1:D:303:VAL:HA	2.01	0.42
1:A:16:LYS:HE3	1:A:112:GLU:OE2	2.19	0.42
1:A:364:ASP:O	1:A:368:ARG:HG3	2.19	0.42
1:A:255:PRO:O	1:A:257:LYS:HD2	2.19	0.42
1:C:3:ILE:HD11	1:C:57:GLY:O	2.19	0.42
1:D:397:ILE:HD12	1:D:397:ILE:HA	1.90	0.42
1:C:78:ALA:HB2	1:C:269:ALA:HA	2.00	0.42
1:C:347:ALA:HB2	1:C:365:ALA:HB2	2.00	0.42
1:A:13:ASN:ND2	1:A:15:ASP:OD1	2.52	0.42
1:C:123:LEU:HA	1:C:124:PRO:HD3	1.92	0.42
1:C:292:GLU:O	1:C:296:LYS:HG2	2.20	0.42
1:A:259:PHE:CD1	1:A:331:MET:HG2	2.55	0.42
1:D:47:LYS:O	1:D:51:VAL:HG22	2.20	0.42
1:C:387:GLY:O	1:C:415:ARG:NH1	2.47	0.42
1:B:394:VAL:O	1:B:397:ILE:HG22	2.20	0.41
1:B:259:PHE:CD1	1:B:331:MET:HG2	2.55	0.41
1:D:341:TRP:CE3	2:H:2:GLC:H61	2.54	0.41
1:D:436:ALA:O	1:D:440:ARG:HG3	2.20	0.41
1:A:431:PHE:CE1	1:D:403:PHE:HB3	2.55	0.41
1:C:234:SER:OG	1:C:299:PRO:HD3	2.20	0.41
1:C:356:GLN:O	3:C:606:HOH:O	2.22	0.41
1:C:296:LYS:HA	1:C:296:LYS:HD3	1.94	0.41
1:C:199:LEU:HD12	1:C:199:LEU:HA	1.95	0.41
1:A:422:LYS:O	1:A:422:LYS:HG3	2.21	0.41
1:A:21:LEU:O	1:A:24:VAL:HG22	2.21	0.41
1:A:250:THR:OG1	1:A:254:GLN:O	2.29	0.41
1:C:377:LEU:HG	1:C:379:ARG:O	2.21	0.41
1:D:10:ILE:HB	1:D:38:VAL:HG12	2.02	0.41
1:C:129:THR:HG22	1:C:132:GLU:CG	2.40	0.40
1:D:59:ASP:OD1	1:D:271:SER:OG	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:LEU:HB3	1:A:384:LEU:HD23	2.03	0.40
1:B:268:ASP:HB3	1:B:271:SER:HB2	2.03	0.40
1:B:378:LYS:C	1:B:379:ARG:HG3	2.42	0.40
1:D:240:LYS:HB2	1:D:240:LYS:HE3	1.93	0.40
1:B:332:PRO:O	1:B:337:MET:HG3	2.21	0.40
1:D:259:PHE:CD1	1:D:331:MET:HG2	2.57	0.40
1:D:379:ARG:HD3	1:D:380:LEU:N	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	442/445 (99%)	437 (99%)	5 (1%)	0	100	100
1	B	439/445 (99%)	433 (99%)	6 (1%)	0	100	100
1	C	442/445 (99%)	437 (99%)	5 (1%)	0	100	100
1	D	438/445 (98%)	434 (99%)	4 (1%)	0	100	100
All	All	1761/1780 (99%)	1741 (99%)	20 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	355/356 (100%)	352 (99%)	3 (1%)	81	93
1	B	352/356 (99%)	349 (99%)	3 (1%)	78	92
1	C	355/356 (100%)	351 (99%)	4 (1%)	73	91
1	D	351/356 (99%)	350 (100%)	1 (0%)	92	98
All	All	1413/1424 (99%)	1402 (99%)	11 (1%)	81	93

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	257	LYS
1	A	259	PHE
1	A	363	LYS
1	B	193	LEU
1	B	259	PHE
1	B	392	ARG
1	C	196	LEU
1	C	240	LYS
1	C	259	PHE
1	C	379	ARG
1	D	259	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	204	HIS
1	C	101	ASN
1	D	101	ASN

