



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:59 PM UTC

PDB ID : 8DF1 / pdb_00008df1
Title : Chi3l1 bound by antibody C59
Authors : Wrapp, D.; McLellan, J.S.
Deposited on : 2022-06-21
Resolution : 3.30 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

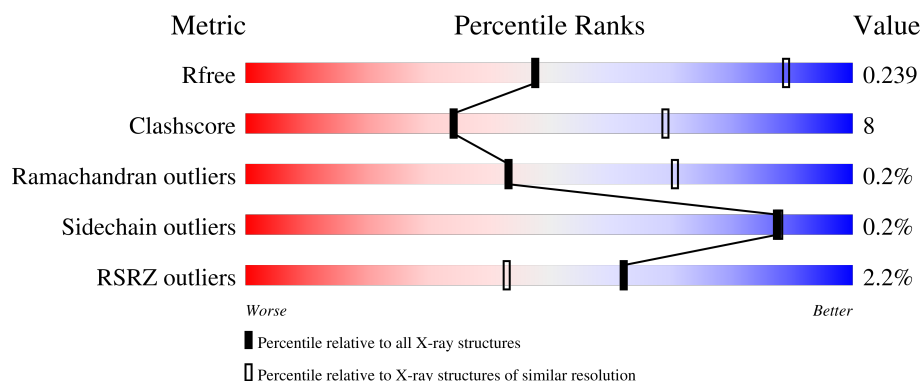
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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.



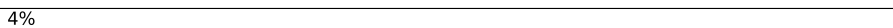
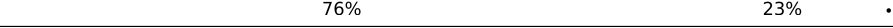
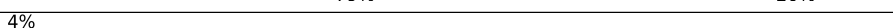
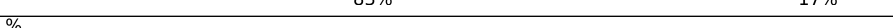
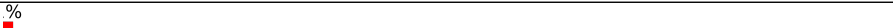
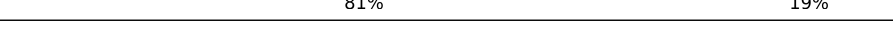
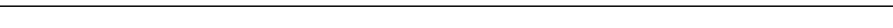
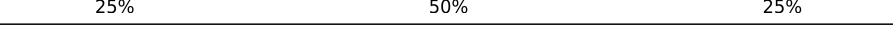
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	362	 82%18%
1	B	362	 86%14%
1	C	362	 84%16%
1	D	362	 81%19%
1	E	362	 %83%17%

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Mol	Chain	Length	Quality of chain
1	F	362	
2	H	223	
2	M	223	
2	O	223	
2	Q	223	
2	S	223	
2	U	223	
3	L	214	
3	N	214	
3	P	214	
3	R	214	
3	T	214	
3	V	214	
4	G	4	
4	W	4	
5	I	3	
5	K	3	
5	X	3	
6	J	5	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 36890 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
			2860	1827	493	529	11			
1	B	362	Total	C	N	O	S	0	0	0
			2860	1827	493	529	11			
1	C	362	Total	C	N	O	S	0	0	0
			2860	1827	493	529	11			
1	D	362	Total	C	N	O	S	0	0	0
			2860	1827	493	529	11			
1	E	362	Total	C	N	O	S	0	0	0
			2860	1827	493	529	11			
1	F	362	Total	C	N	O	S	0	0	0
			2860	1827	493	529	11			

- Molecule 2 is a protein called C59 Fab Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	223	Total	C	N	O	S	0	0	0
			1647	1042	267	327	11			
2	H	223	Total	C	N	O	S	0	0	0
			1647	1042	267	327	11			
2	O	223	Total	C	N	O	S	0	0	0
			1647	1042	267	327	11			
2	Q	223	Total	C	N	O	S	0	0	0
			1647	1042	267	327	11			
2	S	223	Total	C	N	O	S	0	0	0
			1647	1042	267	327	11			
2	U	223	Total	C	N	O	S	0	0	0
			1647	1042	267	327	11			

- Molecule 3 is a protein called C59 Fab Light chain.

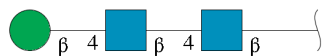
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	N	214	Total	C	N	O	S	0	0	0
			1595	999	257	332	7			
3	L	214	Total	C	N	O	S	0	0	0
			1595	999	257	332	7			
3	P	214	Total	C	N	O	S	0	0	0
			1595	999	257	332	7			
3	R	214	Total	C	N	O	S	0	0	0
			1595	999	257	332	7			
3	T	214	Total	C	N	O	S	0	0	0
			1595	999	257	332	7			
3	V	214	Total	C	N	O	S	0	0	0
			1595	999	257	332	7			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	4	Total	C	N	O	0	0	0
			50	28	2	20			
4	W	4	Total	C	N	O	0	0	0
			50	28	2	20			

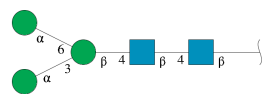
- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	3	Total	C	N	O	0	0	0
			39	22	2	15			
5	K	3	Total	C	N	O	0	0	0
			39	22	2	15			
5	X	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyran

ose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

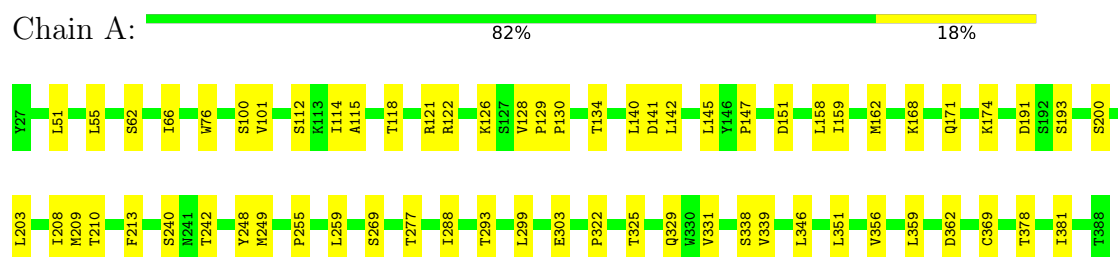


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	J	5	Total	C	N	O	0	0	0
			61	34	2	25			

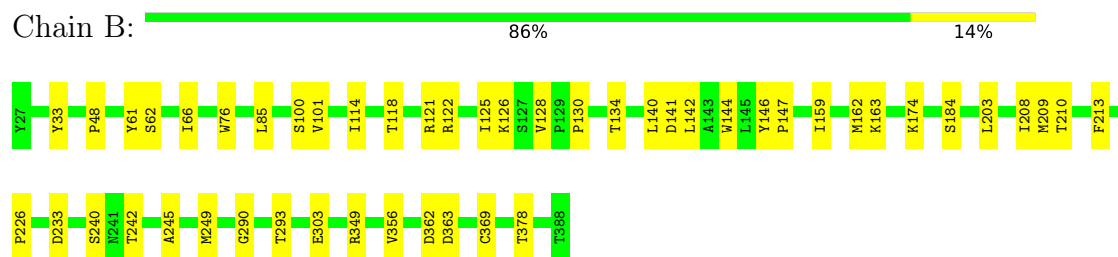
3 Residue-property plots

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.

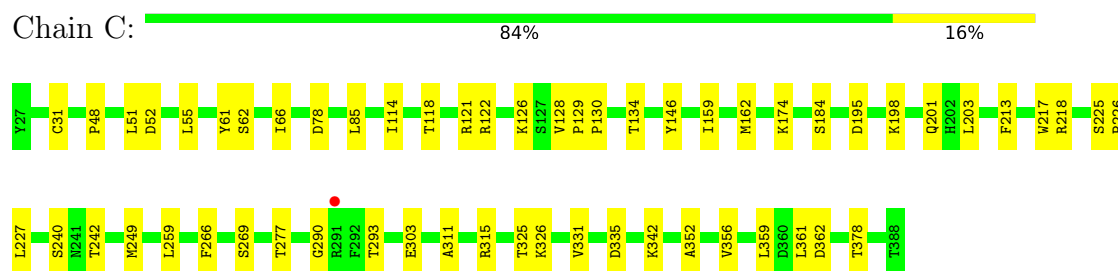
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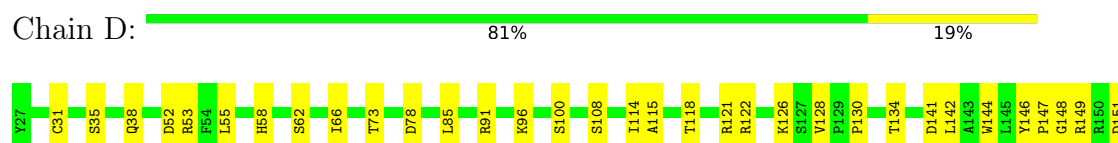
• Molecule 1: Chitinase-3-like protein 1

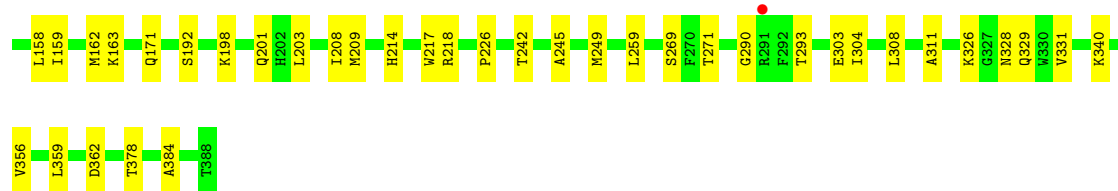


• Molecule 1: Chitinase-3-like protein 1

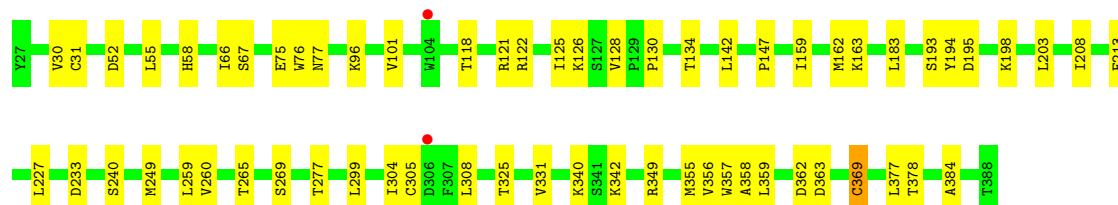
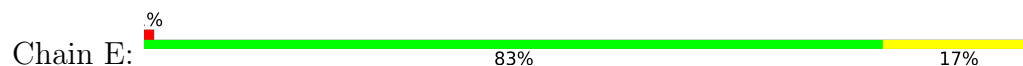


• Molecule 1: Chitinase-3-like protein 1

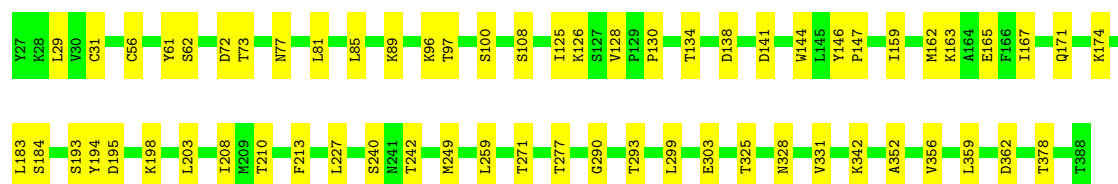
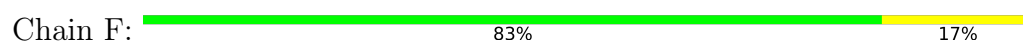




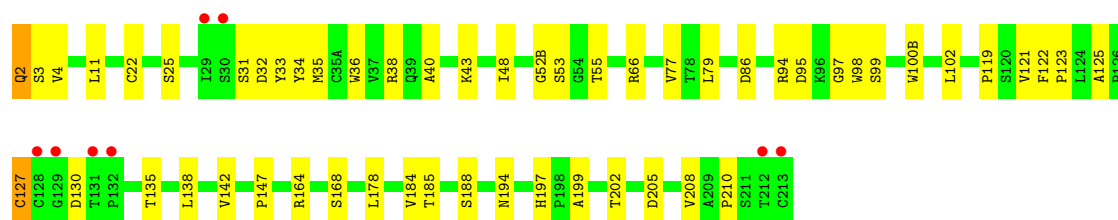
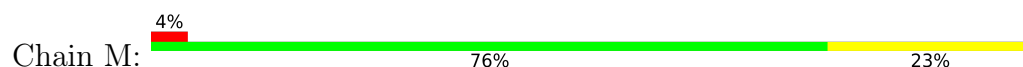
• Molecule 1: Chitinase-3-like protein 1



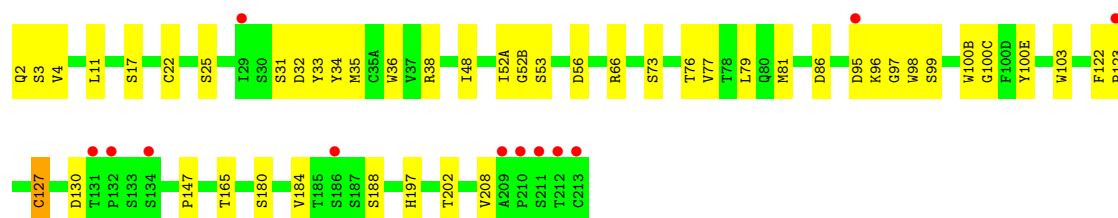
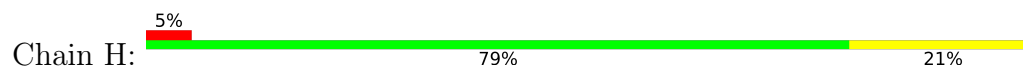
• Molecule 1: Chitinase-3-like protein 1



