Full wwPDB X-ray Structure Validation Report (i)

Oct 3, 2023 – 04:57 PM EDT

PDB ID	:	6OZ4
Title	:	Crystal structure of broadly neutralizing antibody N49P6 Fab in complex with HIV-1 BG505 SOSIP.664 Env trimer ectodomain.
Authors	:	Tolbert, W.D.; Pazgier, M.
Deposited on	:	2019-05-15
Resolution	:	4.05 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 (i) 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 (i)) were used in the production of this report:

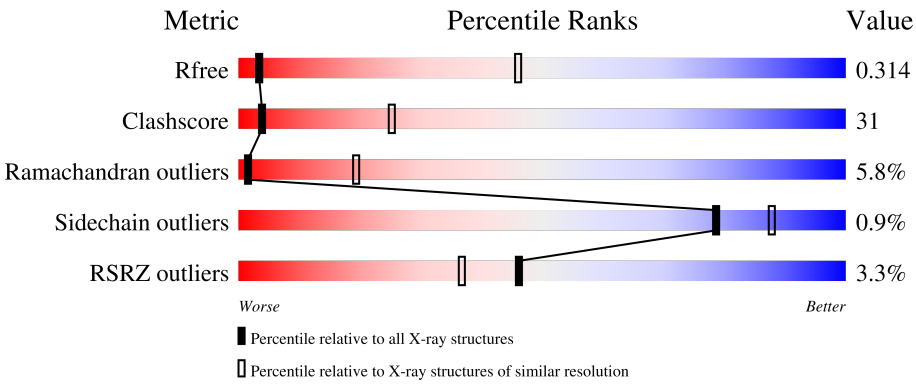
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 4.05 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	1127 (4.42-3.70)
Clashscore	141614	1033 (4.40-3.72)
Ramachandran outliers	138981	1145 (4.42-3.70)
Sidechain outliers	138945	1133 (4.42-3.70)
RSRZ outliers	127900	1005 (4.44-3.68)

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 <=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	G	481	2% 39% 51% • 7%
2	B	153	5% 41% 36% • 20%
3	H	229	2% 44% 51% • •
4	L	205	5% 41% 52% • •
5	A	3	33% 67%

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Mol	Chain	Length	Quality of chain
5	F	3	100%
5	M	3	33% 67%
6	C	2	50% 50%
6	E	2	100%
6	I	2	50% 50%
6	J	2	100%
6	N	2	100%
6	O	2	100%
6	P	2	100%
7	D	4	25% 75%
8	K	7	29% 71%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	BMA	A	3	-	-	-	X
6	NAG	C	2	-	-	-	X
6	NAG	E	2	-	-	-	X
6	NAG	I	1	-	-	-	X
7	BMA	D	3	-	-	-	X
7	MAN	D	4	-	-	-	X
9	NAG	G	603	-	-	-	X

2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 8269 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 Envelope glycoprotein gp160.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	G	449	3536	2219	625	664	28	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	332	ASN	THR	engineered mutation	UNP Q2N0S6
G	501	CYS	ALA	engineered mutation	UNP Q2N0S6
G	509	ARG	-	expression tag	UNP Q2N0S6
G	510	ARG	-	expression tag	UNP Q2N0S6
G	511	ARG	-	expression tag	UNP Q2N0S6
G	512	ARG	-	expression tag	UNP Q2N0S6
G	513	ARG	-	expression tag	UNP Q2N0S6

- Molecule 2 is a protein called Envelope glycoprotein gp41.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	122	971	611	168	186	6	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	559	PRO	ILE	engineered mutation	UNP Q2N0S6
B	605	CYS	THR	engineered mutation	UNP Q2N0S6

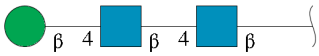
- Molecule 3 is a protein called N49P6 antibody Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	H	225	1722	1082	303	329	8	0	0	0

- Molecule 4 is a protein called N49P6 antibody light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	L	200	1496	936	254	301	5	0	0	0

- 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
			Total	C	N	O			
5	A	3	39	22	2	15	0	0	0
5	F	3	39	22	2	15	0	0	0
5	M	3	39	22	2	15	0	0	0

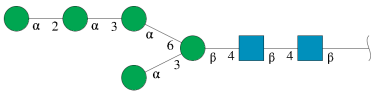
- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
6	C	2	28	16	2	10	0	0	0
6	E	2	28	16	2	10	0	0	0
6	I	2	28	16	2	10	0	0	0
6	J	2	28	16	2	10	0	0	0
6	N	2	28	16	2	10	0	0	0
6	O	2	28	16	2	10	0	0	0
6	P	2	28	16	2	10	0	0	0

- Molecule 7 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.

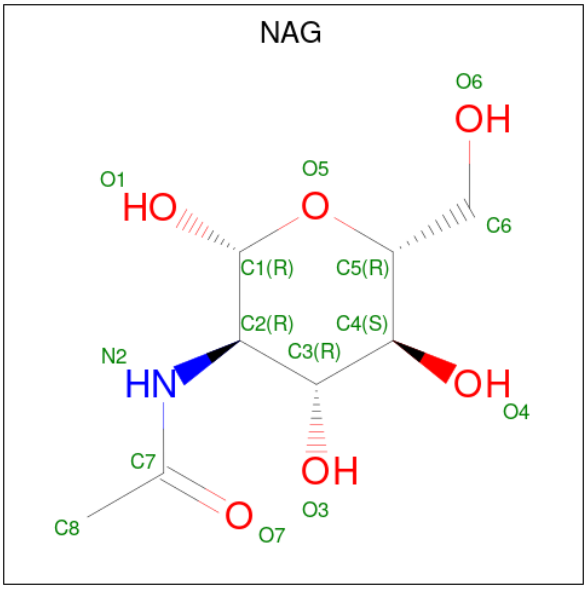
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
7	D	4	50	28	2	20	0	0	0

- Molecule 8 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]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
			Total	C	N	O			
8	K	7	83	46	2	35	0	0	0