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

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	GLC	E	1	2	12,12,12	0.56	0	17,17,17	0.53	0
2	GLC	E	2	2	11,11,12	0.66	0	15,15,17	0.79	0
2	GLC	F	1	2	12,12,12	0.54	0	17,17,17	0.53	0
2	GLC	F	2	2	11,11,12	0.67	0	15,15,17	0.78	0
2	GLC	G	1	2	12,12,12	0.57	0	17,17,17	0.51	0
2	GLC	G	2	2	11,11,12	0.68	0	15,15,17	1.01	1 (6%)
2	GLC	H	1	2	12,12,12	0.54	0	17,17,17	0.54	0
2	GLC	H	2	2	11,11,12	0.67	0	15,15,17	0.73	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	E	1	2	-	0/2/22/22	0/1/1/1
2	GLC	E	2	2	-	0/2/19/22	0/1/1/1
2	GLC	F	1	2	-	0/2/22/22	0/1/1/1
2	GLC	F	2	2	-	0/2/19/22	0/1/1/1
2	GLC	G	1	2	-	0/2/22/22	0/1/1/1
2	GLC	G	2	2	-	0/2/19/22	0/1/1/1
2	GLC	H	1	2	-	0/2/22/22	0/1/1/1
2	GLC	H	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	GLC	C1-O5-C5	2.87	116.08	112.19

There are no chirality outliers.

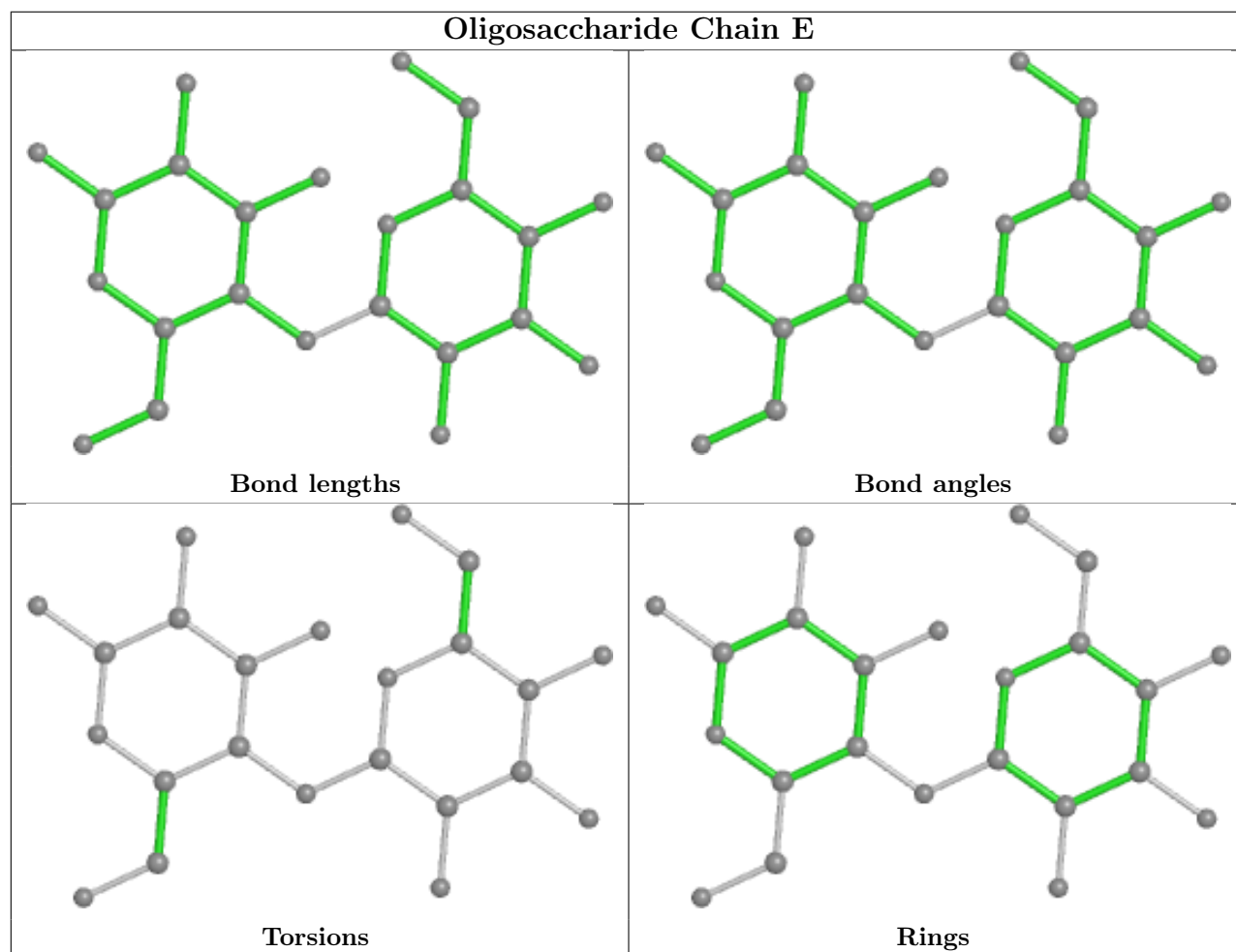
There are no torsion outliers.