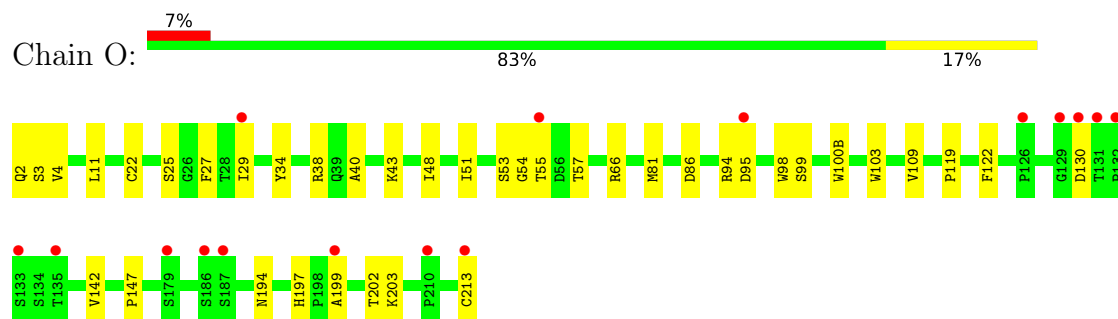
• Molecule 2: C59 Fab Heavy chain



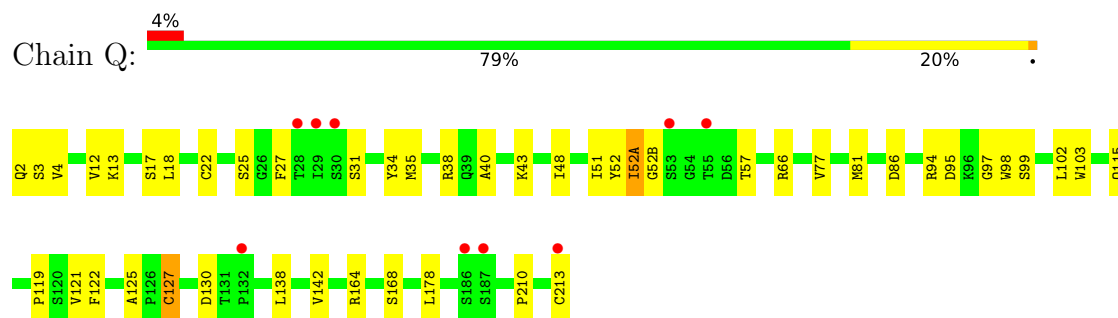
• Molecule 2: C59 Fab Heavy chain



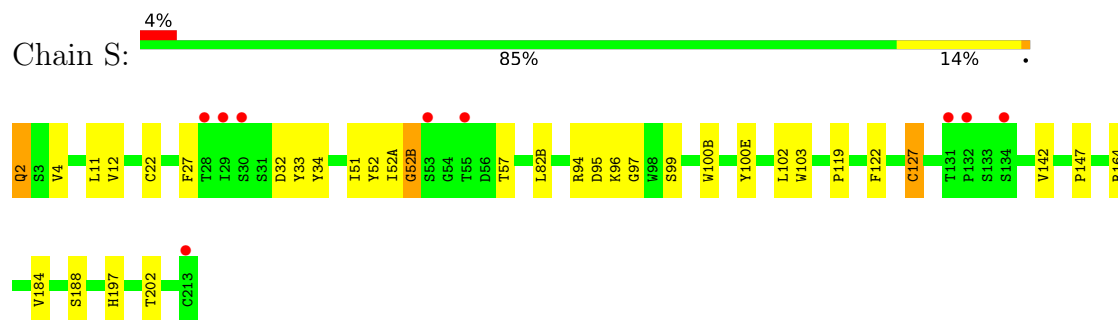
- Molecule 2: C59 Fab Heavy chain



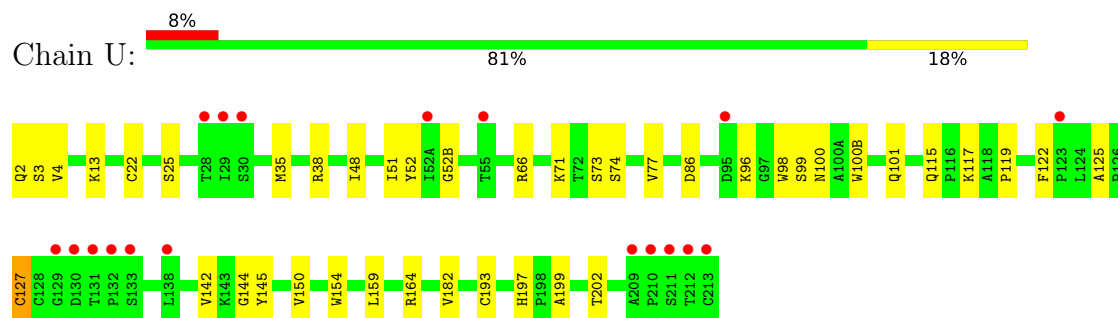
- Molecule 2: C59 Fab Heavy chain



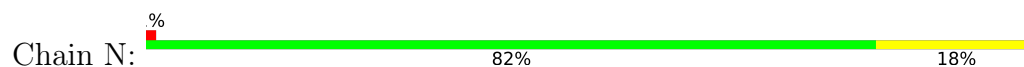
- Molecule 2: C59 Fab Heavy chain



- Molecule 2: C59 Fab Heavy chain

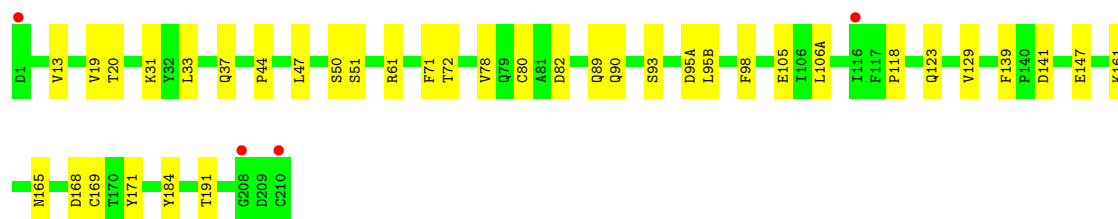
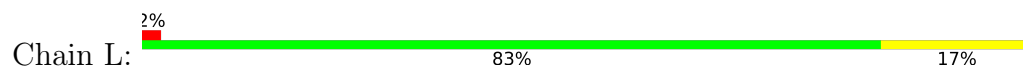


- Molecule 3: C59 Fab Light chain

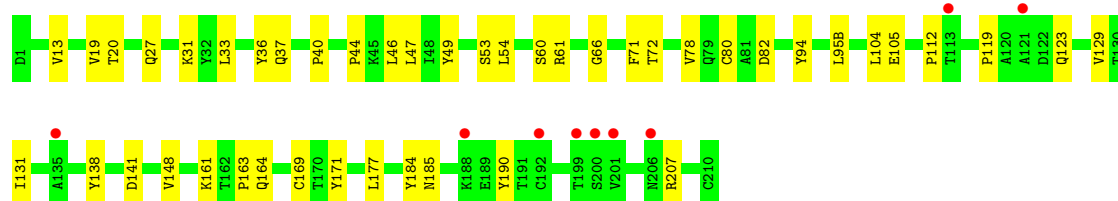
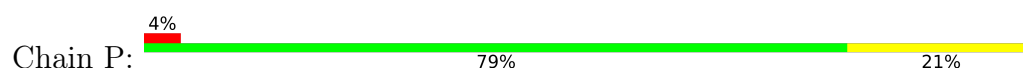




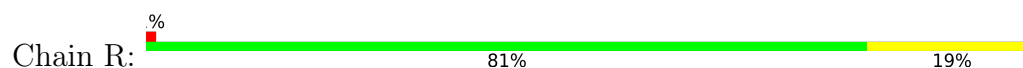
• Molecule 3: C59 Fab Light chain



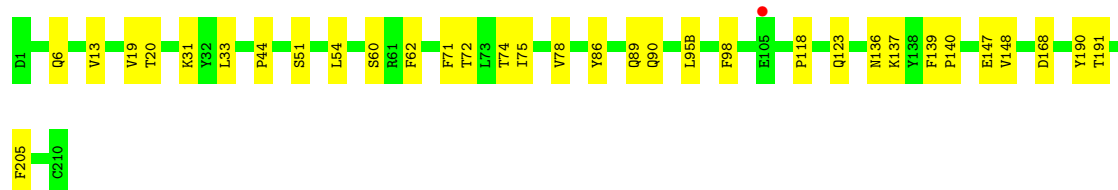
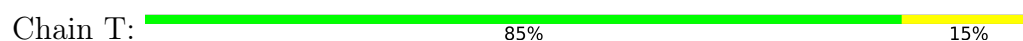
• Molecule 3: C59 Fab Light chain



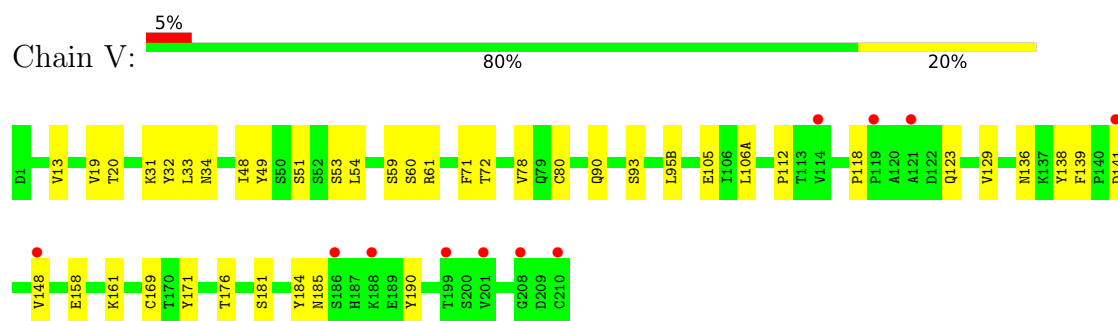
• Molecule 3: C59 Fab Light chain



• Molecule 3: C59 Fab Light chain



- Molecule 3: C59 Fab Light chain



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



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



- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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

Chain J:  80% 20%

MAG1
MAG2
EMJ3
MAN4
MAN5

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.10Å 172.92Å 404.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	87.29 – 3.30 87.29 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (87.29-3.30) 99.9 (87.29-3.30)	Depositor EDS
R_{merge}	0.80	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.20.1-4487	Depositor
R, R_{free}	0.206 , 0.243 0.203 , 0.239	Depositor DCC
R_{free} test set	5506 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.666	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	36890	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.46% 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

5.1 Standard geometry

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

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	0.35	0/2936	0.58	2/3978 (0.1%)
1	B	0.34	0/2936	0.57	2/3978 (0.1%)
1	C	0.34	0/2936	0.58	2/3978 (0.1%)
1	D	0.35	0/2936	0.61	2/3978 (0.1%)
1	E	0.34	0/2936	0.58	0/3978
1	F	0.36	0/2936	0.58	2/3978 (0.1%)
2	H	0.28	0/1684	0.51	0/2311
2	M	0.30	0/1684	0.54	0/2311
2	O	0.27	0/1684	0.50	0/2311
2	Q	0.27	0/1684	0.51	0/2311
2	S	0.27	0/1684	0.51	0/2311
2	U	0.31	0/1684	0.53	0/2311
3	L	0.30	0/1629	0.53	0/2231
3	N	0.32	0/1629	0.53	0/2231
3	P	0.28	0/1629	0.51	0/2231
3	R	0.30	0/1629	0.53	0/2231
3	T	0.28	0/1629	0.52	0/2231
3	V	0.30	0/1629	0.51	0/2231
All	All	0.32	0/37494	0.55	10/51120 (0.0%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	62	SER	CA-C-N	6.30	133.03	121.70
1	B	62	SER	C-N-CA	6.30	133.03	121.70
1	F	62	SER	CA-C-N	5.81	132.16	121.70
1	F	62	SER	C-N-CA	5.81	132.16	121.70
1	C	62	SER	CA-C-N	5.78	132.10	121.70
1	C	62	SER	C-N-CA	5.78	132.10	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	62	SER	CA-C-N	5.28	131.20	121.70
1	D	62	SER	C-N-CA	5.28	131.20	121.70
1	A	62	SER	CA-C-N	5.06	130.81	121.70
1	A	62	SER	C-N-CA	5.06	130.81	121.70