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



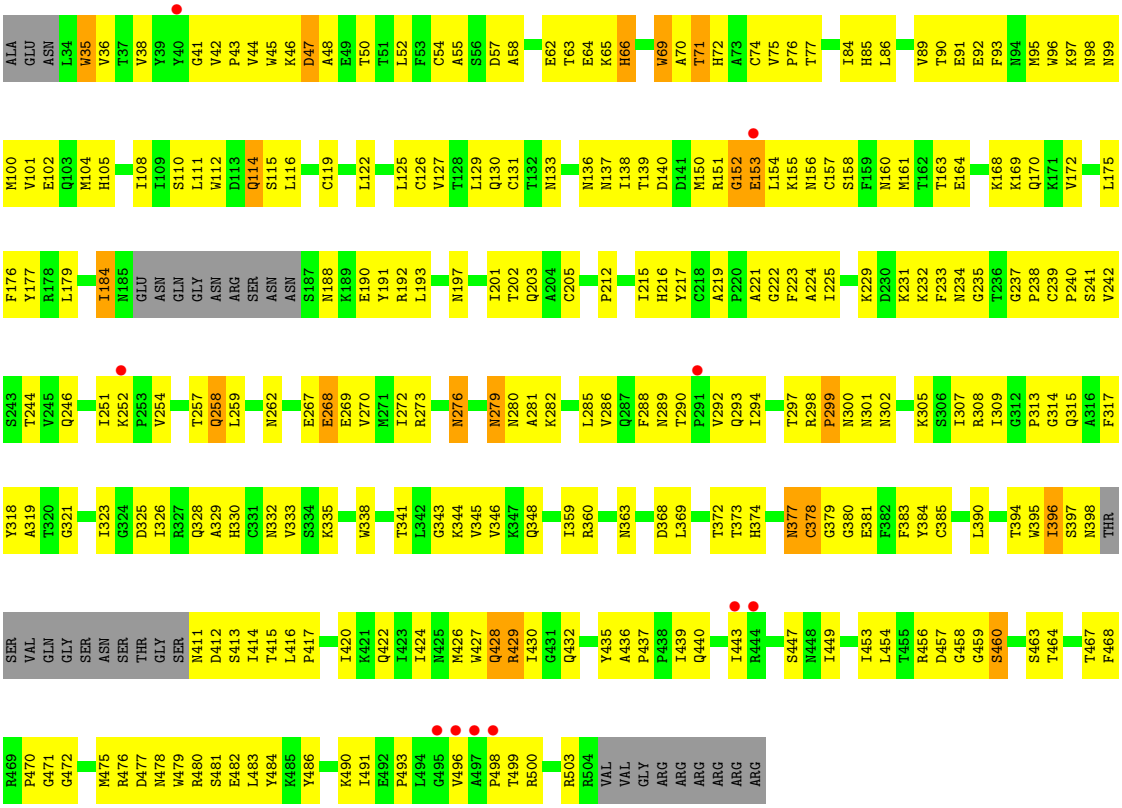
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
9	G	1	14	8	1	5	0	0
			Total	C	N	O		
9	G	1	14	8	1	5	0	0
			Total	C	N	O		
9	G	1	14	8	1	5	0	0
			Total	C	N	O		
9	G	1	14	8	1	5	0	0
			Total	C	N	O		
9	B	1	14	8	1	5	0	0
			Total	C	N	O		
9	B	1	14	8	1	5	0	0
			Total	C	N	O		
9	B	1	14	8	1	5	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

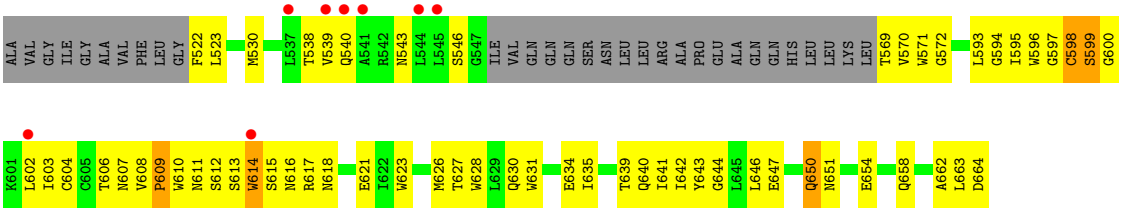
• Molecule 1: Envelope glycoprotein gp160

Chain G: 2% 39% 51% • 7%



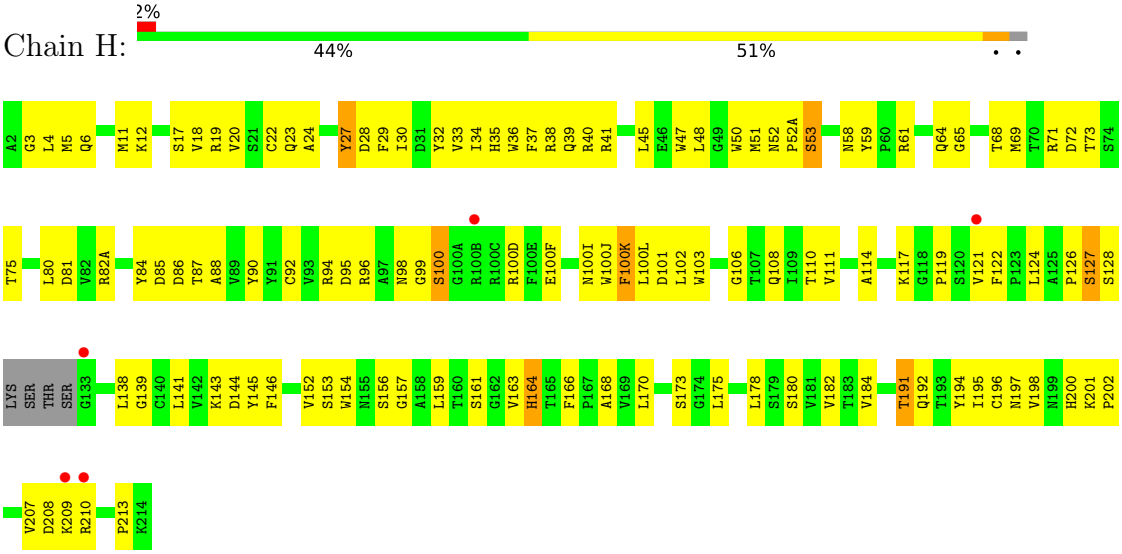
• Molecule 2: Envelope glycoprotein gp41

Chain B: 5% 41% 36% • 20%



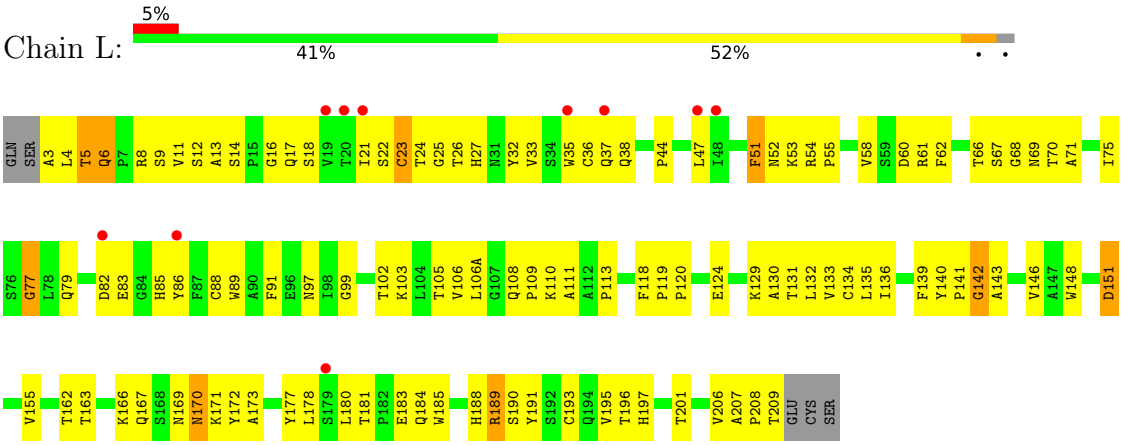
• Molecule 3: N49P6 antibody Fab heavy chain