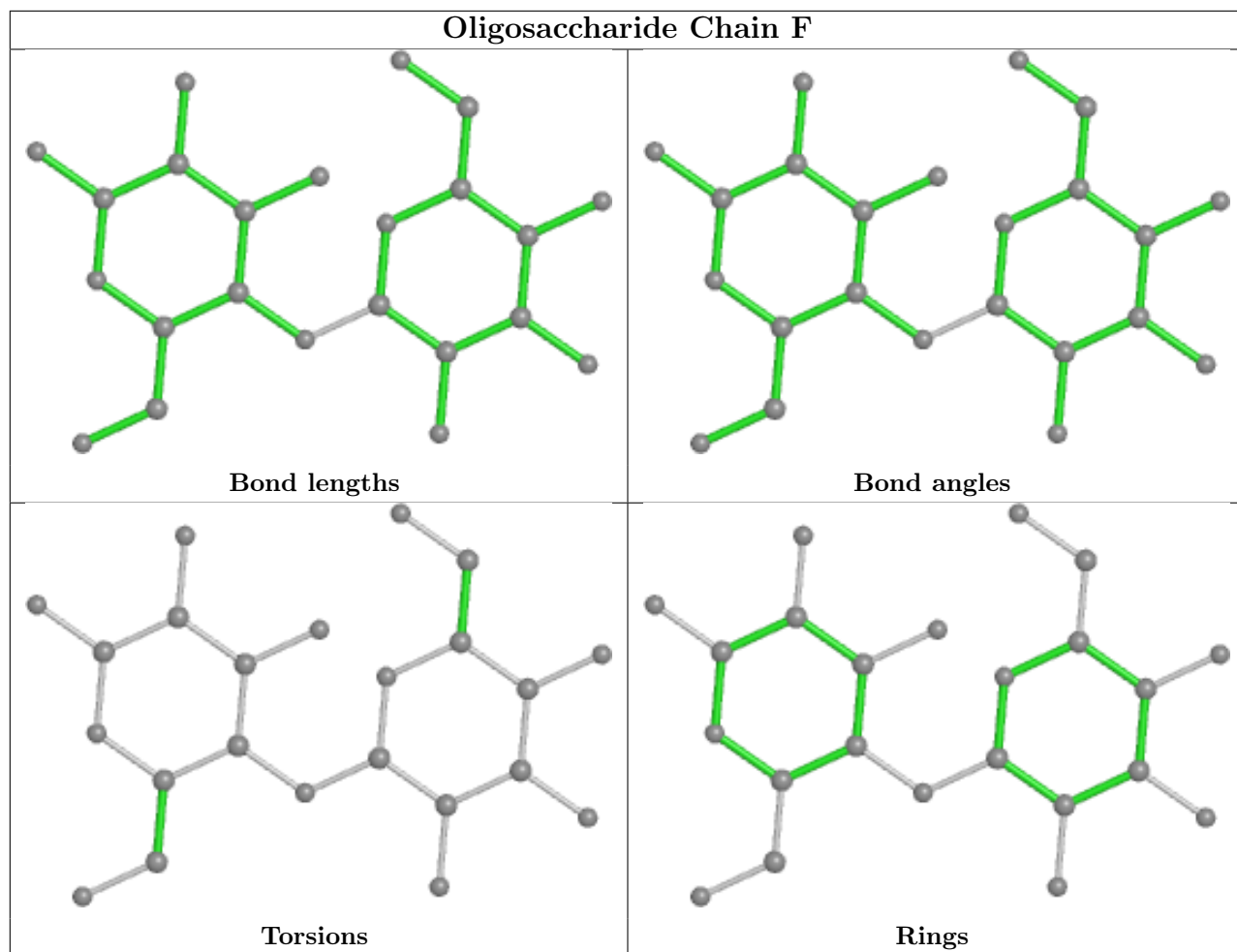
There are no ring outliers.