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	2860	0	2781	45	0
1	B	2860	0	2781	36	2
1	C	2860	0	2781	39	0
1	D	2860	0	2781	47	0
1	E	2860	0	2781	45	2
1	F	2860	0	2781	39	0
2	H	1647	0	1597	31	0
2	M	1647	0	1597	37	0
2	O	1647	0	1597	28	0
2	Q	1647	0	1597	31	0
2	S	1647	0	1597	22	0
2	U	1647	0	1597	26	0
3	L	1595	0	1532	26	0
3	N	1595	0	1532	27	0
3	P	1595	0	1532	32	0
3	R	1595	0	1532	35	0
3	T	1595	0	1532	26	0
3	V	1595	0	1532	34	0
4	G	50	0	43	1	0
4	W	50	0	43	1	0
5	I	39	0	34	1	0
5	K	39	0	34	1	0
5	X	39	0	34	4	0
6	J	61	0	52	1	0
All	All	36890	0	35700	551	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (551) 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)
3:L:13:VAL:HG21	3:L:19:VAL:HG22	1.46	0.98
3:N:13:VAL:HG21	3:N:19:VAL:HG22	1.55	0.89
3:T:13:VAL:HG21	3:T:19:VAL:HG22	1.57	0.87
2:M:164:ARG:NH2	3:N:136:ASN:OD1	2.08	0.86
3:P:13:VAL:HG21	3:P:19:VAL:HG22	1.56	0.86
3:V:13:VAL:HG21	3:V:19:VAL:HG22	1.56	0.85
3:R:13:VAL:HG21	3:R:19:VAL:HG22	1.59	0.84
1:A:128:VAL:HG21	1:A:162:MET:HE1	1.64	0.77
2:U:38:ARG:HB3	2:U:48:ILE:HD11	1.66	0.77
1:B:208:ILE:HD12	1:B:249:MET:HE2	1.68	0.75
1:D:159:ILE:HD12	1:D:203:LEU:HD21	1.70	0.73
1:F:249:MET:HE3	1:F:259:LEU:HD21	1.71	0.73
1:A:249:MET:HE3	1:A:259:LEU:HD21	1.72	0.71
1:F:147:PRO:HG2	1:F:193:SER:HB3	1.73	0.71
1:A:213:PHE:HB3	1:A:240:SER:HB3	1.74	0.70
3:V:105:GLU:OE2	3:V:171:TYR:OH	2.08	0.69
1:A:147:PRO:HB2	1:A:193:SER:HB3	1.74	0.69
5:X:1:NAG:H3	5:X:1:NAG:H83	1.74	0.69
3:L:90:GLN:HE21	3:L:95(B):LEU:HD11	1.58	0.69
2:S:33:TYR:CZ	2:S:52(A):ILE:HG13	2.28	0.69
1:D:362:ASP:OD2	1:D:378:THR:HG23	1.93	0.68
3:P:119:PRO:HD3	3:P:131:ILE:HG22	1.76	0.68
1:B:118:THR:HG22	1:B:121:ARG:HH22	1.59	0.68
3:P:20:THR:HG23	3:P:72:THR:HG23	1.75	0.68
2:M:94:ARG:HD3	2:M:102:LEU:HD12	1.76	0.67
1:F:293:THR:HG23	1:F:303:GLU:OE2	1.93	0.67
1:A:130:PRO:O	1:A:134:THR:HG23	1.94	0.67
3:P:131:ILE:HD11	3:P:177:LEU:HD23	1.76	0.67
1:E:134:THR:HG22	3:T:31:LYS:HG2	1.77	0.67
1:C:293:THR:HG23	1:C:303:GLU:OE2	1.95	0.66
1:B:362:ASP:OD2	1:B:378:THR:HG23	1.94	0.66
1:B:293:THR:HG23	1:B:303:GLU:OE2	1.96	0.66
2:O:213:CYS:O	3:R:207:ARG:NH1	2.29	0.66
2:Q:95:ASP:HB3	2:Q:97:GLY:H	1.60	0.66
2:O:99:SER:HG	2:O:100(B):TRP:CD1	2.15	0.66
1:A:159:ILE:HD12	1:A:203:LEU:HD21	1.78	0.65
1:E:128:VAL:HG21	1:E:162:MET:HE1	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:38:ARG:HB3	2:M:48:ILE:HD11	1.79	0.65
1:C:290:GLY:HA3	1:C:303:GLU:OE1	1.97	0.65
2:U:3:SER:HB3	2:U:25:SER:HB2	1.79	0.65
1:F:159:ILE:HD12	1:F:203:LEU:HD21	1.79	0.65
2:O:197:HIS:HB3	2:O:202:THR:HB	1.77	0.65
3:L:20:THR:HG23	3:L:72:THR:HG23	1.79	0.64
3:V:13:VAL:HG11	3:V:78:VAL:HG21	1.80	0.64
2:U:66:ARG:HH22	2:U:86:ASP:CG	2.05	0.64
1:C:266:PHE:HB3	1:C:361:LEU:HD13	1.78	0.64
2:O:213:CYS:HB2	3:R:185:ASN:HB3	1.79	0.64
3:V:20:THR:HG23	3:V:72:THR:HG23	1.79	0.64
2:H:99:SER:HG	2:H:100(B):TRP:CD1	2.15	0.64
2:H:3:SER:HB3	2:H:25:SER:HB2	1.80	0.63
3:P:37:GLN:HB2	3:P:47:LEU:HD11	1.80	0.63
1:A:210:THR:HB	1:A:242:THR:HG22	1.80	0.63
2:M:168:SER:HB3	2:M:178:LEU:HD23	1.79	0.63
3:P:207:ARG:NH1	2:Q:213:CYS:O	2.32	0.63
1:D:214:HIS:CE1	1:D:218:ARG:HD3	2.34	0.62
2:Q:121:VAL:HG22	2:Q:142:VAL:HG22	1.80	0.62
1:C:277:THR:HB	1:C:325:THR:HG21	1.81	0.62
1:D:208:ILE:HD12	1:D:249:MET:HE2	1.81	0.62
2:S:52:TYR:CE2	2:S:52(B):GLY:HA3	2.35	0.62
2:M:121:VAL:HG22	2:M:142:VAL:HG22	1.81	0.62
1:B:128:VAL:HG21	1:B:162:MET:HE1	1.82	0.62
5:K:1:NAG:H83	5:K:1:NAG:H3	1.81	0.61
3:N:20:THR:HG23	3:N:72:THR:HG23	1.82	0.61
1:B:130:PRO:O	1:B:134:THR:HG23	1.99	0.61
1:A:208:ILE:HD12	1:A:249:MET:HE2	1.83	0.61
1:D:130:PRO:O	1:D:134:THR:HG23	2.00	0.60
2:S:99:SER:HG	2:S:100(B):TRP:CD1	2.18	0.60
2:M:34:TYR:HB2	2:M:95:ASP:OD2	2.01	0.60
2:Q:52:TYR:CE2	2:Q:52(B):GLY:HA3	2.36	0.60
1:C:362:ASP:OD2	1:C:378:THR:HG23	2.02	0.60
1:C:159:ILE:HD12	1:C:203:LEU:HD21	1.83	0.60
1:B:159:ILE:HD12	1:B:203:LEU:HD21	1.83	0.60
1:D:142:LEU:HD12	1:D:159:ILE:HD13	1.83	0.60
1:D:249:MET:HE3	1:D:259:LEU:HD21	1.83	0.60
3:V:90:GLN:HE21	3:V:95(B):LEU:HD11	1.67	0.60
1:A:293:THR:HG23	1:A:303:GLU:OE2	2.02	0.59
1:B:213:PHE:HB3	1:B:240:SER:HB3	1.83	0.59
3:L:90:GLN:NE2	3:L:95(B):LEU:HD11	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:R:20:THR:HG23	3:R:72:THR:HG23	1.84	0.59
1:E:195:ASP:OD2	1:E:198:LYS:HD3	2.03	0.59
2:O:53:SER:O	2:O:55:THR:N	2.36	0.59
2:U:66:ARG:NH2	2:U:86:ASP:OD2	2.35	0.59
1:E:134:THR:HA	3:T:31:LYS:HE2	1.85	0.59
2:O:197:HIS:CE1	2:O:199:ALA:HB3	2.38	0.59
1:C:195:ASP:OD2	1:C:198:LYS:HD3	2.02	0.59
3:N:37:GLN:HB2	3:N:47:LEU:HD11	1.84	0.59
2:H:38:ARG:HB3	2:H:48:ILE:HD11	1.84	0.59
2:S:95:ASP:HB3	2:S:97:GLY:H	1.68	0.59
2:Q:38:ARG:HB3	2:Q:48:ILE:HD11	1.85	0.58
2:U:13:LYS:NZ	2:U:115:GLN:OE1	2.24	0.58
1:C:130:PRO:O	1:C:134:THR:HG23	2.03	0.58
1:E:249:MET:HE3	1:E:259:LEU:HD21	1.85	0.58
2:Q:34:TYR:H	2:Q:95:ASP:HB2	1.68	0.58
1:E:362:ASP:OD2	1:E:378:THR:HG23	2.03	0.58
2:M:53:SER:O	2:M:55:THR:N	2.32	0.58
1:D:217:TRP:CZ3	1:D:218:ARG:HG2	2.39	0.58
1:F:210:THR:HB	1:F:242:THR:HG22	1.85	0.58
1:A:134:THR:HA	3:L:31:LYS:HE2	1.84	0.58
1:C:213:PHE:HB3	1:C:240:SER:HB3	1.84	0.58
4:G:1:NAG:H3	4:G:1:NAG:H83	1.86	0.58
2:Q:127:CYS:HB3	3:R:118:PRO:HG2	1.86	0.58
1:B:174:LYS:HG2	2:M:98:TRP:CD1	2.39	0.57
1:B:290:GLY:HA3	1:B:303:GLU:OE1	2.04	0.57
1:C:128:VAL:HG21	1:C:162:MET:HE1	1.87	0.57
1:F:126:LYS:NZ	1:F:165:GLU:OE2	2.34	0.57
2:M:99:SER:HG	2:M:100(B):TRP:CD1	2.23	0.57
2:U:164:ARG:NH2	3:V:136:ASN:OD1	2.37	0.57
1:F:213:PHE:HB3	1:F:240:SER:HB3	1.86	0.57
2:O:27:PHE:CE2	2:O:94:ARG:HD2	2.40	0.57
3:N:147:GLU:HB2	3:N:191:THR:HB	1.86	0.57
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.85	0.57
3:L:61:ARG:NH1	3:L:82:ASP:OD2	2.31	0.57
2:U:4:VAL:HG13	2:U:22:CYS:SG	2.45	0.57
1:D:293:THR:HG23	1:D:303:GLU:OE2	2.05	0.57
1:E:122:ARG:NH2	1:E:126:LYS:HD3	2.20	0.57
1:E:208:ILE:HD12	1:E:249:MET:HE2	1.87	0.57
1:A:134:THR:HG22	3:L:31:LYS:HG2	1.86	0.56
1:A:118:THR:HG22	1:A:121:ARG:HH22	1.70	0.56
1:F:134:THR:HG22	3:V:31:LYS:HG2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:GLN:OE1	3:L:93:SER:HB2	2.05	0.56
3:L:13:VAL:HG11	3:L:78:VAL:HG21	1.87	0.56
2:O:34:TYR:HB2	2:O:95:ASP:OD2	2.05	0.56
3:N:54:LEU:HD11	3:N:60:SER:HA	1.88	0.56
3:T:54:LEU:HD11	3:T:60:SER:HA	1.88	0.56
2:M:127:CYS:HB3	3:N:118:PRO:HG2	1.88	0.56
1:D:118:THR:HG22	1:D:121:ARG:HH22	1.71	0.56
1:B:122:ARG:NH1	1:B:126:LYS:HD3	2.21	0.56
1:C:122:ARG:NH2	1:C:126:LYS:HD3	2.21	0.56
3:P:61:ARG:NH1	3:P:82:ASP:OD2	2.32	0.55
3:T:20:THR:HG23	3:T:72:THR:HG23	1.88	0.55
1:D:35:SER:O	1:D:38:GLN:HG2	2.06	0.55
1:F:277:THR:HB	1:F:325:THR:HG21	1.88	0.55
2:Q:122:PHE:CD2	3:R:123:GLN:HG3	2.42	0.55
1:B:174:LYS:HG2	2:M:98:TRP:CG	2.41	0.55
1:C:227:LEU:HD13	1:C:342:LYS:HG2	1.87	0.55
2:H:66:ARG:NH2	2:H:86:ASP:OD2	2.40	0.55
2:Q:51:ILE:HG13	2:Q:57:THR:HG22	1.89	0.55
3:R:12:GLU:CD	3:R:106(A):LEU:HD21	2.32	0.55
5:I:1:NAG:H83	5:I:1:NAG:H3	1.87	0.55
3:L:105:GLU:OE2	3:L:171:TYR:OH	2.25	0.55
1:A:142:LEU:HD12	1:A:159:ILE:HD13	1.89	0.54
1:D:328:ASN:OD1	1:D:329:GLN:NE2	2.37	0.54
1:B:118:THR:HG22	1:B:121:ARG:NH2	2.22	0.54
1:F:227:LEU:HD13	1:F:342:LYS:HG2	1.90	0.54
2:O:4:VAL:HG13	2:O:22:CYS:SG	2.47	0.54
2:Q:13:LYS:NZ	2:Q:115:GLN:OE1	2.33	0.54
3:V:80:CYS:SG	3:V:169:CYS:N	2.80	0.54
2:M:123:PRO:HB3	2:M:208:VAL:HG22	1.90	0.54
2:Q:3:SER:HB3	2:Q:25:SER:HB2	1.90	0.54
3:R:184:TYR:O	3:R:190:TYR:OH	2.24	0.54
1:D:128:VAL:HG21	1:D:162:MET:HE1	1.89	0.54
1:D:134:THR:HG21	3:R:31:LYS:H	1.73	0.54
2:H:31:SER:HB2	2:H:52(B):GLY:HA2	1.89	0.54
2:M:184:VAL:HB	2:M:188:SER:HB2	1.90	0.54
1:E:183:LEU:HD22	1:E:194:TYR:CZ	2.43	0.53
2:U:117:LYS:HE3	2:U:144:GLY:O	2.08	0.53
1:B:61:TYR:HD2	1:B:85:LEU:HD12	1.72	0.53
1:C:134:THR:HG22	3:P:31:LYS:HG2	1.90	0.53
2:S:34:TYR:H	2:S:95:ASP:HB2	1.72	0.53
2:U:197:HIS:HB3	2:U:202:THR:HB	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:27:PHE:CE2	2:S:94:ARG:HD3	2.43	0.53
1:C:118:THR:HG22	1:C:121:ARG:HH22	1.73	0.53
1:C:134:THR:HA	3:P:31:LYS:HE2	1.90	0.53
1:C:225:SER:OG	1:C:342:LYS:NZ	2.39	0.53
1:C:55:LEU:HD21	1:C:378:THR:OG1	2.09	0.53
3:N:105:GLU:OE2	3:N:171:TYR:OH	2.26	0.53
2:O:40:ALA:HB3	2:O:43:LYS:HE2	1.90	0.53
1:A:277:THR:HB	1:A:325:THR:HG21	1.91	0.53
1:E:213:PHE:HB3	1:E:240:SER:HB3	1.90	0.53
1:F:128:VAL:HG21	1:F:162:MET:HE1	1.90	0.52
2:M:119:PRO:HB2	2:M:142:VAL:HG13	1.90	0.52
3:P:33:LEU:HD22	3:P:71:PHE:CG	2.44	0.52
2:Q:119:PRO:HB2	2:Q:142:VAL:HG13	1.91	0.52
2:Q:138:LEU:HD11	2:Q:210:PRO:HA	1.89	0.52
1:A:356:VAL:HG11	1:A:378:THR:HG22	1.92	0.52
3:P:112:PRO:HB3	3:P:138:TYR:HB3	1.91	0.52
2:S:96:LYS:HD2	2:S:100(E):TYR:OH	2.10	0.52
3:P:94:TYR:HA	3:P:95(B):LEU:HD12	1.91	0.52
3:V:31:LYS:O	3:V:51:SER:N	2.43	0.52
1:F:31:CYS:HB3	1:F:359:LEU:HG	1.92	0.52
2:U:154:TRP:CZ3	2:U:193:CYS:HB3	2.44	0.52
2:M:3:SER:HB3	2:M:25:SER:HB2	1.91	0.52
1:B:134:THR:HA	3:N:31:LYS:HE2	1.92	0.52
2:Q:94:ARG:HE	2:Q:102:LEU:HD12	1.75	0.52
1:E:130:PRO:O	1:E:134:THR:HG23	2.10	0.52
2:M:194:ASN:ND2	2:M:205:ASP:OD1	2.43	0.52
2:H:103:TRP:CE3	3:L:44:PRO:HD2	2.45	0.52
5:X:1:NAG:H3	5:X:1:NAG:C8	2.38	0.51
1:D:290:GLY:HA3	1:D:303:GLU:OE1	2.11	0.51
2:Q:103:TRP:CE3	3:R:44:PRO:HD2	2.45	0.51
1:E:76:TRP:CZ2	1:E:77:ASN:ND2	2.79	0.51
2:U:154:TRP:HB3	2:U:159:LEU:HD23	1.93	0.51
1:B:134:THR:HG22	3:N:31:LYS:HG2	1.93	0.51
1:F:130:PRO:O	1:F:134:THR:HG23	2.11	0.51
1:F:362:ASP:OD2	1:F:378:THR:HG23	2.10	0.51
2:M:4:VAL:HG13	2:M:22:CYS:SG	2.51	0.51
1:C:356:VAL:HG11	1:C:378:THR:HG22	1.91	0.51
3:P:33:LEU:HD22	3:P:71:PHE:CB	2.41	0.51
2:H:127:CYS:HB3	3:L:118:PRO:HG2	1.93	0.51
3:T:33:LEU:HD22	3:T:71:PHE:CB	2.41	0.51
1:A:66:ILE:HG21	1:A:114:ILE:HD13	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:66:ARG:NH2	2:M:86:ASP:OD2	2.44	0.51
2:S:4:VAL:HG13	2:S:22:CYS:SG	2.51	0.51
1:E:340:LYS:HD3	1:E:384:ALA:HB2	1.93	0.51
1:D:146:TYR:HE1	1:D:209:MET:HE1	1.75	0.51
2:M:122:PHE:CD2	3:N:123:GLN:HG3	2.46	0.51
2:H:96:LYS:HD2	2:H:100(E):TYR:HE1	1.76	0.51
3:P:80:CYS:SG	3:P:169:CYS:N	2.83	0.50
2:Q:40:ALA:HB3	2:Q:43:LYS:HE2	1.93	0.50
2:Q:168:SER:HB3	2:Q:178:LEU:HD23	1.92	0.50
2:S:12:VAL:HG21	2:S:82(B):LEU:HD13	1.94	0.50
2:U:96:LYS:HE3	2:U:101:GLN:OE1	2.11	0.50
1:E:147:PRO:HB2	1:E:193:SER:HB3	1.92	0.50
2:M:95:ASP:HB3	2:M:97:GLY:H	1.76	0.50
1:E:67:SER:HA	4:W:1:NAG:H62	1.93	0.50
1:F:290:GLY:HA3	1:F:303:GLU:OE1	2.11	0.50
2:O:213:CYS:HB2	3:R:185:ASN:O	2.11	0.50
2:S:164:ARG:NH2	3:T:136:ASN:OD1	2.35	0.50
6:J:1:NAG:H3	6:J:1:NAG:H83	1.93	0.50
1:D:134:THR:HA	3:R:31:LYS:HE2	1.93	0.50
1:E:55:LEU:HD21	1:E:378:THR:OG1	2.12	0.50
3:P:141:ASP:OD2	3:P:161:LYS:NZ	2.45	0.50
3:R:54:LEU:HD11	3:R:60:SER:HA	1.94	0.50
3:R:12:GLU:HG2	3:R:105:GLU:HB3	1.93	0.50
3:T:147:GLU:HB2	3:T:191:THR:HB	1.94	0.50
1:D:134:THR:HG22	3:R:31:LYS:HG2	1.92	0.50
3:T:89:GLN:HB2	3:T:98:PHE:CD1	2.47	0.50
1:C:217:TRP:CZ3	1:C:218:ARG:HG2	2.47	0.50
1:E:118:THR:HG22	1:E:121:ARG:HH22	1.77	0.50
1:C:249:MET:HE3	1:C:259:LEU:HD21	1.94	0.49
3:T:90:GLN:NE2	3:T:95(B):LEU:HD11	2.27	0.49
1:A:51:LEU:HD22	1:A:359:LEU:HD13	1.94	0.49
2:U:99:SER:HG	2:U:100(B):TRP:CD1	2.31	0.49
2:H:4:VAL:HG13	2:H:22:CYS:SG	2.52	0.49
1:E:58:HIS:HA	1:E:96:LYS:O	2.12	0.49
3:N:12:GLU:HG2	3:N:105:GLU:HB3	1.94	0.49
3:P:105:GLU:OE2	3:P:171:TYR:OH	2.30	0.49
3:V:90:GLN:NE2	3:V:95(B):LEU:HD11	2.26	0.49
1:A:362:ASP:OD2	1:A:378:THR:HG23	2.12	0.49
3:L:89:GLN:HB2	3:L:98:PHE:CD1	2.48	0.49
3:R:147:GLU:HB2	3:R:191:THR:HB	1.94	0.49
1:D:146:TYR:CE1	1:D:209:MET:HE1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:35:MET:HG2	2:M:77:VAL:HG21	1.95	0.49
2:O:29:ILE:O	2:O:29:ILE:HD12	2.12	0.49
1:D:122:ARG:NH2	1:D:126:LYS:HD3	2.28	0.49
1:F:61:TYR:HD2	1:F:85:LEU:HD12	1.76	0.49
2:M:31:SER:HB2	2:M:52(B):GLY:HA2	1.95	0.49
2:O:51:ILE:HG13	2:O:57:THR:HG22	1.93	0.49
3:P:54:LEU:HD11	3:P:60:SER:HA	1.95	0.49
2:Q:17:SER:HB2	2:Q:81:MET:O	2.13	0.49
1:E:277:THR:HB	1:E:325:THR:HG21	1.94	0.48
1:C:134:THR:HG21	3:P:31:LYS:H	1.79	0.48
3:V:141:ASP:OD2	3:V:161:LYS:NZ	2.43	0.48
1:C:126:LYS:HD2	3:P:27:GLN:OE1	2.13	0.48
3:V:106(A):LEU:HA	3:V:139:PHE:CZ	2.48	0.48
1:C:311:ALA:HB2	1:C:326:LYS:HB2	1.95	0.48
3:N:13:VAL:HG11	3:N:78:VAL:HG21	1.95	0.48
3:V:49:TYR:O	3:V:53:SER:HB2	2.12	0.48
1:F:134:THR:HG21	3:V:31:LYS:H	1.78	0.48
2:Q:130:ASP:N	2:Q:130:ASP:OD1	2.46	0.48
3:R:144:VAL:HG22	3:R:194:VAL:HG22	1.94	0.48
2:S:11:LEU:HB2	2:S:147:PRO:HG3	1.94	0.48
1:A:346:LEU:HB2	1:A:351:LEU:HD12	1.95	0.48
1:E:299:LEU:HD12	1:E:331:VAL:HG21	1.95	0.48
3:N:190:TYR:HB2	3:N:205:PHE:CE1	2.49	0.48
3:T:137:LYS:NZ	3:T:168:ASP:OD2	2.33	0.48
2:M:168:SER:HA	2:M:178:LEU:HB3	1.95	0.48
2:H:35:MET:HG2	2:H:77:VAL:HG21	1.96	0.48
1:D:66:ILE:HG21	1:D:114:ILE:HD13	1.95	0.48
1:D:304:ILE:O	1:D:308:LEU:HG	2.14	0.48
2:U:127:CYS:HB3	3:V:118:PRO:HG2	1.96	0.48
1:A:174:LYS:HG2	2:H:98:TRP:CD1	2.49	0.47
1:D:149:ARG:HA	1:D:192:SER:O	2.13	0.47
2:H:197:HIS:HB3	2:H:202:THR:HB	1.95	0.47
2:O:38:ARG:HB3	2:O:48:ILE:HD11	1.96	0.47
3:T:139:PHE:CD1	3:T:140:PRO:HA	2.49	0.47
1:C:61:TYR:HD2	1:C:85:LEU:HD12	1.79	0.47
2:O:3:SER:HB3	2:O:25:SER:HB2	1.95	0.47
1:A:118:THR:HG22	1:A:121:ARG:NH2	2.29	0.47
1:D:58:HIS:HA	1:D:96:LYS:O	2.14	0.47
1:E:134:THR:HG21	3:T:31:LYS:H	1.80	0.47
2:H:66:ARG:HH22	2:H:86:ASP:CG	2.22	0.47
3:V:59:SER:OG	3:V:61:ARG:HG3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:V:148:VAL:HG22	3:V:190:TYR:HD1	1.79	0.47
2:H:96:LYS:HD2	2:H:100(E):TYR:CE1	2.49	0.47
3:L:80:CYS:SG	3:L:169:CYS:N	2.87	0.47
1:B:208:ILE:HD11	1:B:245:ALA:HB1	1.96	0.47
1:F:174:LYS:HG2	2:U:98:TRP:CG	2.49	0.47
2:O:81:MET:HE1	2:O:109:VAL:HG21	1.97	0.47
2:Q:35:MET:HG2	2:Q:77:VAL:HG21	1.96	0.47
3:R:49:TYR:O	3:R:53:SER:HB2	2.15	0.47
1:C:146:TYR:CE1	1:C:184:SER:HB2	2.50	0.47
1:E:76:TRP:CE2	1:E:77:ASN:ND2	2.82	0.47
1:E:356:VAL:HG11	1:E:378:THR:HG22	1.96	0.47
1:F:77:ASN:HB2	1:F:81:LEU:HD13	1.97	0.47
2:O:103:TRP:CE3	3:P:44:PRO:HD2	2.50	0.47
3:T:62:PHE:CD1	3:T:75:ILE:HG12	2.50	0.47
1:E:269:SER:HB3	1:E:331:VAL:HG22	1.97	0.47
2:M:11:LEU:HB2	2:M:147:PRO:HG3	1.96	0.47
2:S:33:TYR:CE2	2:S:52(A):ILE:HG13	2.50	0.47
1:B:142:LEU:HD12	1:B:159:ILE:HD13	1.97	0.46
1:E:30:VAL:O	1:E:355:MET:HA	2.15	0.46
1:E:159:ILE:HD12	1:E:203:LEU:HD21	1.97	0.46
2:H:33:TYR:CE2	2:H:52(A):ILE:HG13	2.48	0.46
3:L:106(A):LEU:HA	3:L:139:PHE:CZ	2.50	0.46
3:P:66:GLY:HA3	3:P:71:PHE:HA	1.97	0.46
5:X:1:NAG:H83	5:X:1:NAG:C3	2.44	0.46
1:F:171:GLN:OE1	3:V:93:SER:HB2	2.15	0.46
3:V:33:LEU:HG	3:V:34:ASN:N	2.30	0.46
1:A:55:LEU:HD21	1:A:378:THR:OG1	2.15	0.46
1:C:315:ARG:NH1	1:C:335:ASP:OD2	2.40	0.46
1:B:210:THR:HB	1:B:242:THR:HG22	1.97	0.46
1:F:125:ILE:HA	1:F:162:MET:HE2	1.97	0.46
1:F:259:LEU:HD23	1:F:259:LEU:HA	1.71	0.46
3:N:90:GLN:NE2	3:N:95(B):LEU:HD11	2.31	0.46
2:M:66:ARG:HH22	2:M:86:ASP:CG	2.24	0.46
3:N:129:VAL:HG21	3:N:184:TYR:HB2	1.96	0.46
1:A:145:LEU:HD22	1:A:209:MET:HE1	1.98	0.46
1:C:52:ASP:HB3	1:C:55:LEU:HB2	1.97	0.46
1:F:134:THR:CG2	3:V:31:LYS:H	2.28	0.46
3:P:13:VAL:HG11	3:P:78:VAL:HG21	1.98	0.46
2:S:2:PCA:C	2:S:102:LEU:HD11	2.41	0.46
3:V:181:SER:O	3:V:185:ASN:ND2	2.49	0.46
1:B:134:THR:HG21	3:N:31:LYS:H	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:299:LEU:HD12	1:F:331:VAL:HG21	1.98	0.45
1:A:269:SER:HB3	1:A:331:VAL:HG22	1.98	0.45
1:B:33:TYR:CE2	1:B:48:PRO:HG3	2.51	0.45
1:D:226:PRO:O	1:D:242:THR:HG23	2.16	0.45
2:Q:52:TYR:C	2:Q:52(B):GLY:H	2.25	0.45
2:S:51:ILE:HG13	2:S:57:THR:HG22	1.97	0.45
3:N:141:ASP:OD2	3:N:161:LYS:NZ	2.42	0.45
1:A:76:TRP:H	1:A:76:TRP:CD1	2.33	0.45
1:C:198:LYS:O	1:C:201:GLN:HG2	2.16	0.45
1:A:112:SER:HB2	1:A:151:ASP:OD2	2.17	0.45
1:C:174:LYS:HG2	2:O:98:TRP:CD1	2.51	0.45
1:D:340:LYS:HD3	1:D:384:ALA:HB2	1.98	0.45
1:A:259:LEU:HA	1:A:259:LEU:HD23	1.68	0.45
1:F:29:LEU:O	1:F:56:CYS:HB3	2.17	0.45
1:F:96:LYS:HA	1:F:138:ASP:OD2	2.17	0.45
2:M:22:CYS:HB3	2:M:77:VAL:HG22	1.99	0.45
2:H:31:SER:O	2:H:33:TYR:N	2.50	0.45
2:S:127:CYS:HB3	3:T:118:PRO:HG2	1.99	0.45
1:F:163:LYS:O	1:F:167:ILE:HG13	2.17	0.45
2:O:194:ASN:HB3	2:O:203:LYS:HE3	1.99	0.45
1:F:146:TYR:CE1	1:F:184:SER:HB2	2.52	0.45
1:F:183:LEU:HB3	1:F:194:TYR:CE2	2.52	0.45
2:M:122:PHE:CE2	3:N:123:GLN:HG3	2.52	0.44
2:O:119:PRO:HB2	2:O:142:VAL:HG13	1.99	0.44
3:P:129:VAL:HG21	3:P:184:TYR:HB2	1.99	0.44
1:D:271:THR:HG23	1:D:328:ASN:OD1	2.17	0.44
3:L:147:GLU:HB2	3:L:191:THR:HB	1.98	0.44
2:Q:4:VAL:HG13	2:Q:22:CYS:SG	2.57	0.44
2:H:123:PRO:HB3	2:H:208:VAL:HG22	1.99	0.44
3:P:40:PRO:HG2	3:P:163:PRO:HB3	1.98	0.44
2:Q:27:PHE:CE2	2:Q:94:ARG:HD3	2.52	0.44
2:S:103:TRP:CE3	3:T:44:PRO:HD2	2.53	0.44
2:S:122:PHE:CD2	3:T:123:GLN:HG3	2.53	0.44
2:U:100:ASN:O	3:V:32:TYR:OH	2.23	0.44
1:D:148:GLY:H	1:D:151:ASP:HB2	1.81	0.44
2:H:73:SER:HB3	2:H:76:THR:OG1	2.18	0.44
2:H:165:THR:HG23	2:H:180:SER:HB2	1.99	0.44
3:R:34:ASN:OD1	3:R:49:TYR:HA	2.17	0.44
1:A:128:VAL:CG2	1:A:162:MET:HE1	2.41	0.44
1:B:101:VAL:HG23	1:B:140:LEU:HD11	1.98	0.44
1:B:128:VAL:CG2	1:B:162:MET:HE1	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:75:GLU:HG3	1:E:76:TRP:H	1.83	0.44
1:E:265:THR:HG22	1:E:377:LEU:HD12	1.99	0.44
2:O:130:ASP:N	2:O:130:ASP:OD1	2.50	0.44
1:A:101:VAL:HG23	1:A:140:LEU:HD11	1.99	0.44
1:F:271:THR:HG23	1:F:328:ASN:OD1	2.18	0.44
3:P:36:TYR:CD1	3:P:46:LEU:HA	2.52	0.44
3:P:49:TYR:O	3:P:53:SER:HB2	2.17	0.44
1:D:134:THR:CG2	3:R:31:LYS:H	2.31	0.44
2:O:11:LEU:HB2	2:O:147:PRO:HG3	2.00	0.44
3:P:148:VAL:HG22	3:P:190:TYR:CD1	2.53	0.44
3:R:13:VAL:HG11	3:R:78:VAL:HG21	2.00	0.44
1:A:322:PRO:HD3	1:A:338:SER:HB3	1.99	0.44
1:B:208:ILE:CD1	1:B:249:MET:HE2	2.43	0.44
3:N:106(A):LEU:HA	3:N:139:PHE:CZ	2.53	0.44
2:O:66:ARG:HH22	2:O:86:ASP:CG	2.25	0.44
3:T:6:GLN:NE2	3:T:86:TYR:O	2.51	0.44
1:C:269:SER:HB3	1:C:331:VAL:HG22	1.98	0.44
1:D:100:SER:HA	1:D:141:ASP:O	2.18	0.44
1:D:144:TRP:O	1:D:147:PRO:HD3	2.18	0.44
2:H:130:ASP:OD1	2:H:130:ASP:N	2.50	0.44
2:O:53:SER:C	2:O:55:THR:N	2.76	0.44
3:R:137:LYS:NZ	3:R:168:ASP:OD2	2.36	0.44
1:A:200:SER:HB3	1:A:255:PRO:HD3	2.00	0.43
1:B:66:ILE:HG21	1:B:114:ILE:HD13	2.00	0.43
2:U:145:TYR:CE2	2:U:150:VAL:HG13	2.53	0.43
1:B:33:TYR:OH	1:B:48:PRO:HD3	2.18	0.43
1:D:55:LEU:HD21	1:D:378:THR:OG1	2.18	0.43
2:M:2:PCA:HB3	2:M:102:LEU:HD11	1.99	0.43
3:P:148:VAL:HG22	3:P:190:TYR:HD1	1.83	0.43
3:R:104:LEU:HD23	3:R:104:LEU:HA	1.91	0.43
2:U:197:HIS:CE1	2:U:199:ALA:HB3	2.53	0.43
1:E:362:ASP:OD1	1:E:363:ASP:N	2.50	0.43
1:D:311:ALA:HB2	1:D:326:LYS:HB2	2.00	0.43
2:M:125:ALA:O	3:N:118:PRO:HD2	2.17	0.43
2:M:135:THR:OG1	2:M:185:THR:HG22	2.18	0.43
3:N:40:PRO:HG2	3:N:163:PRO:HB3	1.98	0.43
2:S:119:PRO:HB2	2:S:142:VAL:HG13	1.99	0.43
2:M:197:HIS:CE1	2:M:199:ALA:HB3	2.53	0.43
2:M:197:HIS:HB3	2:M:202:THR:HB	1.99	0.43
3:L:141:ASP:OD2	3:L:161:LYS:NZ	2.39	0.43
1:B:146:TYR:CE1	1:B:184:SER:HB2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:31:CYS:HB3	1:C:359:LEU:HG	2.01	0.43
1:E:76:TRP:CE2	3:R:61:ARG:HG2	2.53	0.43
2:S:197:HIS:HB3	2:S:202:THR:HB	2.01	0.43
3:V:129:VAL:HG21	3:V:184:TYR:HB2	1.99	0.43
1:A:168:LYS:HD3	3:L:95(A):ASP:OD2	2.19	0.43
1:A:208:ILE:CD1	1:A:249:MET:HE2	2.48	0.43
1:B:163:LYS:HA	1:B:163:LYS:HD3	1.91	0.43
3:N:206:ASN:HB2	3:N:209:ASP:CG	2.44	0.43
2:M:31:SER:O	2:M:33:TYR:N	2.52	0.43
2:H:36:TRP:CE2	2:H:79:LEU:HB2	2.54	0.43
3:P:185:ASN:HB3	2:Q:213:CYS:C	2.44	0.43
2:U:35:MET:HG2	2:U:77:VAL:HG21	2.01	0.43
1:B:226:PRO:O	1:B:242:THR:HG23	2.18	0.43
1:D:31:CYS:HB3	1:D:359:LEU:HG	2.01	0.43
1:E:208:ILE:HG22	1:E:260:VAL:O	2.18	0.43
1:F:72:ASP:OD1	1:F:73:THR:N	2.51	0.43
3:R:33:LEU:HD22	3:R:71:PHE:CB	2.49	0.43
2:U:122:PHE:CD2	3:V:123:GLN:HG3	2.54	0.43
1:A:288:ILE:HG12	1:A:329:GLN:NE2	2.33	0.43
1:C:48:PRO:O	1:C:51:LEU:HG	2.19	0.43
1:C:128:VAL:HB	1:C:129:PRO:HD3	2.01	0.43
1:C:226:PRO:O	1:C:242:THR:HG23	2.18	0.43
5:X:1:NAG:C8	5:X:1:NAG:C3	2.97	0.43
1:B:356:VAL:HG11	1:B:378:THR:HG22	2.00	0.42
3:T:13:VAL:HG11	3:T:78:VAL:HG21	2.01	0.42
1:C:259:LEU:O	1:C:352:ALA:HB3	2.19	0.42
2:O:29:ILE:HD12	2:O:29:ILE:C	2.45	0.42
2:S:94:ARG:HG2	2:S:95:ASP:O	2.19	0.42
3:V:54:LEU:HD23	3:V:54:LEU:HA	1.91	0.42
1:B:100:SER:HA	1:B:141:ASP:O	2.19	0.42
3:N:33:LEU:HD22	3:N:71:PHE:CB	2.49	0.42
3:R:54:LEU:HD23	3:R:54:LEU:HA	1.79	0.42
3:T:62:PHE:HD1	3:T:75:ILE:HG12	1.83	0.42
3:T:139:PHE:CG	3:T:140:PRO:HA	2.54	0.42
2:U:125:ALA:O	3:V:118:PRO:HD2	2.19	0.42
1:B:125:ILE:HA	1:B:162:MET:HE2	2.00	0.42
1:C:159:ILE:CD1	1:C:203:LEU:HD21	2.49	0.42
2:O:53:SER:C	2:O:55:THR:H	2.28	0.42
1:D:53:ARG:CZ	1:D:91:ARG:HD2	2.49	0.42
1:D:198:LYS:O	1:D:201:GLN:HG2	2.19	0.42
1:D:269:SER:HB3	1:D:331:VAL:HG22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:122:PHE:CD2	3:P:123:GLN:HG3	2.55	0.42
2:Q:12:VAL:HG11	2:Q:18:LEU:HB2	2.02	0.42
2:Q:125:ALA:O	3:R:118:PRO:HD2	2.19	0.42
2:U:51:ILE:HD13	2:U:71:LYS:HG3	2.01	0.42
3:V:158:GLU:HB3	3:V:176:THR:OG1	2.20	0.42
1:D:208:ILE:HD11	1:D:245:ALA:HB1	2.01	0.42
1:E:31:CYS:HB3	1:E:359:LEU:HG	2.02	0.42
1:F:89:LYS:HD2	1:F:97:THR:HG23	2.02	0.42
3:R:31:LYS:O	3:R:51:SER:N	2.53	0.42
1:B:363:ASP:OD2	1:B:369:CYS:HB2	2.20	0.42
1:C:66:ILE:HG21	1:C:114:ILE:HD13	2.02	0.42
1:F:183:LEU:HD22	1:F:194:TYR:CZ	2.55	0.42
2:H:122:PHE:CD2	3:L:123:GLN:HG3	2.55	0.42
1:A:100:SER:HA	1:A:141:ASP:O	2.19	0.42
1:E:259:LEU:HD23	1:E:259:LEU:HA	1.81	0.42
1:F:108:SER:HA	1:F:144:TRP:HE1	1.85	0.42
1:F:208:ILE:HB	1:F:249:MET:HE2	2.02	0.42
3:T:190:TYR:HB2	3:T:205:PHE:CE1	2.54	0.42
2:U:73:SER:OG	2:U:74:SER:N	2.53	0.42
1:F:100:SER:HA	1:F:141:ASP:O	2.19	0.41
2:H:17:SER:HB2	2:H:81:MET:O	2.20	0.41
3:L:129:VAL:HG21	3:L:184:TYR:HB2	2.02	0.41
3:P:104:LEU:HA	3:P:104:LEU:HD23	1.74	0.41
2:Q:31:SER:OG	2:Q:52(A):ILE:O	2.34	0.41
3:V:184:TYR:O	3:V:190:TYR:OH	2.38	0.41
1:A:339:VAL:CG1	1:A:381:ILE:HG13	2.50	0.41
1:D:85:LEU:C	1:D:85:LEU:HD23	2.45	0.41
1:D:171:GLN:OE1	3:R:93:SER:HB2	2.20	0.41
1:E:304:ILE:O	1:E:308:LEU:HG	2.20	0.41
3:L:33:LEU:HD22	3:L:71:PHE:CG	2.55	0.41
3:T:19:VAL:O	3:T:74:THR:HA	2.20	0.41
1:A:208:ILE:CG1	1:A:249:MET:HE2	2.50	0.41
1:B:144:TRP:O	1:B:147:PRO:HD3	2.19	0.41
1:D:73:THR:OG1	1:D:78:ASP:HB2	2.21	0.41
2:H:184:VAL:HB	2:H:188:SER:HB2	2.02	0.41
2:Q:164:ARG:NH2	3:R:136:ASN:OD1	2.43	0.41
3:R:80:CYS:SG	3:R:169:CYS:N	2.94	0.41
3:V:33:LEU:HD22	3:V:71:PHE:CB	2.50	0.41
3:V:148:VAL:HG22	3:V:190:TYR:CD1	2.54	0.41
1:A:122:ARG:NH2	1:A:126:LYS:HD3	2.35	0.41
2:M:40:ALA:HB3	2:M:43:LYS:HE2	2.00	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:33:LEU:HD22	3:L:71:PHE:CB	2.51	0.41
2:Q:66:ARG:NH2	2:Q:86:ASP:OD2	2.53	0.41
1:D:163:LYS:HA	1:D:163:LYS:HD3	1.89	0.41
1:D:356:VAL:HG11	1:D:378:THR:HG22	2.03	0.41
1:E:357:TRP:HA	1:E:358:ALA:HA	1.87	0.41
3:N:122:ASP:OD1	3:N:122:ASP:N	2.52	0.41
2:O:213:CYS:CB	3:R:185:ASN:HB3	2.50	0.41
2:S:184:VAL:HB	2:S:188:SER:HB2	2.02	0.41
3:T:31:LYS:O	3:T:51:SER:N	2.54	0.41
1:D:52:ASP:HB3	1:D:55:LEU:HB2	2.02	0.41
1:E:52:ASP:HB3	1:E:55:LEU:HB2	2.03	0.41
1:A:115:ALA:HB2	1:A:158:LEU:HD12	2.02	0.41
1:A:128:VAL:HB	1:A:129:PRO:HD3	2.03	0.41
1:E:134:THR:CG2	3:T:31:LYS:H	2.33	0.41
2:Q:122:PHE:CE2	3:R:123:GLN:HG3	2.56	0.41
1:E:66:ILE:HD11	1:E:101:VAL:HG11	2.03	0.41
3:N:104:LEU:HD23	3:N:104:LEU:HA	1.78	0.41
2:H:11:LEU:HB2	2:H:147:PRO:HG3	2.02	0.41
2:H:36:TRP:O	2:H:48:ILE:HB	2.20	0.41
3:P:164:GLN:HG2	3:P:171:TYR:CE1	2.56	0.41
2:U:52:TYR:CE2	2:U:52(B):GLY:HA3	2.56	0.41
1:B:146:TYR:CE1	1:B:209:MET:HE1	2.56	0.41
1:C:61:TYR:OH	1:C:78:ASP:OD1	2.39	0.41
1:D:118:THR:HG22	1:D:121:ARG:NH2	2.34	0.41
1:E:208:ILE:CG1	1:E:249:MET:HE2	2.51	0.41
3:L:31:LYS:O	3:L:51:SER:N	2.54	0.41
3:R:164:GLN:HG2	3:R:171:TYR:CE1	2.55	0.41
3:V:112:PRO:HB3	3:V:138:TYR:HB3	2.03	0.41
1:E:128:VAL:CG2	1:E:162:MET:HE1	2.49	0.41
1:F:356:VAL:HG11	1:F:378:THR:HG22	2.04	0.41
2:M:36:TRP:CE2	2:M:79:LEU:HB2	2.56	0.41
2:H:53:SER:HB2	2:H:56:ASP:OD1	2.21	0.41
2:Q:98:TRP:CG	2:Q:99:SER:N	2.89	0.41
3:V:48:ILE:HA	3:V:53:SER:O	2.21	0.41
1:A:101:VAL:CG2	1:A:140:LEU:HD11	2.51	0.40
1:E:142:LEU:HD12	1:E:159:ILE:HD13	2.02	0.40
1:E:227:LEU:HD13	1:E:342:LYS:HG2	2.02	0.40
2:H:34:TYR:HB2	2:H:95:ASP:OD2	2.20	0.40
1:A:299:LEU:HD12	1:A:331:VAL:HG21	2.03	0.40
1:C:259:LEU:HA	1:C:259:LEU:HD23	1.70	0.40
1:D:115:ALA:HB2	1:D:158:LEU:HD12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:125:ILE:HA	1:E:162:MET:HE2	2.03	0.40
1:E:163:LYS:HA	1:E:163:LYS:HD3	1.86	0.40
1:F:259:LEU:O	1:F:352:ALA:HB3	2.20	0.40
2:M:130:ASP:N	2:M:130:ASP:OD1	2.53	0.40
2:M:138:LEU:HD21	2:M:210:PRO:HA	2.03	0.40
3:T:148:VAL:HG22	3:T:190:TYR:CD1	2.56	0.40
1:B:76:TRP:HH2	3:V:60:SER:HB2	1.87	0.40
1:D:108:SER:HB2	1:D:144:TRP:HE1	1.85	0.40
1:F:195:ASP:OD2	1:F:198:LYS:HD3	2.22	0.40
3:N:62:PHE:CD1	3:N:75:ILE:HG12	2.57	0.40
2:H:95:ASP:HB3	2:H:97:GLY:H	1.87	0.40
2:H:99:SER:HB3	2:H:100(C):GLY:H	1.87	0.40
3:L:165:ASN:HB3	3:L:168:ASP:OD1	2.20	0.40
2:U:119:PRO:HB2	2:U:142:VAL:HG13	2.04	0.40
1:A:174:LYS:HG2	2:H:98:TRP:CG	2.57	0.40
3:L:61:ARG:HH22	3:L:82:ASP:CG	2.29	0.40
3:R:21:ILE:HD13	3:R:21:ILE:HG21	1.76	0.40
2:S:34:TYR:HB2	2:S:95:ASP:OD2	2.22	0.40
2:U:159:LEU:HD21	2:U:182:VAL:HG21	2.03	0.40
3:V:33:LEU:HD22	3:V:71:PHE:CG	2.56	0.40
1:A:191:ASP:OD1	1:A:248:TYR:OH	2.36	0.40
1:E:305:CYS:CB	1:E:369:CYS:HA	2.52	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:349:ARG:NH2	1:E:233:ASP:OD2[1_655]	2.01	0.19
1:B:233:ASP:OD2	1:E:349:ARG:NH2[1_655]	2.14	0.06