Chain H:



• Molecule 4: N49P6 antibody light chain

Chain L:

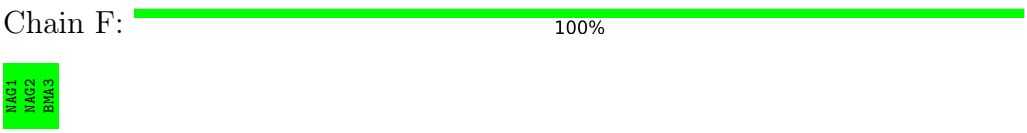


• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

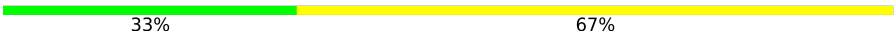
Chain A:

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

Chain F:



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

Chain M: 

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

Chain C:

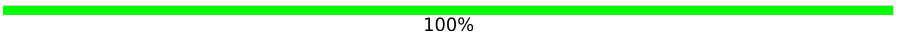
• Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:

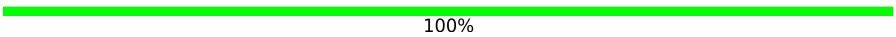
• Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:

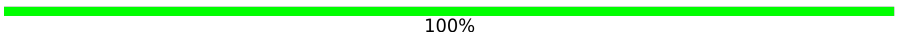
• Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J: 

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

Chain N: 

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

Chain O: 

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

Chain P: 100%

NAG1
NAG2

• Molecule 7: 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

Chain D: 25% 75%

NAG1
NAG2
BMA3
MAN4

• Molecule 8: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K: 29% 71%

NAG1
NAG2
BMA3
MAN4
MAN5
MAN6
MAN7

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α, β, γ	164.70Å 164.70Å 164.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.95 – 4.05 49.66 – 4.03	Depositor EDS
% Data completeness (in resolution range)	99.4 (39.95-4.05) 99.7 (49.66-4.03)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.15 (at 4.00Å)	Xtriage
Refinement program	PHENIX (1.15.2_3472: ???), REFMAC	Depositor
R, R_{free}	0.255 , 0.315 0.254 , 0.314	Depositor DCC
R_{free} test set	580 reflections (4.62%)	wwPDB-VP
Wilson B-factor (Å^2)	208.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 245.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	0.072 for l,-k,h	Xtriage
F_o,F_c correlation	0.87	EDS
Total number of atoms	8269	wwPDB-VP
Average B, all atoms (Å^2)	225.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.99% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: BMA, NAG, MAN

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	G	0.42	0/3610	0.68	0/4901
2	B	0.38	0/988	0.63	0/1340
3	H	0.44	0/1765	0.70	0/2401
4	L	0.40	0/1534	0.67	0/2094
All	All	0.42	0/7897	0.68	0/10736

There are no bond length outliers.

There are no bond angle outliers.

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	G	3536	0	3471	251	0
2	B	971	0	945	65	0
3	H	1722	0	1664	105	0
4	L	1496	0	1447	114	0
5	A	39	0	34	0	0
5	F	39	0	34	0	0
5	M	39	0	34	2	0
6	C	28	0	25	2	0
6	E	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	I	28	0	25	0	0
6	J	28	0	25	0	0
6	N	28	0	25	0	0
6	O	28	0	25	0	0
6	P	28	0	25	0	0
7	D	50	0	43	1	0
8	K	83	0	70	2	0
9	B	42	0	39	2	0
9	G	56	0	52	4	0
All	All	8269	0	8008	499	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 31.