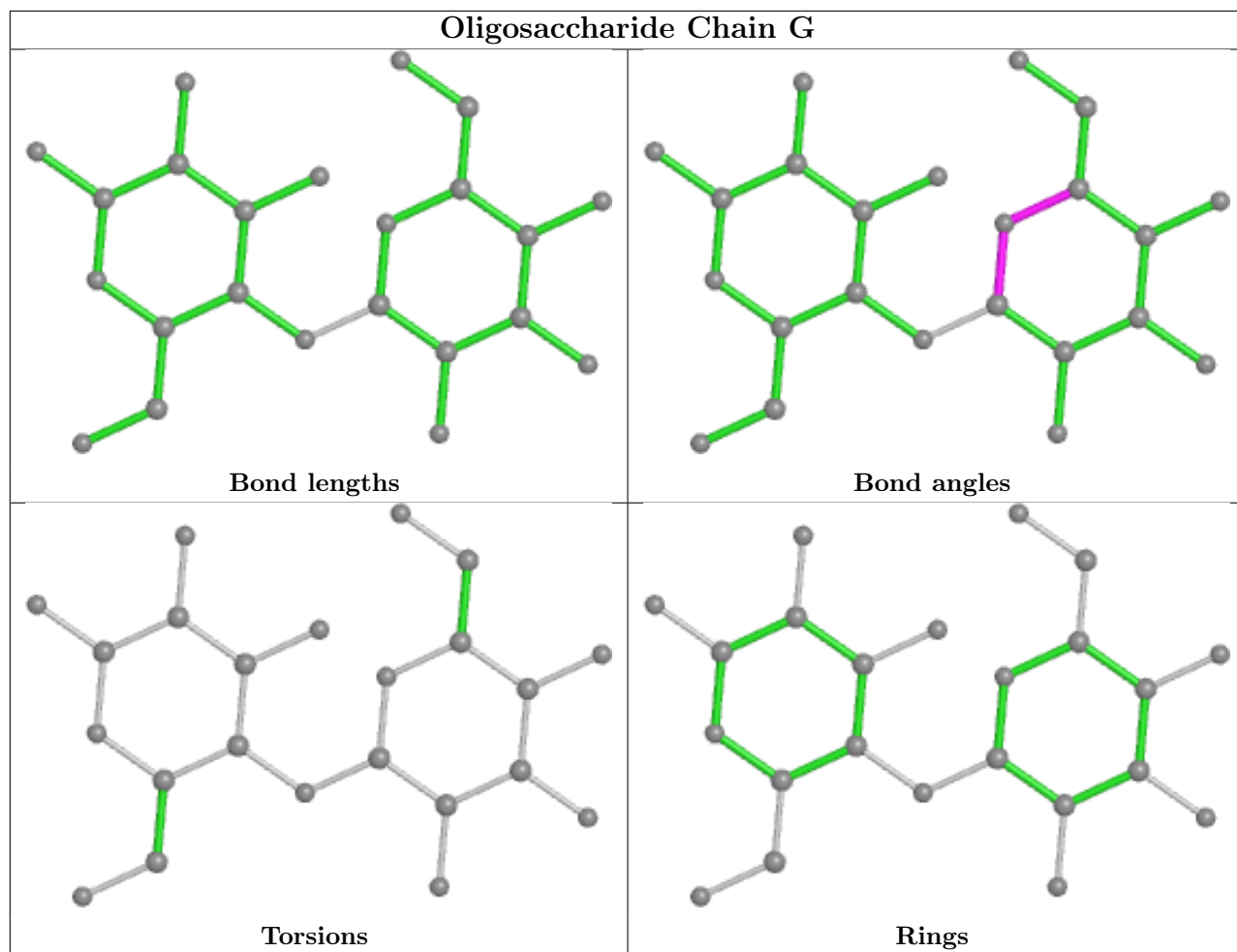
1 monomer is involved in 1 short contact:

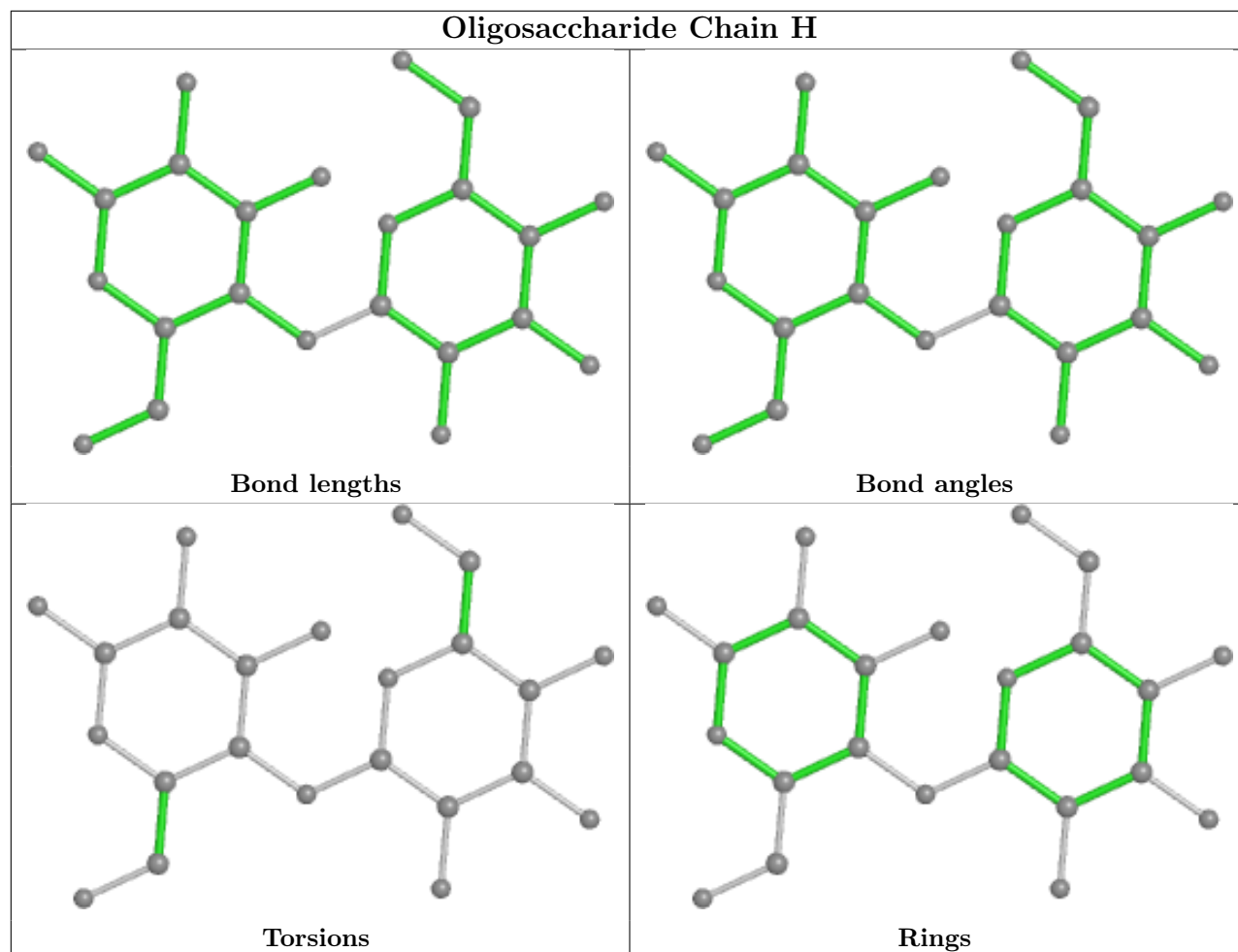
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	2	GLC	1	0