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	360/362 (99%)	352 (98%)	8 (2%)	0	100	100
1	B	360/362 (99%)	351 (98%)	9 (2%)	0	100	100
1	C	360/362 (99%)	351 (98%)	9 (2%)	0	100	100
1	D	360/362 (99%)	353 (98%)	7 (2%)	0	100	100
1	E	360/362 (99%)	352 (98%)	8 (2%)	0	100	100
1	F	360/362 (99%)	350 (97%)	10 (3%)	0	100	100
2	H	221/223 (99%)	211 (96%)	9 (4%)	1 (0%)	24	55
2	M	221/223 (99%)	210 (95%)	10 (4%)	1 (0%)	24	55
2	O	221/223 (99%)	210 (95%)	10 (4%)	1 (0%)	24	55
2	Q	221/223 (99%)	209 (95%)	11 (5%)	1 (0%)	24	55
2	S	221/223 (99%)	209 (95%)	10 (4%)	2 (1%)	14	43
2	U	221/223 (99%)	211 (96%)	10 (4%)	0	100	100
3	L	212/214 (99%)	207 (98%)	4 (2%)	1 (0%)	24	55
3	N	212/214 (99%)	205 (97%)	7 (3%)	0	100	100
3	P	212/214 (99%)	206 (97%)	6 (3%)	0	100	100
3	R	212/214 (99%)	205 (97%)	6 (3%)	1 (0%)	24	55
3	T	212/214 (99%)	206 (97%)	6 (3%)	0	100	100
3	V	212/214 (99%)	205 (97%)	7 (3%)	0	100	100
All	All	4758/4794 (99%)	4603 (97%)	147 (3%)	8 (0%)	43	71