All (499) 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:G:43:PRO:CG	2:B:628:TRP:CD1	2.09	1.33
1:G:43:PRO:HG2	2:B:628:TRP:CD1	1.72	1.19
1:G:101:VAL:HB	1:G:483:LEU:HD12	1.21	1.13
1:G:43:PRO:HG2	2:B:628:TRP:CG	1.84	1.10
1:G:184:ILE:HD11	1:G:190:GLU:HG2	1.27	1.09
1:G:43:PRO:HG3	2:B:628:TRP:CD1	1.82	1.06
2:B:598:CYS:HA	2:B:604:CYS:SG	2.01	1.00
3:H:6:GLN:HG3	3:H:106:GLY:H	1.27	0.99
1:G:43:PRO:CG	2:B:628:TRP:CG	2.44	0.97
1:G:101:VAL:CB	1:G:483:LEU:HD12	1.95	0.96
1:G:43:PRO:HD2	2:B:628:TRP:CD2	2.07	0.90
1:G:503:ARG:NH2	2:B:604:CYS:SG	2.44	0.90
4:L:180:LEU:HD11	4:L:185:TRP:CB	2.01	0.90
1:G:232:LYS:HE3	1:G:269:GLU:HB2	1.53	0.90
4:L:180:LEU:HD11	4:L:185:TRP:HB2	1.54	0.87
3:H:163:VAL:HG12	3:H:182:VAL:HG22	1.54	0.87
1:G:301:ASN:HB3	1:G:323:ILE:HB	1.57	0.86
1:G:140:ASP:HA	1:G:151:ARG:HB2	1.56	0.85
3:H:99:GLY:O	3:H:100(D):ARG:NH1	2.09	0.85
1:G:476:ARG:HA	1:G:479:TRP:HE3	1.39	0.85
4:L:32:TYR:CE1	4:L:51:PHE:HE2	1.95	0.84
1:G:35:TRP:HA	2:B:609:PRO:HA	1.59	0.84
1:G:43:PRO:HD2	2:B:628:TRP:CG	2.14	0.83
3:H:47:TRP:HZ2	3:H:50:TRP:HD1	1.24	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:270:VAL:HG12	1:G:289:ASN:H	1.43	0.82
2:B:530:MET:HB2	2:B:623:TRP:HA	1.61	0.82
2:B:598:CYS:CA	2:B:604:CYS:SG	2.68	0.82
3:H:94:ARG:NH2	3:H:101:ASP:OD2	2.10	0.81
4:L:146:VAL:HG22	4:L:195:VAL:HG12	1.63	0.81
1:G:35:TRP:HB2	1:G:499:THR:O	1.80	0.81
3:H:119:PRO:HB3	3:H:145:TYR:HB3	1.63	0.80
1:G:476:ARG:HA	1:G:479:TRP:CE3	2.16	0.80
3:H:191:THR:HG23	3:H:192:GLN:H	1.47	0.80
1:G:57:ASP:HA	1:G:77:THR:H	1.46	0.79
4:L:196:THR:HG22	4:L:201:THR:HG22	1.64	0.79
1:G:43:PRO:CD	2:B:628:TRP:CG	2.66	0.79
1:G:377:ASN:HD22	1:G:378:CYS:N	1.81	0.79
1:G:90:THR:HG22	1:G:240:PRO:HA	1.65	0.79
4:L:32:TYR:CE1	4:L:51:PHE:CE2	2.70	0.79
3:H:68:THR:HB	3:H:81:ASP:HB2	1.65	0.78
4:L:32:TYR:HD1	4:L:51:PHE:CD2	2.01	0.78
4:L:189:ARG:HD3	4:L:208:PRO:HD2	1.66	0.77
3:H:47:TRP:CZ2	3:H:50:TRP:HD1	2.02	0.77
1:G:335:LYS:HG3	1:G:413:SER:HA	1.66	0.77
1:G:116:LEU:HD13	1:G:202:THR:HG21	1.65	0.77
3:H:156:SER:H	3:H:197:ASN:ND2	1.83	0.76
9:G:602:NAG:H4	4:L:32:TYR:HE2	1.51	0.75
1:G:292:VAL:HB	1:G:449:ILE:HB	1.69	0.74
4:L:162:THR:HG22	4:L:163:THR:H	1.50	0.74
3:H:100(I):ASN:OD1	3:H:100(J):TRP:N	2.20	0.74
1:G:43:PRO:HG2	2:B:628:TRP:CB	2.18	0.73
3:H:5:MET:HB2	3:H:23:GLN:HB3	1.71	0.73
3:H:124:LEU:HD11	3:H:141:LEU:HD23	1.70	0.72
9:G:602:NAG:H4	4:L:32:TYR:CE2	2.24	0.72
3:H:6:GLN:HG3	3:H:106:GLY:N	2.04	0.72
1:G:152:GLY:O	1:G:154:LEU:N	2.23	0.72
4:L:32:TYR:CD1	4:L:51:PHE:CE2	2.78	0.72
1:G:285:LEU:HD11	1:G:481:SER:HB3	1.72	0.71
1:G:457:ASP:OD2	1:G:467:THR:OG1	2.07	0.71
4:L:6:GLN:HB3	4:L:102:THR:CG2	2.21	0.71
1:G:232:LYS:NZ	1:G:268:GLU:O	2.24	0.70
1:G:101:VAL:CA	1:G:483:LEU:CD1	2.68	0.70
1:G:168:LYS:HG2	1:G:169:LYS:O	1.91	0.70
1:G:101:VAL:CA	1:G:483:LEU:HD12	2.22	0.69
4:L:32:TYR:HE1	4:L:51:PHE:CE2	2.10	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:254:VAL:HG23	1:G:478:ASN:HD21	1.56	0.69
2:B:522:PHE:HD1	2:B:539:VAL:HG12	1.56	0.69
3:H:143:LYS:NZ	4:L:131:THR:OG1	2.26	0.68
3:H:30:ILE:HD11	3:H:53:SER:HA	1.75	0.68
1:G:35:TRP:O	1:G:498:PRO:HA	1.92	0.68
1:G:456:ARG:HG3	1:G:468:PHE:HE1	1.58	0.68
2:B:594:GLY:HA2	2:B:599:SER:HB2	1.75	0.68
1:G:101:VAL:N	1:G:483:LEU:CD1	2.57	0.68
2:B:626:MET:SD	2:B:630:GLN:NE2	2.67	0.68
2:B:658:GLN:O	2:B:662:ALA:N	2.24	0.68
4:L:11:VAL:HG21	4:L:21:ILE:HD11	1.75	0.68
4:L:32:TYR:HD1	4:L:51:PHE:HD2	1.41	0.68
4:L:32:TYR:CD1	4:L:51:PHE:CD2	2.82	0.67
1:G:92:GLU:HA	1:G:238:PRO:HA	1.74	0.67
1:G:457:ASP:HA	3:H:58:ASN:HD21	1.59	0.67
1:G:332:ASN:HA	1:G:414:ILE:O	1.95	0.67