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









5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	444/445 (99%)	0.05	8 (1%) 68 67	30, 49, 74, 110	0
1	B	441/445 (99%)	-0.03	3 (0%) 87 88	30, 43, 62, 92	0
1	C	444/445 (99%)	0.03	10 (2%) 60 59	29, 48, 72, 96	0
1	D	440/445 (98%)	-0.13	1 (0%) 95 95	28, 40, 55, 79	0
All	All	1769/1780 (99%)	-0.02	22 (1%) 79 79	28, 44, 69, 110	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	174	ASN	5.4
1	A	173	GLU	3.8
1	C	175	GLY	3.3
1	A	237	ASP	3.0
1	A	174	ASN	2.9
1	C	173	GLU	2.8
1	C	185	ASP	2.7
1	C	176	LYS	2.7
1	B	5	GLU	2.6
1	C	422	LYS	2.5
1	D	238	THR	2.5
1	A	239	SER	2.5
1	B	173	GLU	2.5
1	B	374	PHE	2.5
1	C	237	ASP	2.4
1	C	242	ASN	2.4
1	A	176	LYS	2.3
1	A	241	VAL	2.3
1	A	175	GLY	2.1
1	C	241	VAL	2.0
1	A	422	LYS	2.0
1	C	244	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

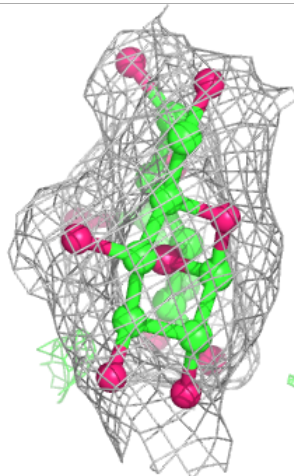
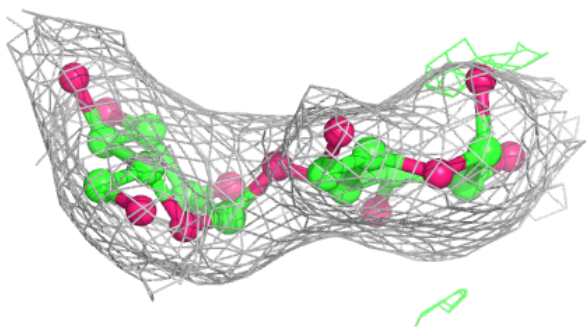
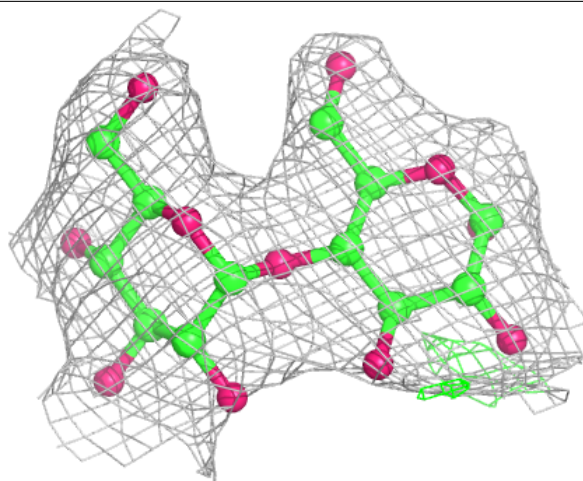
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	E	1	12/12	0.96	0.19	33,36,41,45	0
2	GLC	G	1	12/12	0.96	0.17	28,34,43,44	0
2	GLC	H	1	12/12	0.96	0.15	28,32,34,39	0
2	GLC	F	1	12/12	0.97	0.16	27,30,35,37	0
2	GLC	H	2	11/12	0.97	0.16	28,31,34,34	0
2	GLC	E	2	11/12	0.98	0.15	32,37,46,47	0
2	GLC	G	2	11/12	0.98	0.16	30,36,38,38	0
2	GLC	F	2	11/12	0.99	0.15	28,33,36,36	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.