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	M	32	ASP
2	H	32	ASP
2	O	54	GLY
2	S	32	ASP
2	Q	52(A)	ILE
3	L	50	SER
3	R	50	SER
2	S	52(B)	GLY

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	302/302 (100%)	301 (100%)	1 (0%)	86	86
1	B	302/302 (100%)	302 (100%)	0	100	100
1	C	302/302 (100%)	302 (100%)	0	100	100
1	D	302/302 (100%)	302 (100%)	0	100	100
1	E	302/302 (100%)	301 (100%)	1 (0%)	86	86
1	F	302/302 (100%)	302 (100%)	0	100	100
2	H	186/186 (100%)	185 (100%)	1 (0%)	81	83
2	M	186/186 (100%)	185 (100%)	1 (0%)	81	83
2	O	186/186 (100%)	186 (100%)	0	100	100
2	Q	186/186 (100%)	185 (100%)	1 (0%)	81	83
2	S	186/186 (100%)	185 (100%)	1 (0%)	81	83
2	U	186/186 (100%)	185 (100%)	1 (0%)	81	83
3	L	182/182 (100%)	182 (100%)	0	100	100
3	N	182/182 (100%)	182 (100%)	0	100	100
3	P	182/182 (100%)	182 (100%)	0	100	100
3	R	182/182 (100%)	182 (100%)	0	100	100
3	T	182/182 (100%)	182 (100%)	0	100	100
3	V	182/182 (100%)	182 (100%)	0	100	100
All	All	4020/4020 (100%)	4013 (100%)	7 (0%)	87	88

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	369	CYS
1	E	369	CYS
2	M	127	CYS
2	H	127	CYS
2	Q	127	CYS
2	S	127	CYS
2	U	127	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	ASN
1	A	319	GLN
1	B	117	ASN
1	C	319	GLN
1	D	153	GLN
1	D	314	HIS
1	D	319	GLN
1	E	38	GLN
1	E	77	ASN
1	E	117	ASN
1	E	153	GLN
1	E	319	GLN
1	E	379	ASN
1	F	38	GLN
1	F	314	HIS
1	F	344	GLN
1	F	379	ASN
2	M	189	GLN
3	N	79	GLN
3	N	153	GLN
3	N	172	ASN
2	H	39	GLN
3	L	38	GLN
3	L	153	GLN
3	L	172	ASN
2	O	39	GLN
3	P	38	GLN
3	P	172	ASN
3	R	153	GLN
3	R	187	HIS
3	T	206	ASN
3	V	27	GLN
3	V	159	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues 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	PCA	M	2	2	7,8,9	1.96	1 (14%)	9,10,12	2.69	5 (55%)
2	PCA	U	2	2	7,8,9	1.85	1 (14%)	9,10,12	2.23	5 (55%)
2	PCA	H	2	2	7,8,9	1.86	1 (14%)	9,10,12	2.30	5 (55%)
2	PCA	O	2	2	7,8,9	1.95	1 (14%)	9,10,12	2.54	5 (55%)
2	PCA	Q	2	2	7,8,9	1.90	1 (14%)	9,10,12	2.00	4 (44%)
2	PCA	S	2	2	7,8,9	1.90	1 (14%)	9,10,12	2.46	5 (55%)

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	PCA	M	2	2	-	0/0/11/13	0/1/1/1
2	PCA	U	2	2	-	0/0/11/13	0/1/1/1
2	PCA	H	2	2	-	0/0/11/13	0/1/1/1
2	PCA	O	2	2	-	0/0/11/13	0/1/1/1
2	PCA	Q	2	2	-	0/0/11/13	0/1/1/1
2	PCA	S	2	2	-	0/0/11/13	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	2	PCA	CD-N	5.03	1.47	1.34
2	O	2	PCA	CD-N	4.94	1.46	1.34
2	Q	2	PCA	CD-N	4.92	1.46	1.34
2	S	2	PCA	CD-N	4.87	1.46	1.34
2	H	2	PCA	CD-N	4.74	1.46	1.34
2	U	2	PCA	CD-N	4.73	1.46	1.34

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	2	PCA	CB-CA-C	-4.42	106.59	112.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	2	PCA	OE-CD-CG	-3.82	119.91	126.72
2	O	2	PCA	OE-CD-CG	-3.57	120.35	126.72
2	O	2	PCA	CB-CA-C	-3.47	107.89	112.66
2	S	2	PCA	CB-CA-C	-3.31	108.12	112.66
2	H	2	PCA	OE-CD-CG	-3.25	120.92	126.72
2	S	2	PCA	CB-CA-N	3.16	111.94	103.24
2	S	2	PCA	CA-N-CD	-3.16	102.77	113.58
2	U	2	PCA	OE-CD-CG	-3.15	121.09	126.72
2	S	2	PCA	OE-CD-CG	-3.12	121.16	126.72
2	H	2	PCA	CA-N-CD	-3.09	102.99	113.58
2	O	2	PCA	CG-CD-N	3.09	115.94	108.39
2	O	2	PCA	CA-N-CD	-3.03	103.19	113.58
2	M	2	PCA	CA-N-CD	-3.03	103.20	113.58
2	H	2	PCA	CB-CA-N	3.03	111.56	103.24
2	U	2	PCA	CA-N-CD	-3.00	103.30	113.58
2	M	2	PCA	CG-CD-N	2.98	115.68	108.39
2	H	2	PCA	CG-CD-N	2.90	115.49	108.39
2	U	2	PCA	CB-CA-N	2.90	111.22	103.24
2	Q	2	PCA	OE-CD-CG	-2.87	121.59	126.72
2	S	2	PCA	CG-CD-N	2.86	115.40	108.39
2	U	2	PCA	CG-CD-N	2.79	115.22	108.39
2	Q	2	PCA	CA-N-CD	-2.68	104.41	113.58
2	M	2	PCA	CB-CA-N	2.63	110.47	103.24
2	O	2	PCA	CB-CA-N	2.59	110.35	103.24
2	Q	2	PCA	CG-CD-N	2.53	114.58	108.39
2	Q	2	PCA	CB-CA-N	2.49	110.11	103.24
2	U	2	PCA	CB-CA-C	-2.38	109.39	112.66
2	H	2	PCA	CB-CA-C	-2.23	109.59	112.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	M	2	PCA	1	0
2	S	2	PCA	1	0