4:L:188:HIS:HB2	4:L:191:TYR:CZ	2.30	0.67
1:G:254:VAL:O	1:G:478:ASN:ND2	2.28	0.66
1:G:86:LEU:HD23	1:G:89:VAL:HG21	1.77	0.66
4:L:180:LEU:HD11	4:L:185:TRP:HB3	1.77	0.66
1:G:360:ARG:HB3	1:G:467:THR:HG22	1.78	0.66
1:G:42:VAL:HG23	1:G:44:VAL:HG12	1.78	0.66
1:G:155:LYS:O	1:G:175:LEU:HA	1.96	0.65
4:L:32:TYR:HE1	4:L:51:PHE:HE2	1.39	0.65
2:B:602:LEU:HB3	2:B:603:ILE:HD12	1.79	0.65
4:L:166:LYS:HD2	4:L:170:ASN:HA	1.79	0.65
3:H:100(I):ASN:HA	4:L:91:PHE:HE1	1.61	0.65
4:L:51:PHE:CE1	4:L:52:ASN:ND2	2.66	0.64
4:L:180:LEU:HD12	4:L:180:LEU:C	2.17	0.64
4:L:130:ALA:HB3	4:L:180:LEU:O	1.97	0.64
2:B:651:ASN:HA	2:B:654:GLU:HB2	1.80	0.64
1:G:101:VAL:HB	1:G:483:LEU:CD1	2.14	0.63
1:G:122:LEU:HB3	1:G:125:LEU:HD12	1.80	0.63
1:G:232:LYS:CE	1:G:269:GLU:HB2	2.28	0.63
4:L:162:THR:HG22	4:L:163:THR:N	2.14	0.63
1:G:41:GLY:N	1:G:493:PRO:O	2.30	0.63
3:H:126:PRO:HD3	3:H:138:LEU:HB3	1.80	0.63
1:G:98:ASN:OD1	1:G:99:ASN:N	2.28	0.63
1:G:307:ILE:HB	1:G:317:PHE:HB3	1.80	0.63
2:B:593:LEU:HG	2:B:599:SER:HA	1.81	0.63
1:G:300:ASN:ND2	1:G:326:ILE:HG23	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:35:HIS:CD2	3:H:50:TRP:HB3	2.34	0.62
1:G:43:PRO:CD	2:B:628:TRP:CD2	2.81	0.62
2:B:543:ASN:HA	2:B:546:SER:HB3	1.80	0.62
1:G:54:CYS:HB2	1:G:215:ILE:HD11	1.81	0.62
4:L:136:ILE:HG22	4:L:139:PHE:CD2	2.34	0.62
1:G:344:LYS:O	1:G:348:GLN:HG2	2.00	0.62
1:G:427:TRP:C	1:G:429:ARG:H	2.04	0.61
1:G:57:ASP:OD1	1:G:58:ALA:N	2.30	0.61
1:G:224:ALA:HA	1:G:246:GLN:OE1	2.01	0.61
3:H:195:ILE:HG22	3:H:210:ARG:HA	1.80	0.61
3:H:3:GLY:HA2	3:H:102:LEU:HD11	1.80	0.61
3:H:159:LEU:HD11	3:H:161:SER:HB2	1.83	0.61
1:G:43:PRO:CD	2:B:628:TRP:CD1	2.84	0.61
1:G:374:HIS:HB3	1:G:385:CYS:HB2	1.82	0.61
1:G:43:PRO:HG2	2:B:628:TRP:HB2	1.83	0.61
3:H:11:MET:HA	3:H:110:THR:O	2.00	0.61
4:L:38:GLN:HB3	4:L:85:HIS:HB2	1.82	0.61
4:L:68:GLY:O	4:L:70:THR:N	2.34	0.61
1:G:396:ILE:HD12	1:G:397:SER:N	2.16	0.60
3:H:50:TRP:NE1	3:H:58:ASN:HB3	2.16	0.60
1:G:317:PHE:HE2	1:G:319:ALA:HB2	1.64	0.60
2:B:643:TYR:O	2:B:647:GLU:HG2	2.01	0.60
2:B:644:GLY:HA2	2:B:647:GLU:OE2	2.02	0.60
3:H:51:MET:HE1	3:H:71:ARG:HB3	1.84	0.60
1:G:369:LEU:HA	1:G:372:THR:HG22	1.84	0.60
2:B:663:LEU:HD12	2:B:664:ASP:HB3	1.83	0.60
1:G:160:ASN:HD22	6:C:1:NAG:H83	1.64	0.60
2:B:598:CYS:O	2:B:600:GLY:N	2.33	0.59
2:B:616:ASN:O	2:B:617:ARG:NH1	2.30	0.59
3:H:100(K):PHE:CD2	3:H:101:ASP:HB2	2.37	0.59
4:L:140:TYR:HD1	4:L:172:TYR:HE2	1.47	0.59
1:G:325:ASP:OD1	1:G:326:ILE:N	2.36	0.59
3:H:65:GLY:O	3:H:82(A):ARG:NH2	2.35	0.59
1:G:221:ALA:C	1:G:223:PHE:H	2.06	0.59
3:H:52(A):PRO:HA	3:H:71:ARG:HD2	1.85	0.59
3:H:152:VAL:HG11	3:H:180:SER:HB2	1.84	0.59
3:H:100(I):ASN:HA	4:L:91:PHE:CE1	2.37	0.59
3:H:59:TYR:HE1	3:H:69:MET:HG2	1.68	0.58
4:L:120:PRO:HD3	4:L:132:LEU:HD21	1.83	0.58
1:G:91:GLU:HG3	1:G:92:GLU:H	1.68	0.58
1:G:93:PHE:HB2	1:G:233:PHE:CZ	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:84:ILE:HB	1:G:244:THR:HG22	1.84	0.58
1:G:480:ARG:HA	1:G:483:LEU:HB2	1.86	0.58
1:G:74:CYS:HB2	2:B:571:TRP:HH2	1.68	0.58
1:G:300:ASN:HD21	1:G:326:ILE:HG23	1.68	0.58
2:B:614:TRP:O	2:B:615:SER:OG	2.20	0.58
1:G:57:ASP:CG	1:G:58:ALA:H	2.08	0.57
1:G:454:LEU:HA	1:G:471:GLY:H	1.68	0.57
1:G:369:LEU:HG	1:G:384:TYR:CD1	2.39	0.57
4:L:6:GLN:OE1	4:L:88:CYS:N	2.33	0.57
1:G:280:ASN:OD1	1:G:281:ALA:N	2.37	0.57
1:G:279:ASN:CG	1:G:282:LYS:HG2	2.25	0.57
3:H:47:TRP:HZ2	3:H:50:TRP:CD1	2.15	0.57
4:L:105:THR:HG21	4:L:141:PRO:CB	2.34	0.57
1:G:86:LEU:HB3	1:G:89:VAL:CG2	2.34	0.57
1:G:119:CYS:SG	1:G:205:CYS:N	2.74	0.57
1:G:305:LYS:HB2	1:G:319:ALA:HB3	1.87	0.57
3:H:143:LYS:HZ1	4:L:129:LYS:HD3	1.70	0.57