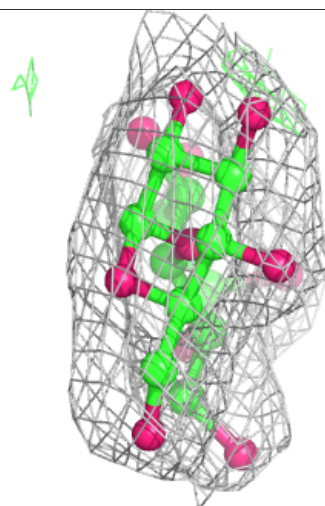
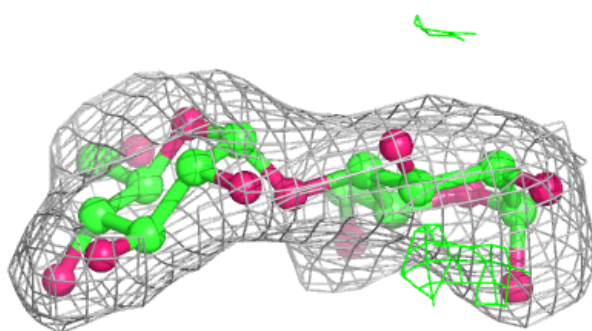
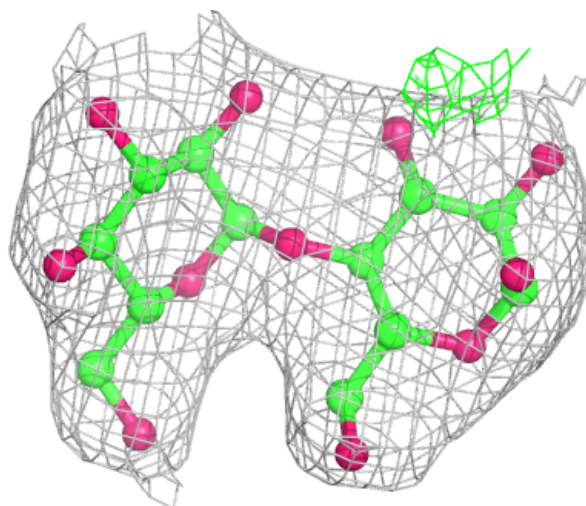
Electron density around Chain E:

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



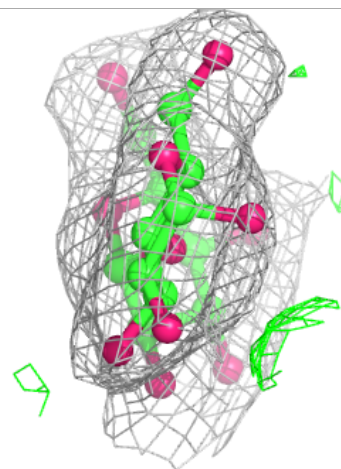
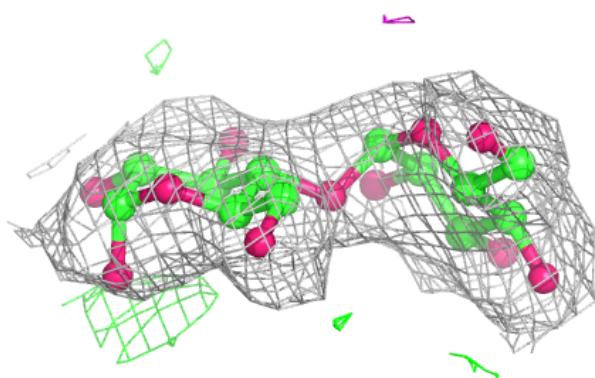
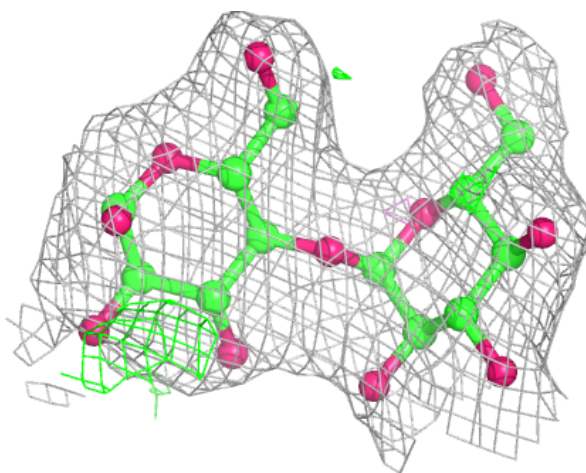
Electron density around Chain F:

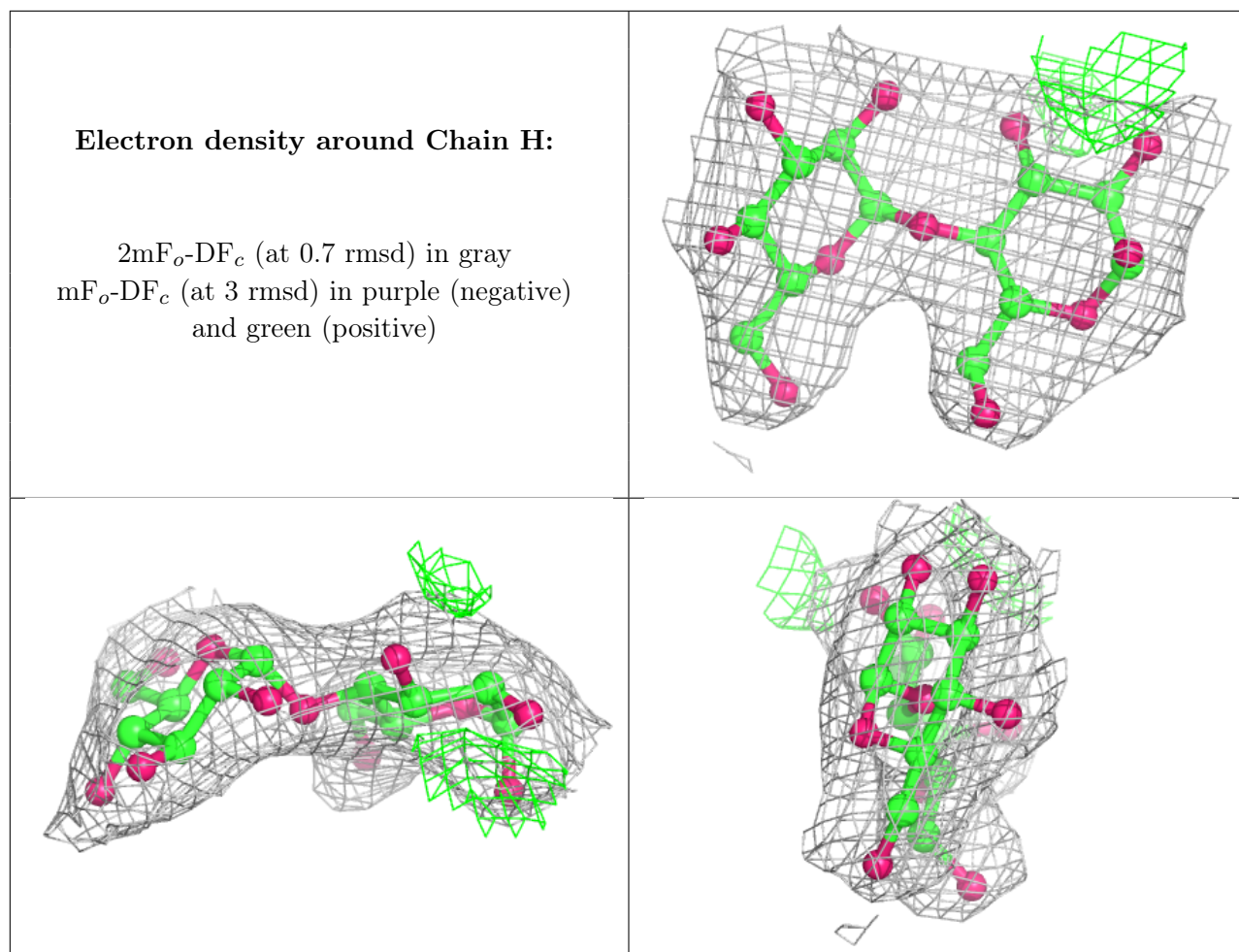
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.