5.5 Carbohydrates

22 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$
4	NAG	G	1	1,4	14,14,15	0.51	0	17,19,21	1.24	1 (5%)
4	NAG	G	2	4	14,14,15	0.49	0	17,19,21	0.57	0
4	BMA	G	3	4	11,11,12	1.25	1 (9%)	15,15,17	0.96	1 (6%)
4	MAN	G	4	4	11,11,12	1.24	2 (18%)	15,15,17	1.00	1 (6%)
5	NAG	I	1	1,5	14,14,15	0.46	0	17,19,21	1.45	2 (11%)
5	NAG	I	2	5	14,14,15	0.38	0	17,19,21	0.44	0
5	BMA	I	3	5	11,11,12	0.87	0	15,15,17	0.87	0
6	NAG	J	1	1,6	14,14,15	0.38	0	17,19,21	1.40	2 (11%)
6	NAG	J	2	6	14,14,15	0.62	1 (7%)	17,19,21	0.58	0
6	BMA	J	3	6	11,11,12	0.88	0	15,15,17	0.90	1 (6%)
6	MAN	J	4	6	11,11,12	1.34	1 (9%)	15,15,17	0.98	1 (6%)
6	MAN	J	5	6	11,11,12	0.99	0	15,15,17	0.82	1 (6%)
5	NAG	K	1	1,5	14,14,15	0.58	0	17,19,21	1.26	2 (11%)
5	NAG	K	2	5	14,14,15	0.79	1 (7%)	17,19,21	0.42	0
5	BMA	K	3	5	11,11,12	1.04	1 (9%)	15,15,17	0.83	0
4	NAG	W	1	1,4	14,14,15	0.61	1 (7%)	17,19,21	0.57	0
4	NAG	W	2	4	14,14,15	0.33	0	17,19,21	0.41	0
4	BMA	W	3	4	11,11,12	0.97	0	15,15,17	1.02	1 (6%)
4	MAN	W	4	4	11,11,12	1.45	2 (18%)	15,15,17	1.31	1 (6%)
5	NAG	X	1	1,5	14,14,15	0.81	1 (7%)	17,19,21	1.34	1 (5%)
5	NAG	X	2	5	14,14,15	0.46	0	17,19,21	0.52	0
5	BMA	X	3	5	11,11,12	1.14	1 (9%)	15,15,17	0.77	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	G	1	1,4	-	4/6/23/26	0/1/1/1
4	NAG	G	2	4	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	BMA	G	3	4	-	0/2/19/22	0/1/1/1
4	MAN	G	4	4	-	0/2/19/22	0/1/1/1
5	NAG	I	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	I	2	5	-	1/6/23/26	0/1/1/1
5	BMA	I	3	5	-	0/2/19/22	0/1/1/1
6	NAG	J	1	1,6	-	4/6/23/26	0/1/1/1
6	NAG	J	2	6	-	1/6/23/26	0/1/1/1
6	BMA	J	3	6	-	0/2/19/22	0/1/1/1
6	MAN	J	4	6	-	0/2/19/22	0/1/1/1
6	MAN	J	5	6	-	0/2/19/22	0/1/1/1
5	NAG	K	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	K	2	5	-	1/6/23/26	0/1/1/1
5	BMA	K	3	5	-	0/2/19/22	0/1/1/1
4	NAG	W	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	W	2	4	-	1/6/23/26	0/1/1/1
4	BMA	W	3	4	-	0/2/19/22	0/1/1/1
4	MAN	W	4	4	-	2/2/19/22	0/1/1/1
5	NAG	X	1	1,5	-	3/6/23/26	0/1/1/1
5	NAG	X	2	5	-	1/6/23/26	0/1/1/1
5	BMA	X	3	5	-	0/2/19/22	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	J	4	MAN	C2-C3	3.05	1.57	1.52
4	W	4	MAN	C1-C2	3.03	1.59	1.52
5	K	2	NAG	O5-C1	-2.89	1.38	1.43
5	X	1	NAG	C1-C2	-2.70	1.48	1.52
4	G	4	MAN	C4-C5	2.40	1.58	1.53
5	X	3	BMA	C2-C3	2.39	1.56	1.52
4	G	4	MAN	C1-C2	2.17	1.57	1.52
4	W	4	MAN	O5-C1	2.17	1.47	1.43
4	W	1	NAG	O5-C1	-2.10	1.40	1.43
5	K	3	BMA	C1-C2	2.09	1.57	1.52
4	G	3	BMA	C1-C2	2.08	1.57	1.52
6	J	2	NAG	O5-C1	-2.00	1.40	1.43

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	W	4	MAN	C1-O5-C5	4.17	117.77	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	1	NAG	C2-N2-C7	3.93	128.16	122.90
5	X	1	NAG	C2-N2-C7	3.84	128.05	122.90
4	G	1	NAG	C2-N2-C7	3.81	128.01	122.90
5	K	1	NAG	C2-N2-C7	3.75	127.93	122.90
6	J	1	NAG	C2-N2-C7	3.75	127.93	122.90
5	I	1	NAG	C1-C2-N2	3.15	115.40	110.43
6	J	1	NAG	C1-C2-N2	3.04	115.23	110.43
6	J	4	MAN	C1-O5-C5	2.71	115.82	112.19
4	W	3	BMA	C1-O5-C5	2.70	115.81	112.19
4	G	4	MAN	C1-O5-C5	2.68	115.78	112.19
5	K	1	NAG	C1-C2-N2	2.47	114.33	110.43
4	G	3	BMA	C1-O5-C5	2.27	115.23	112.19
6	J	5	MAN	C1-O5-C5	2.27	115.23	112.19
6	J	3	BMA	C1-O5-C5	2.18	115.11	112.19
5	X	3	BMA	C1-O5-C5	2.15	115.06	112.19

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	1	NAG	C8-C7-N2-C2
4	G	1	NAG	O7-C7-N2-C2
5	I	1	NAG	C8-C7-N2-C2
5	I	1	NAG	O7-C7-N2-C2
5	K	1	NAG	C8-C7-N2-C2
5	K	1	NAG	O7-C7-N2-C2
5	X	1	NAG	C8-C7-N2-C2
5	X	1	NAG	O7-C7-N2-C2
6	J	1	NAG	C8-C7-N2-C2
6	J	1	NAG	O7-C7-N2-C2
4	W	4	MAN	O5-C5-C6-O6
4	W	1	NAG	C4-C5-C6-O6
5	K	2	NAG	O5-C5-C6-O6
4	W	2	NAG	O5-C5-C6-O6
6	J	2	NAG	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
5	I	2	NAG	O5-C5-C6-O6
5	X	2	NAG	O5-C5-C6-O6
5	X	1	NAG	C3-C2-N2-C7
4	W	1	NAG	O5-C5-C6-O6
4	G	1	NAG	C1-C2-N2-C7
6	J	1	NAG	C1-C2-N2-C7

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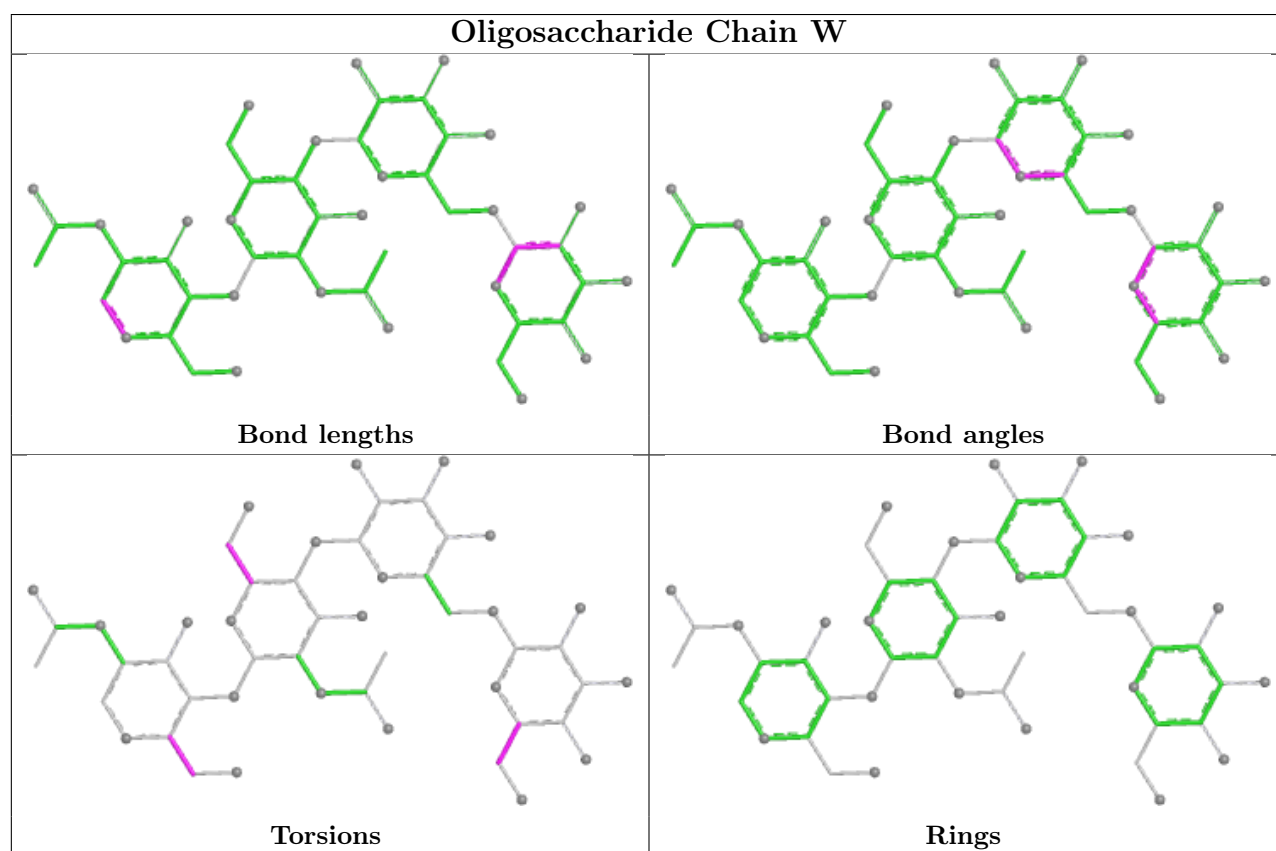
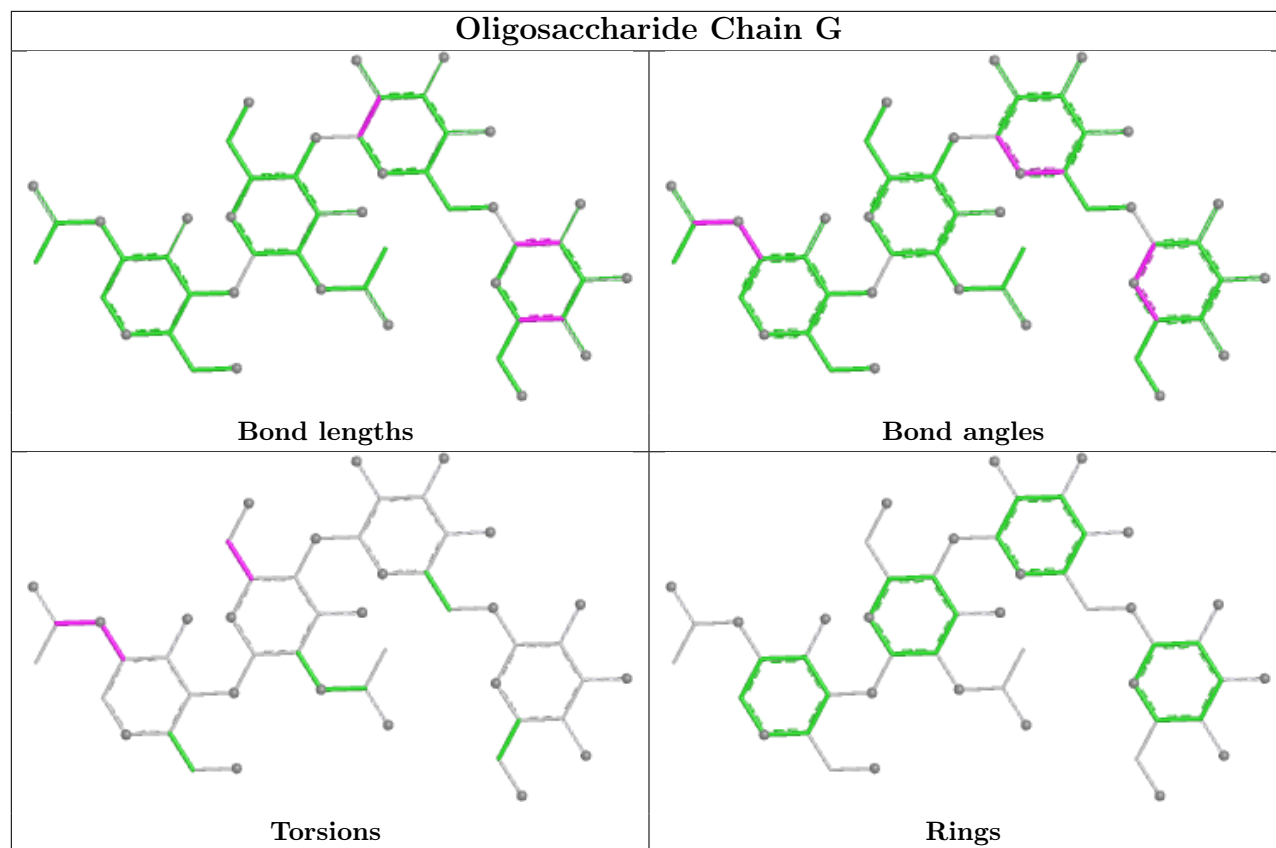
Mol	Chain	Res	Type	Atoms
5	I	1	NAG	C3-C2-N2-C7
5	K	1	NAG	C3-C2-N2-C7
6	J	1	NAG	C3-C2-N2-C7
4	W	4	MAN	C4-C5-C6-O6
5	I	1	NAG	C1-C2-N2-C7
5	K	1	NAG	C1-C2-N2-C7
4	G	1	NAG	C3-C2-N2-C7