2:B:598:CYS:SG	2:B:604:CYS:CB	2.93	0.56
3:H:103:TRP:CD2	4:L:44:PRO:HB2	2.40	0.56
1:G:133:ASN:HD22	1:G:151:ARG:HH22	1.53	0.56
1:G:164:GLU:OE2	1:G:308:ARG:NE	2.34	0.56
4:L:54:ARG:HB2	4:L:58:VAL:HG21	1.86	0.56
1:G:104:MET:HG3	1:G:217:TYR:OH	2.06	0.56
2:B:598:CYS:CB	2:B:604:CYS:SG	2.93	0.56
1:G:138:ILE:O	1:G:139:THR:OG1	2.22	0.56
1:G:294:ILE:HG23	1:G:447:SER:HB2	1.88	0.56
1:G:112:TRP:CZ2	1:G:424:ILE:HD11	2.40	0.56
1:G:177:TYR:CE1	1:G:420:ILE:HB	2.41	0.56
4:L:54:ARG:NE	4:L:60:ASP:HA	2.21	0.56
1:G:175:LEU:HD12	1:G:321:GLY:O	2.06	0.56
1:G:104:MET:SD	1:G:479:TRP:CD1	2.99	0.55
1:G:298:ARG:HE	1:G:443:ILE:HD12	1.71	0.55
1:G:345:VAL:O	1:G:348:GLN:N	2.40	0.55
1:G:360:ARG:HG3	1:G:394:THR:HG22	1.88	0.55
4:L:140:TYR:HB2	4:L:171:LYS:HD2	1.88	0.55
4:L:5:THR:OG1	4:L:24:THR:N	2.40	0.55
4:L:113:PRO:HA	4:L:139:PHE:HB3	1.89	0.55
1:G:138:ILE:HG21	1:G:326:ILE:HB	1.89	0.55
1:G:378:CYS:SG	1:G:379:GLY:N	2.80	0.55
4:L:35:TRP:CZ3	4:L:88:CYS:HB3	2.42	0.55
4:L:113:PRO:HD3	4:L:197:HIS:CD2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:69:TRP:CE3	1:G:71:THR:HG23	2.42	0.54
3:H:32:TYR:CD1	3:H:96:ARG:HA	2.41	0.54
1:G:43:PRO:HG3	2:B:628:TRP:NE1	2.18	0.54
1:G:156:ASN:HA	1:G:175:LEU:HD23	1.89	0.54
2:B:522:PHE:CE1	2:B:543:ASN:HB2	2.41	0.54
1:G:329:ALA:HB2	1:G:420:ILE:HD11	1.88	0.54
3:H:166:PHE:CE1	4:L:173:ALA:HB1	2.42	0.54
1:G:45:TRP:HB3	1:G:491:ILE:HG22	1.89	0.54
1:G:101:VAL:HA	1:G:483:LEU:CD1	2.38	0.54
4:L:52:ASN:OD1	4:L:53:LYS:N	2.41	0.54
1:G:212:PRO:HB2	1:G:252:LYS:HB3	1.88	0.54
1:G:223:PHE:CE1	1:G:490:LYS:HD3	2.42	0.54
9:G:602:NAG:H83	9:G:602:NAG:H3	1.89	0.54
1:G:101:VAL:HG23	1:G:479:TRP:HB2	1.90	0.53
1:G:396:ILE:HD12	1:G:397:SER:H	1.73	0.53
3:H:27:TYR:CD1	3:H:28:ASP:N	2.76	0.53
4:L:4:LEU:HB2	4:L:99:GLY:HA2	1.90	0.53
1:G:122:LEU:HD21	1:G:203:GLN:HB2	1.90	0.53
2:B:611:ASN:HB3	2:B:614:TRP:CD2	2.44	0.53
3:H:29:PHE:CE2	3:H:73:THR:HA	2.43	0.53
4:L:139:PHE:CE1	4:L:142:GLY:HA2	2.43	0.53
1:G:101:VAL:HA	1:G:483:LEU:HD11	1.91	0.53
1:G:136:ASN:OD1	1:G:137:ASN:N	2.41	0.53
1:G:150:MET:SD	1:G:328:GLN:HB3	2.48	0.53
1:G:482:GLU:HA	1:G:484:TYR:CE2	2.44	0.53
4:L:136:ILE:HD11	4:L:195:VAL:HG11	1.90	0.53
1:G:104:MET:SD	1:G:479:TRP:HD1	2.31	0.53
4:L:140:TYR:CD1	4:L:141:PRO:HA	2.43	0.53
1:G:372:THR:HG23	1:G:373:THR:HG23	1.90	0.52
3:H:121:VAL:HG23	3:H:209:LYS:HG3	1.90	0.52
4:L:32:TYR:CD1	4:L:51:PHE:HE2	2.21	0.52
1:G:86:LEU:HB3	1:G:89:VAL:HG21	1.90	0.52
1:G:272:ILE:HG22	1:G:286:VAL:HG12	1.90	0.52
3:H:86:ASP:O	3:H:90:TYR:OH	2.22	0.52
1:G:112:TRP:HZ2	1:G:424:ILE:HD11	1.73	0.52
1:G:154:LEU:HA	1:G:176:PHE:O	2.10	0.52
1:G:397:SER:OG	1:G:398:ASN:N	2.42	0.52
2:B:523:LEU:HA	2:B:540:GLN:HE21	1.75	0.52
4:L:108:GLN:HB3	4:L:109:PRO:CD	2.40	0.52
1:G:317:PHE:CE2	1:G:319:ALA:HB2	2.44	0.52
3:H:38:ARG:HB2	3:H:48:LEU:HD11	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:131:CYS:HA	1:G:157:CYS:HA	1.92	0.52
3:H:196:CYS:O	3:H:208:ASP:HA	2.10	0.52
1:G:428:GLN:O	1:G:430:ILE:N	2.42	0.51
2:B:612:SER:OG	2:B:613:SER:N	2.43	0.51
3:H:154:TRP:H	3:H:157:GLY:HA2	1.76	0.51
4:L:141:PRO:O	4:L:143:ALA:N	2.43	0.51
1:G:86:LEU:HD11	1:G:244:THR:HB	1.92	0.51
2:B:538:THR:HG23	2:B:539:VAL:HG23	1.91	0.51
1:G:385:CYS:HB3	1:G:416:LEU:CD1	2.40	0.51
1:G:129:LEU:O	1:G:190:GLU:HA	2.10	0.51
1:G:480:ARG:HH22	3:H:100(D):ARG:NH2	2.09	0.51
4:L:83:GLU:OE1	4:L:106:VAL:HG23	2.10	0.51
4:L:178:LEU:CD2	4:L:180:LEU:HD23	2.41	0.51
1:G:335:LYS:HZ3	1:G:411:ASN:HA	1.75	0.51
4:L:180:LEU:CD1	4:L:185:TRP:CB	2.84	0.51
1:G:215:ILE:HG22	1:G:251:ILE:O	2.11	0.51
3:H:20:VAL:CG1	3:H:80:LEU:HB3	2.40	0.51
4:L:130:ALA:HB3	4:L:180:LEU:HD12	1.93	0.51
1:G:258:GLN:HG2	1:G:470:PRO:HB2	1.93	0.51
3:H:37:PHE:HA	3:H:48:LEU:HD13	1.92	0.51
4:L:61:ARG:HG3	4:L:62:PHE:CD1	2.46	0.51
1:G:154:LEU:HD23	1:G:177:TYR:HA	1.93	0.50
2:B:646:LEU:O	2:B:650:GLN:HB2	2.12	0.50
4:L:189:ARG:HA	4:L:208:PRO:CD	2.42	0.50
1:G:299:PRO:HD3	1:G:329:ALA:HA	1.92	0.50
3:H:18:VAL:HG22	3:H:19:ARG:H	1.76	0.50
1:G:197:ASN:HD22	7:D:1:NAG:H83	1.76	0.50
1:G:467:THR:HG23	3:H:61:ARG:HH21	1.76	0.50
1:G:119:CYS:HB3	1:G:203:GLN:O	2.12	0.50
2:B:631:TRP:NE1	2:B:635:ILE:HG13	2.26	0.50
6:C:2:NAG:H83	6:C:2:NAG:H3	1.92	0.50
3:H:33:VAL:CG1	3:H:95:ASP:HB3	2.42	0.49
3:H:52:ASN:HD21	3:H:53:SER:HB3	1.77	0.49
3:H:103:TRP:HB3	4:L:44:PRO:HD2	1.94	0.49
3:H:164:HIS:CD2	4:L:167:GLN:HG2	2.47	0.49
3:H:200:HIS:HD2	3:H:202:PRO:HD2	1.77	0.49
4:L:6:GLN:HB3	4:L:102:THR:HG23	1.94	0.49
1:G:38:VAL:HA	1:G:496:VAL:HA	1.95	0.49
1:G:279:ASN:HA	3:H:100(J):TRP:HZ2	1.77	0.49
3:H:198:VAL:CG2	3:H:207:VAL:HB	2.42	0.49
4:L:18:SER:HA	4:L:75:ILE:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:67:SER:OG	4:L:70:THR:OG1	2.31	0.49
3:H:41:ARG:HH22	3:H:170:LEU:H	1.59	0.49
3:H:166:PHE:HE1	4:L:173:ALA:HB1	1.77	0.49
2:B:607:ASN:HB2	2:B:650:GLN:HE22	1.76	0.49
3:H:36:TRP:CD1	3:H:69:MET:HE3	2.47	0.49
3:H:38:ARG:CB	3:H:48:LEU:HD11	2.43	0.49
1:G:184:ILE:CD1	1:G:190:GLU:HG2	2.20	0.49
1:G:293:GLN:O	1:G:333:VAL:HG23	2.13	0.48
2:B:650:GLN:NE2	2:B:654:GLU:OE2	2.46	0.48
1:G:363:ASN:ND2	5:M:1:NAG:O7	2.46	0.48
1:G:422:GLN:HA	1:G:436:ALA:HB3	1.96	0.48
2:B:640:GLN:HG2	2:B:641:ILE:N	2.28	0.48
3:H:127:SER:OG	3:H:128:SER:N	2.45	0.48
1:G:297:THR:O	1:G:329:ALA:HA	2.14	0.48
1:G:36:VAL:HG23	2:B:606:THR:HB	1.95	0.48
1:G:131:CYS:SG	1:G:191:TYR:HD2	2.37	0.48
3:H:184:VAL:HG11	3:H:194:TYR:CE2	2.49	0.48
4:L:33:VAL:CG2	4:L:88:CYS:HB2	2.44	0.48
1:G:416:LEU:HD12	1:G:416:LEU:O	2.13	0.48
1:G:170:GLN:HG2	1:G:172:VAL:HG13	1.96	0.48
1:G:368:ASP:OD2	3:H:71:ARG:NH2	2.42	0.48
3:H:95:ASP:OD2	3:H:98:ASN:N	2.47	0.48
1:G:285:LEU:HD12	1:G:477:ASP:HB3	1.95	0.48
3:H:12:LYS:HD3	3:H:18:VAL:HB	1.96	0.48
4:L:105:THR:HG21	4:L:141:PRO:HB2	1.96	0.48
1:G:381:GLU:HG3	1:G:439:ILE:CD1	2.44	0.47
4:L:118:PHE:HB2	4:L:133:VAL:CG2	2.44	0.47
1:G:257:THR:O	1:G:259:LEU:N	2.41	0.47
4:L:61:ARG:NH2	4:L:77:GLY:O	2.47	0.47
1:G:50:THR:HB	1:G:223:PHE:CE1	2.50	0.47
1:G:231:LYS:HD3	1:G:267:GLU:O	2.13	0.47
2:B:631:TRP:O	2:B:635:ILE:HG12	2.14	0.47
1:G:105:HIS:O	1:G:108:ILE:HG13	2.15	0.47
3:H:34:ILE:HG22	3:H:51:MET:O	2.15	0.47
4:L:151:ASP:HA	4:L:190:SER:HB3	1.95	0.47
2:B:598:CYS:SG	2:B:604:CYS:HB3	2.55	0.47
3:H:121:VAL:CG2	3:H:209:LYS:HG3	2.44	0.47
1:G:101:VAL:CA	1:G:483:LEU:HD11	2.43	0.47
1:G:456:ARG:HG3	1:G:468:PHE:CE1	2.43	0.47
2:B:618:ASN:HB2	2:B:621:GLU:HG2	1.97	0.47
1:G:225:ILE:CG2	1:G:486:TYR:HD1	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:130:ALA:HB3	4:L:180:LEU:CD1	2.44	0.47
1:G:46:LYS:HG3	1:G:47:ASP:N	2.30	0.46
1:G:69:TRP:HE3	1:G:71:THR:H	1.64	0.46
1:G:85:HIS:CD2	1:G:86:LEU:H	2.33	0.46
1:G:110:SER:O	1:G:114:GLN:HG2	2.15	0.46
2:B:618:ASN:ND2	9:B:702:NAG:O7	2.48	0.46
1:G:100:MET:HG3	1:G:483:LEU:HD13	1.98	0.46
1:G:341:THR:O	1:G:345:VAL:HG23	2.16	0.46
1:G:75:VAL:HG13	1:G:76:PRO:HD2	1.97	0.46
1:G:360:ARG:HB3	1:G:467:THR:CG2	2.44	0.46
1:G:475:MET:HG3	1:G:479:TRP:CZ3	2.49	0.46
3:H:20:VAL:HG12	3:H:80:LEU:HB3	1.96	0.46
4:L:37:GLN:O	4:L:44:PRO:HA	2.16	0.46
1:G:482:GLU:HA	1:G:484:TYR:HE2	1.81	0.46
3:H:84:TYR:HA	3:H:111:VAL:CG1	2.45	0.46
1:G:125:LEU:O	1:G:127:VAL:HG12	2.16	0.46
1:G:279:ASN:ND2	1:G:282:LYS:HE2	2.31	0.46
3:H:122:PHE:CD2	4:L:124:GLU:HG2	2.50	0.46
4:L:180:LEU:CD1	4:L:185:TRP:HB2	2.37	0.46