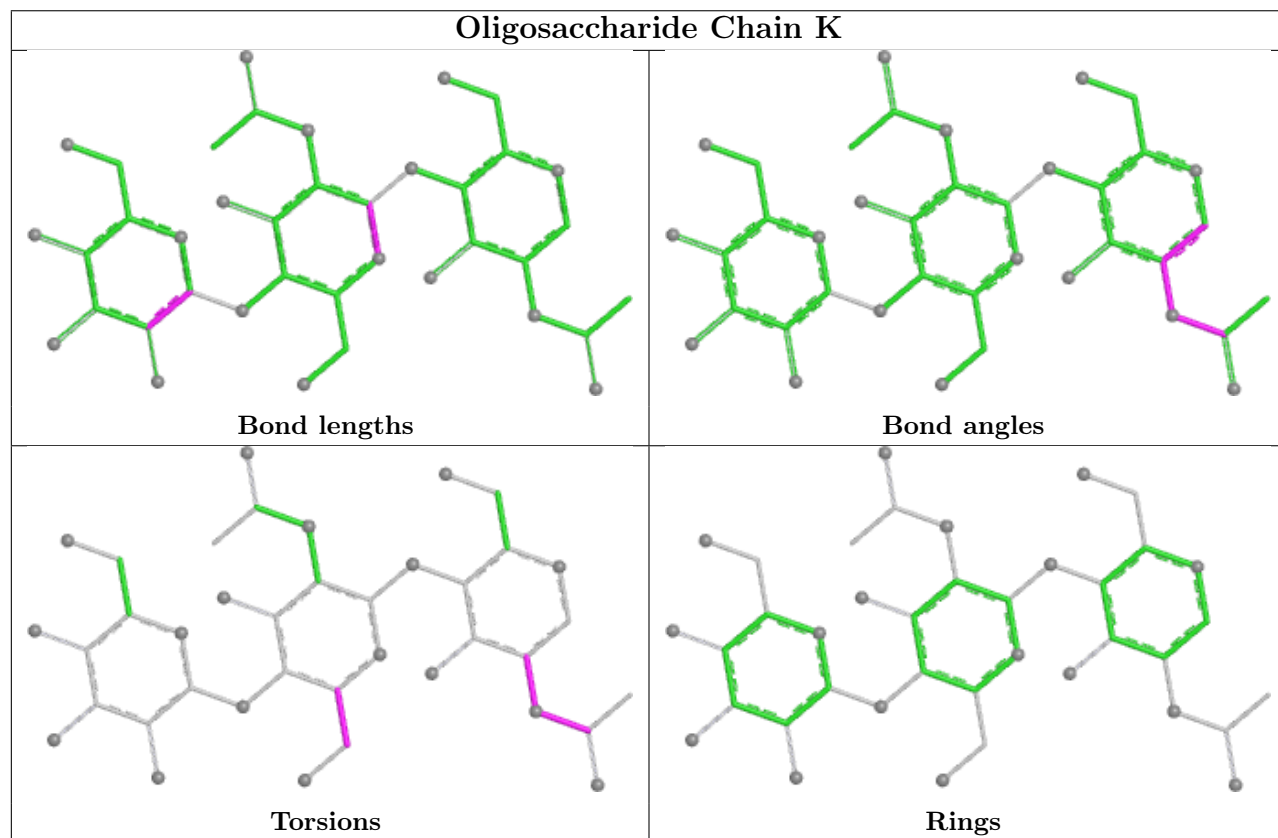
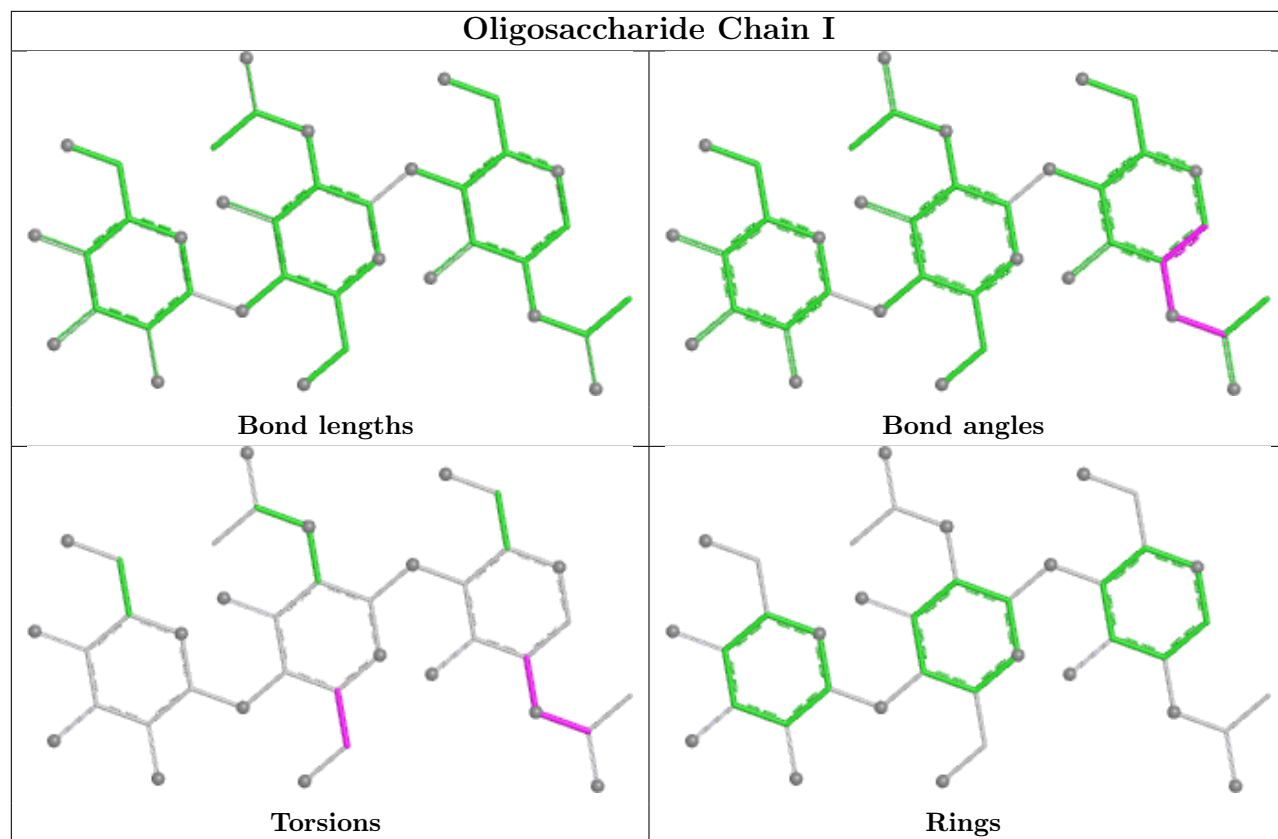
There are no ring outliers.

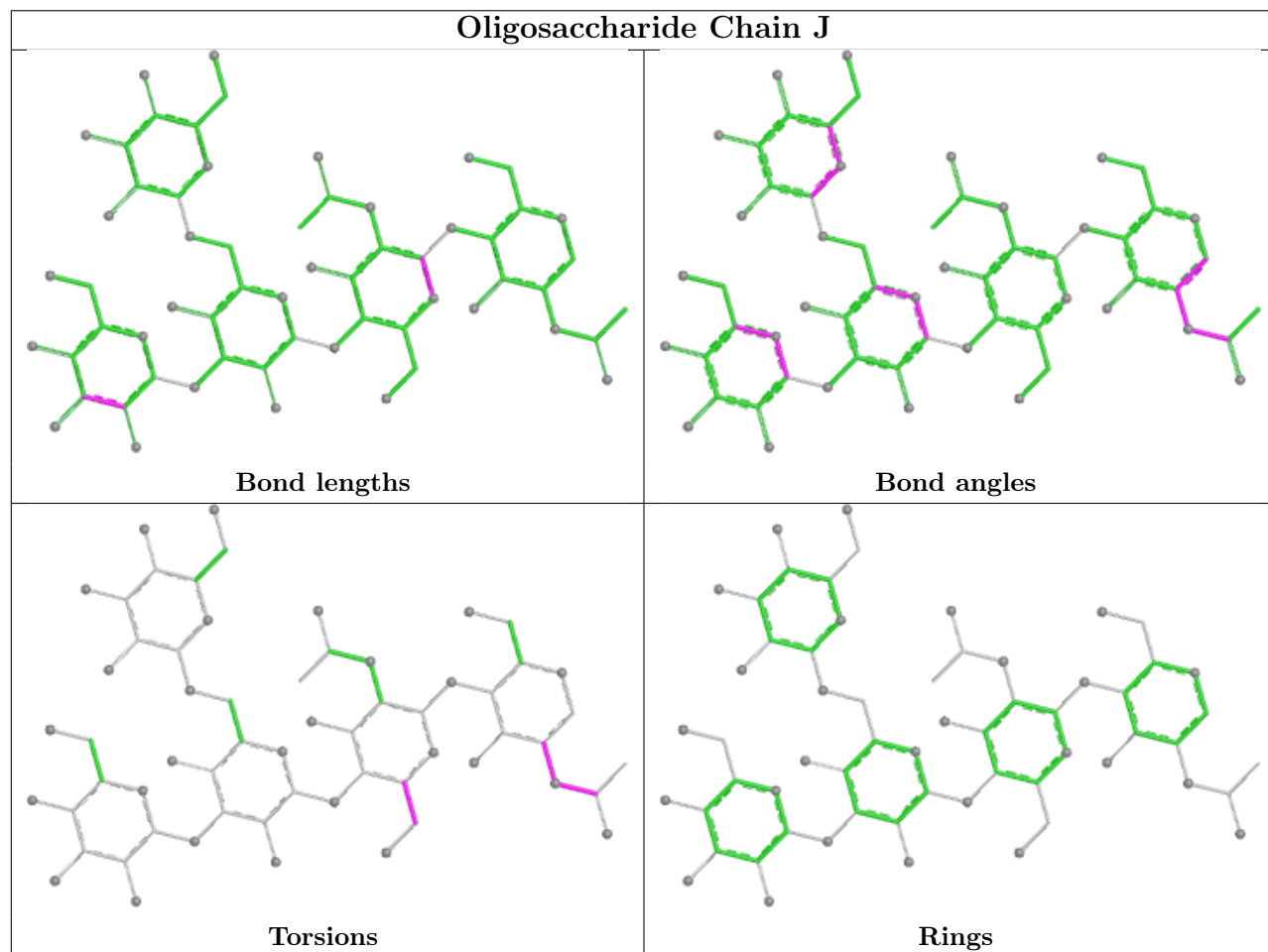
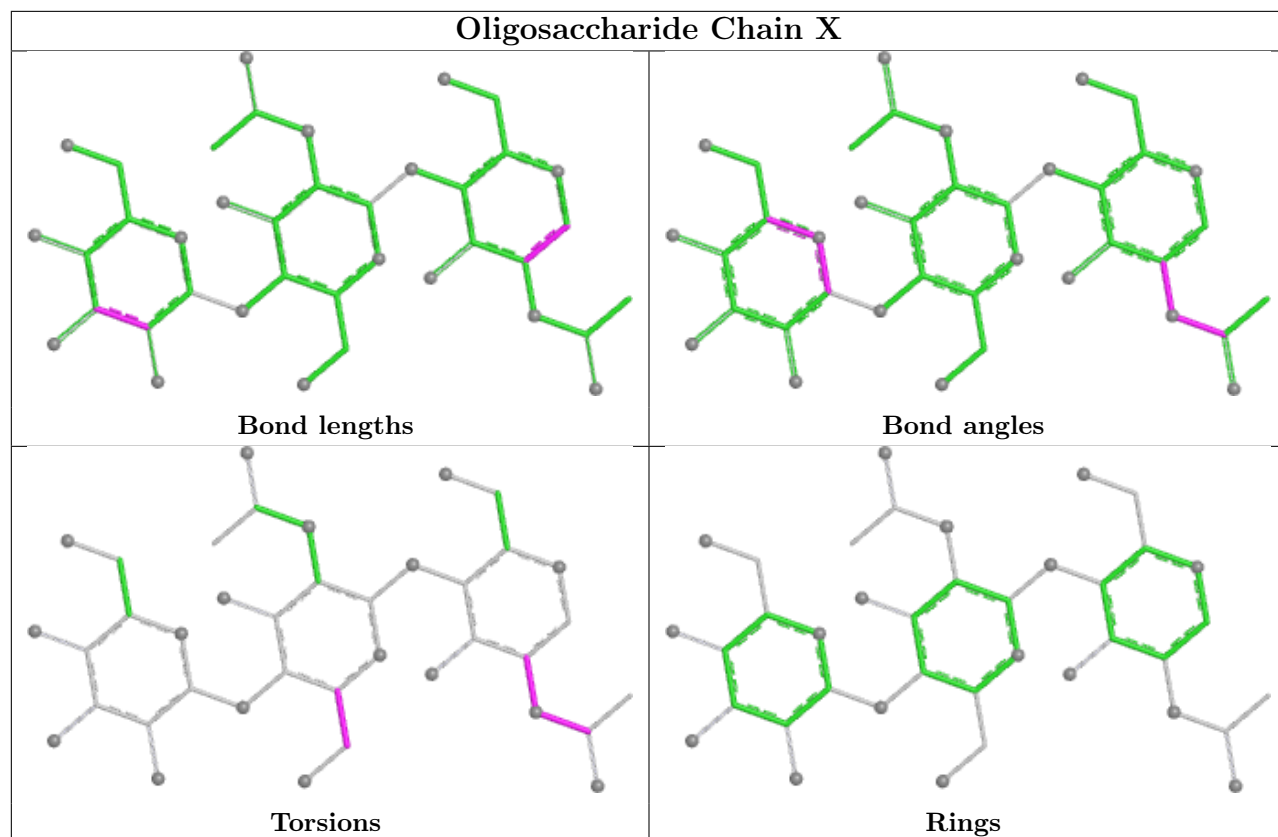
6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	1	NAG	1	0
4	W	1	NAG	1	0
6	J	1	NAG	1	0
5	X	1	NAG	4	0
5	K	1	NAG	1	0
5	I	1	NAG	1	0

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







5.6 Ligand geometry

There are no ligands in this entry.

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	362/362 (100%)	-0.25	0 100 100	35, 52, 81, 133	0
1	B	362/362 (100%)	-0.25	0 100 100	35, 53, 82, 133	0
1	C	362/362 (100%)	-0.26	1 (0%) 90 82	40, 55, 86, 143	0
1	D	362/362 (100%)	-0.21	1 (0%) 90 82	39, 52, 80, 133	0
1	E	362/362 (100%)	-0.20	2 (0%) 85 73	39, 55, 84, 124	0
1	F	362/362 (100%)	-0.32	0 100 100	33, 48, 84, 133	0
2	H	222/223 (99%)	0.22	12 (5%) 31 21	42, 76, 159, 211	0
2	M	222/223 (99%)	-0.01	8 (3%) 46 31	37, 66, 132, 187	0
2	O	222/223 (99%)	0.43	16 (7%) 21 15	48, 84, 191, 218	0
2	Q	222/223 (99%)	0.33	9 (4%) 41 27	48, 84, 152, 212	0
2	S	222/223 (99%)	0.24	9 (4%) 41 27	51, 80, 140, 189	0
2	U	222/223 (99%)	0.26	18 (8%) 18 13	43, 69, 178, 235	0
3	L	214/214 (100%)	0.07	4 (1%) 66 48	37, 69, 150, 178	0
3	N	214/214 (100%)	-0.14	2 (0%) 81 65	34, 59, 114, 155	0
3	P	214/214 (100%)	0.33	9 (4%) 40 26	44, 79, 181, 233	0
3	R	214/214 (100%)	0.05	3 (1%) 73 55	43, 69, 144, 163	0
3	T	214/214 (100%)	-0.07	1 (0%) 87 76	44, 68, 108, 128	0
3	V	214/214 (100%)	0.32	11 (5%) 33 22	39, 75, 165, 199	0
All	All	4788/4794 (99%)	-0.02	106 (2%) 62 43	33, 59, 148, 235	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	S	53	SER	4.4
2	O	132	PRO	4.1
2	O	131	THR	4.1

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Mol	Chain	Res	Type	RSRZ
2	U	55	THR	4.0
2	S	132	PRO	4.0
2	O	210	PRO	3.8
2	U	132	PRO	3.8
3	V	114	VAL	3.7
2	Q	29	ILE	3.7
2	M	129	GLY	3.6
2	U	29	ILE	3.4
2	S	29	ILE	3.3
2	S	55	THR	3.3
2	U	131	THR	3.2
2	M	212	THR	3.1
2	U	30	SER	3.0
2	H	134	SER	2.9
3	P	200	SER	2.9
3	P	199	THR	2.9
2	M	132	PRO	2.9
2	H	209	ALA	2.9
3	N	1	ASP	2.8
3	R	188	LYS	2.8
2	S	30	SER	2.8
2	S	213	CYS	2.8
2	H	132	PRO	2.7
2	O	55	THR	2.7
2	S	28	THR	2.7
3	V	199	THR	2.7
2	Q	186	SER	2.7
3	P	121	ALA	2.7
2	H	29	ILE	2.6
2	H	186	SER	2.6
2	H	211	SER	2.6
2	O	130	ASP	2.6
3	V	208	GLY	2.6
2	U	28	THR	2.6
2	U	133	SER	2.5
3	L	210	CYS	2.5
2	Q	187	SER	2.5
2	O	199	ALA	2.5
2	H	131	THR	2.5
2	M	128	CYS	2.4
2	O	179	SER	2.4
3	P	188	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
2	Q	28	THR	2.4
2	O	29	ILE	2.4
3	L	208	GLY	2.4
3	V	186	SER	2.4
2	U	210	PRO	2.4
3	P	201	VAL	2.4
2	Q	30	SER	2.4
2	Q	53	SER	2.4
2	O	129	GLY	2.3
2	O	187	SER	2.3
3	V	121	ALA	2.3
2	O	95	ASP	2.3
3	V	201	VAL	2.3
2	M	131	THR	2.3
2	Q	55	THR	2.3
2	H	213	CYS	2.3
2	O	213	CYS	2.3
3	P	113	THR	2.3
2	U	209	ALA	2.3
2	M	30	SER	2.3
2	M	213	CYS	2.2
2	U	213	CYS	2.2
3	V	210	CYS	2.2
1	E	306	ASP	2.2
2	H	210	PRO	2.2
2	O	126	PRO	2.2
1	D	291	ARG	2.2
3	P	206	ASN	2.2
3	N	210	CYS	2.2
3	P	192	CYS	2.2
2	Q	132	PRO	2.2
3	V	188	LYS	2.2
1	C	291	ARG	2.2
3	L	1	ASP	2.2
2	U	123	PRO	2.1
3	V	141	ASP	2.1
2	O	135	THR	2.1
2	M	29	ILE	2.1
2	S	134	SER	2.1
2	U	129	GLY	2.1
3	V	119	PRO	2.1
1	E	104	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
2	H	212	THR	2.1
2	U	212	THR	2.1
3	P	135	ALA	2.1
3	R	121	ALA	2.1
2	O	133	SER	2.1
2	U	130	ASP	2.1
2	S	131	THR	2.1
2	H	123	PRO	2.1
3	T	105	GLU	2.1
2	O	186	SER	2.1
3	L	116	ILE	2.1
2	H	95	ASP	2.1
2	Q	213	CYS	2.1
3	V	148	VAL	2.0
2	U	211	SER	2.0
2	U	138	LEU	2.0
2	U	95	ASP	2.0
2	U	52(A)	ILE	2.0
3	R	95(A)	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	PCA	H	2	8/9	0.86	0.12	62,67,70,72	0
2	PCA	S	2	8/9	0.87	0.14	64,69,73,74	0
2	PCA	U	2	8/9	0.90	0.11	63,64,66,67	0
2	PCA	M	2	8/9	0.92	0.10	57,62,64,66	0
2	PCA	O	2	8/9	0.92	0.10	66,68,70,72	0
2	PCA	Q	2	8/9	0.93	0.09	63,65,69,71	0

6.3 Carbohydrates [i](#)

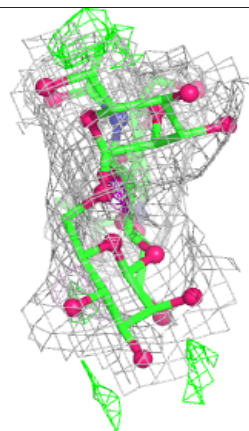
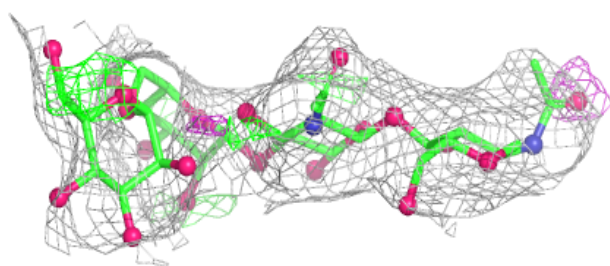
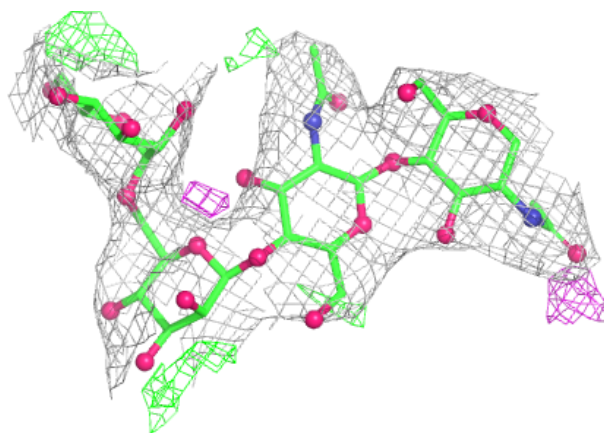
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	G	1	14/15	-	-	50,54,62,70	0
4	NAG	G	2	14/15	-	-	72,77,95,99	0
4	BMA	G	3	11/12	-	-	109,117,120,123	0
4	MAN	G	4	11/12	-	-	114,121,125,126	0
4	MAN	W	4	11/12	0.26	0.15	122,124,128,131	0
4	BMA	W	3	11/12	0.32	0.15	142,149,154,156	0
5	BMA	X	3	11/12	0.34	0.15	129,138,143,143	0
5	NAG	X	2	14/15	0.80	0.12	84,90,107,116	0
5	NAG	I	1	14/15	-	-	49,56,65,75	0
5	NAG	I	2	14/15	-	-	83,85,100,100	0
5	BMA	I	3	11/12	-	-	99,101,102,102	0
5	NAG	K	1	14/15	-	-	56,61,74,88	0
5	NAG	K	2	14/15	-	-	104,108,126,130	0
5	BMA	K	3	11/12	-	-	132,134,136,136	0
4	NAG	W	2	14/15	0.80	0.11	92,102,117,124	0
5	NAG	X	1	14/15	0.87	0.13	51,55,66,77	0
4	NAG	W	1	14/15	0.91	0.10	59,67,77,86	0
6	NAG	J	1	14/15	-	-	60,67,76,84	0
6	NAG	J	2	14/15	-	-	87,95,120,129	0
6	BMA	J	3	11/12	-	-	164,169,177,178	0
6	MAN	J	4	11/12	-	-	151,159,161,162	0
6	MAN	J	5	11/12	-	-	132,141,145,147	0

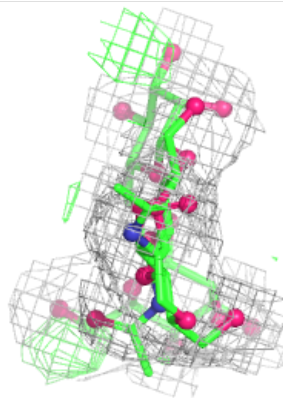
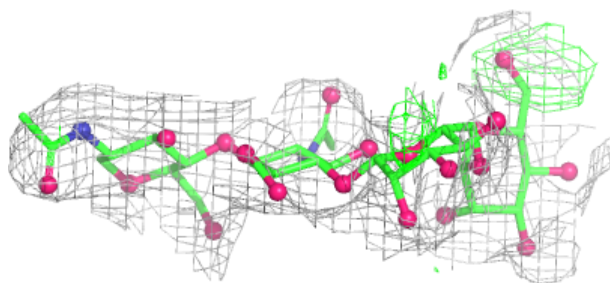
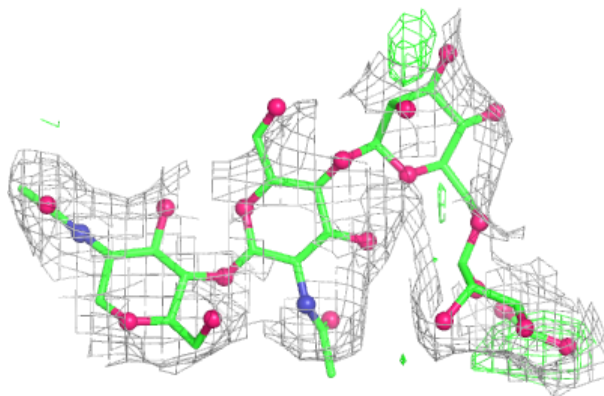
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain G:

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

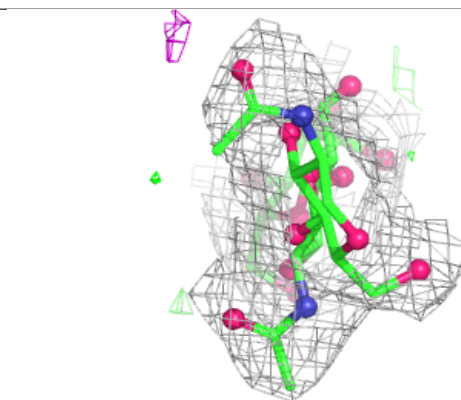
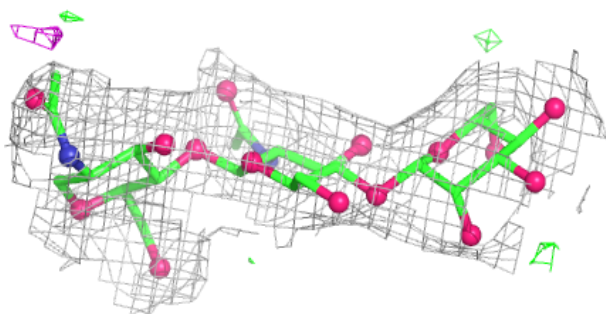
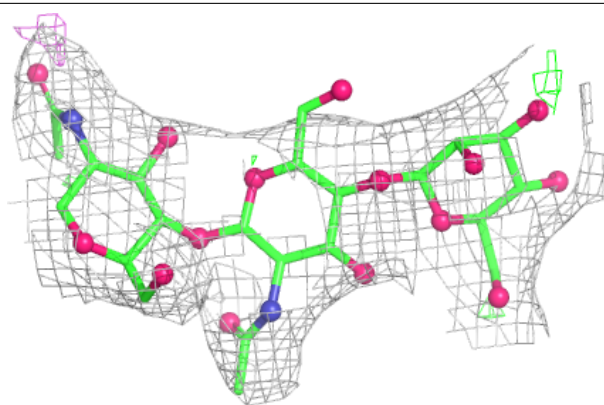
**Electron density around Chain W:**

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

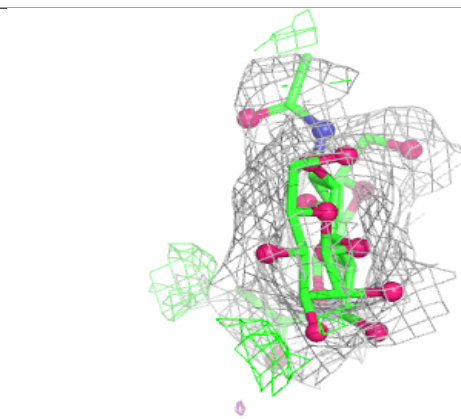
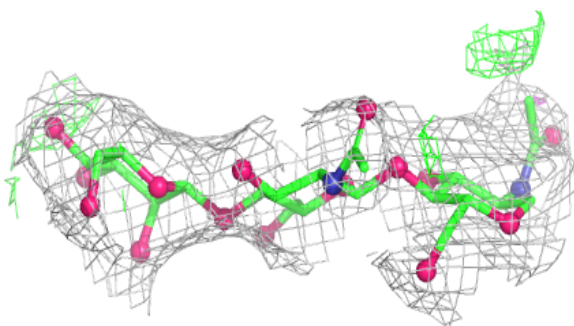
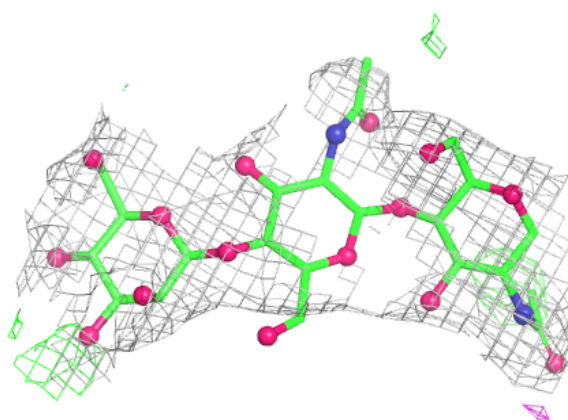


Electron density around Chain I:

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

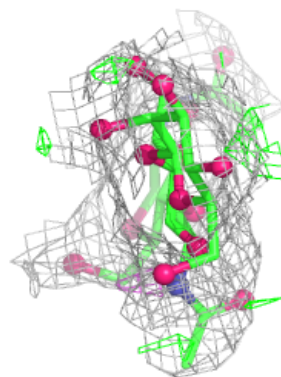
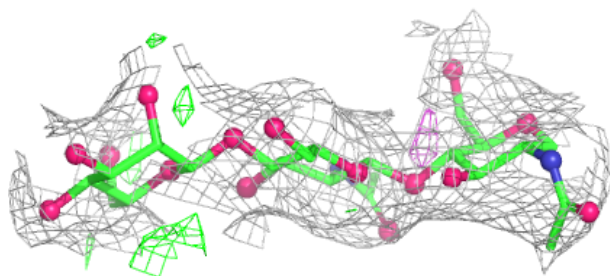
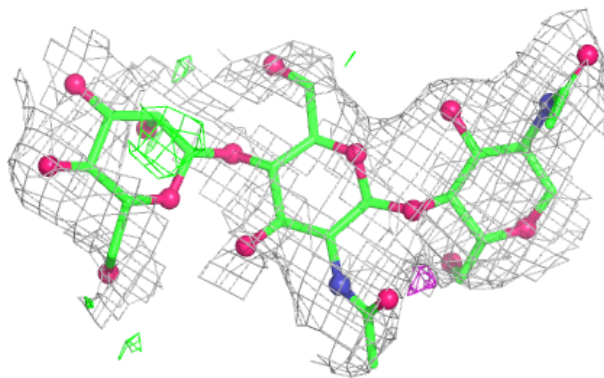
**Electron density around Chain K:**

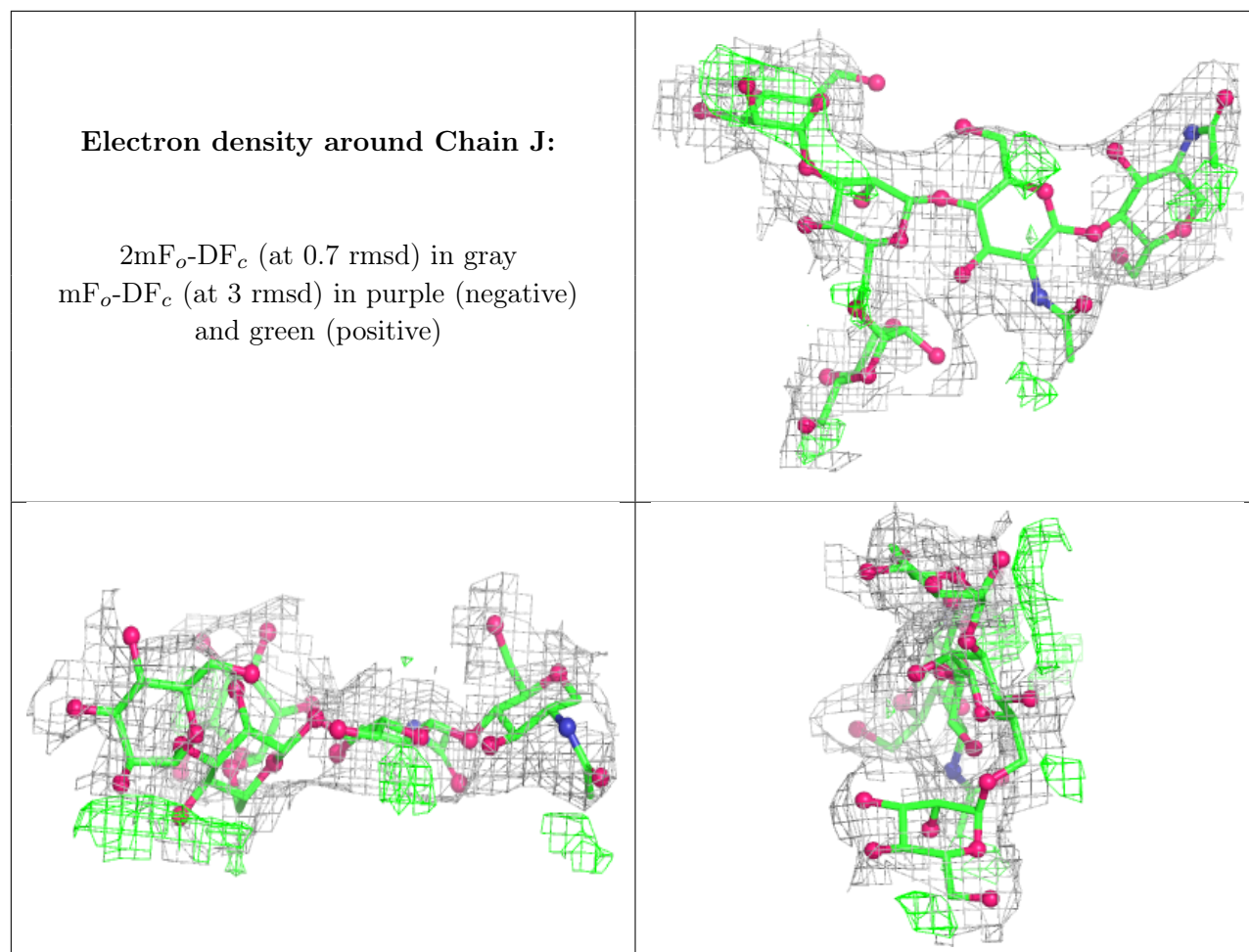
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain X:

$2mF_o-DF_c$ (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.