1:G:424:ILE:HB	1:G:426:MET:HG3	1.98	0.46
3:H:47:TRP:CZ2	3:H:50:TRP:CD1	2.94	0.46
1:G:95:MET:SD	1:G:273:ARG:HD3	2.56	0.46
1:G:99:ASN:O	1:G:102:GLU:N	2.49	0.46
1:G:330:HIS:CB	1:G:417:PRO:HA	2.46	0.46
1:G:427:TRP:O	1:G:429:ARG:N	2.49	0.46
1:G:233:PHE:C	1:G:235:GLY:H	2.19	0.46
1:G:377:ASN:HD21	1:G:380:GLY:C	2.19	0.46
1:G:429:ARG:HB3	1:G:432:GLN:NE2	2.30	0.46
1:G:96:TRP:HD1	1:G:480:ARG:NH1	2.14	0.45
3:H:88:ALA:O	3:H:108:GLN:HG3	2.17	0.45
1:G:225:ILE:HG21	1:G:486:TYR:HD1	1.81	0.45
4:L:13:ALA:HA	4:L:106(A):LEU:HG	1.98	0.45
1:G:276:ASN:OD1	9:G:602:NAG:H82	2.16	0.45
3:H:191:THR:HG23	3:H:192:GLN:N	2.23	0.45
4:L:33:VAL:HA	4:L:89:TRP:O	2.16	0.45
1:G:64:GLU:O	1:G:66:HIS:N	2.50	0.45
1:G:163:THR:HG21	1:G:168:LYS:HB3	1.99	0.45
2:B:595:ILE:HD12	2:B:596:TRP:CD1	2.51	0.45
3:H:87:THR:HG21	3:H:170:LEU:HD23	1.99	0.45
4:L:178:LEU:HD22	4:L:180:LEU:HD23	1.97	0.45
1:G:280:ASN:HB3	1:G:456:ARG:NH2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:596:TRP:C	2:B:651:ASN:HD21	2.19	0.45
1:G:96:TRP:CD1	1:G:480:ARG:NH1	2.85	0.45
1:G:335:LYS:NZ	1:G:411:ASN:HA	2.32	0.45
1:G:276:ASN:O	1:G:279:ASN:HB3	2.16	0.45
1:G:222:GLY:O	1:G:491:ILE:HG13	2.17	0.45
3:H:36:TRP:CH2	3:H:92:CYS:HB3	2.52	0.45
1:G:55:ALA:HB3	1:G:216:HIS:HB2	1.99	0.44
1:G:122:LEU:HD12	1:G:201:ILE:HG12	1.99	0.44
1:G:373:THR:HB	1:G:385:CYS:O	2.17	0.44
1:G:476:ARG:NH2	3:H:100:SER:HA	2.32	0.44
3:H:59:TYR:CE1	3:H:69:MET:HG2	2.51	0.44
3:H:126:PRO:HD2	3:H:213:PRO:HB3	1.98	0.44
4:L:66:THR:HA	4:L:71:ALA:HA	1.98	0.44
1:G:52:LEU:CD2	1:G:217:TYR:HB3	2.48	0.44
1:G:313:PRO:O	1:G:315:GLN:N	2.50	0.44
3:H:40:ARG:NE	3:H:85:ASP:O	2.47	0.44
3:H:119:PRO:HD3	3:H:200:HIS:ND1	2.32	0.44
4:L:134:CYS:C	4:L:135:LEU:HD12	2.38	0.44
1:G:302:ASN:HB2	1:G:440:GLN:HA	2.00	0.44
1:G:338:TRP:CZ2	1:G:390:LEU:HD13	2.52	0.44
1:G:333:VAL:CG1	1:G:414:ILE:HB	2.48	0.44
1:G:92:GLU:HG3	1:G:237:GLY:C	2.38	0.44
3:H:4:LEU:HD13	3:H:92:CYS:SG	2.57	0.44
1:G:223:PHE:CZ	1:G:490:LYS:HD3	2.53	0.44
4:L:12:SER:HA	4:L:105:THR:O	2.18	0.44
4:L:207:ALA:O	4:L:209:THR:HG22	2.18	0.44
1:G:229:LYS:N	1:G:241:SER:O	2.32	0.44
3:H:18:VAL:HG22	3:H:19:ARG:N	2.33	0.44
3:H:201:LYS:N	3:H:202:PRO:CD	2.81	0.44
4:L:14:SER:N	4:L:106(A):LEU:HB2	2.33	0.44
4:L:148:TRP:HB2	4:L:155:VAL:HG22	2.00	0.44
4:L:180:LEU:HD12	4:L:181:THR:N	2.33	0.44
4:L:189:ARG:HA	4:L:208:PRO:HD2	1.98	0.44
1:G:46:LYS:HG3	1:G:47:ASP:H	1.82	0.44
1:G:116:LEU:HA	1:G:119:CYS:SG	2.58	0.44
1:G:153:GLU:OE1	1:G:154:LEU:HG	2.18	0.44
4:L:151:ASP:HB2	4:L:188:HIS:O	2.18	0.44
1:G:435:TYR:CE2	1:G:437:PRO:HA	2.53	0.43
3:H:12:LYS:HE2	3:H:17:SER:O	2.17	0.43
3:H:39:GLN:HB2	3:H:45:LEU:HG	2.00	0.43
4:L:3:ALA:N	4:L:97:ASN:OD1	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:36:TRP:NE1	3:H:69:MET:HE3	2.33	0.43
3:H:146:PHE:HB2	3:H:175:LEU:HD22	2.00	0.43
4:L:14:SER:HB2	4:L:17:GLN:NE2	2.33	0.43
4:L:111:ALA:HB3	4:L:140:TYR:N	2.32	0.43
1:G:52:LEU:HD21	1:G:217:TYR:HB3	1.99	0.43
1:G:133:ASN:OD1	1:G:155:LYS:NZ	2.42	0.43
1:G:427:TRP:C	1:G:429:ARG:N	2.70	0.43
1:G:458:GLY:O	1:G:460:SER:N	2.51	0.43
2:B:597:GLY:N	2:B:651:ASN:HD21	2.17	0.43
1:G:130:GLN:O	1:G:158:SER:N	2.51	0.43
1:G:46:LYS:CG	1:G:47:ASP:H	2.31	0.43
1:G:305:LYS:O	1:G:318:TYR:HA	2.18	0.43
3:H:153:SER:HB3	3:H:157:GLY:N	2.34	0.43
4:L:82:ASP:HB3	4:L:86:TYR:OH	2.18	0.43
4:L:166:LYS:CD	4:L:170:ASN:HA	2.45	0.43
1:G:101:VAL:HG21	1:G:480:ARG:HG3	2.00	0.43
3:H:33:VAL:HG13	3:H:95:ASP:HB3	2.01	0.43
1:G:122:LEU:HD22	1:G:125:LEU:HD11	1.99	0.43
4:L:11:VAL:HG21	4:L:21:ILE:CD1	2.47	0.43
4:L:103:LYS:NZ	4:L:143:ALA:HB2	2.34	0.43
4:L:108:GLN:HB3	4:L:109:PRO:HD2	2.00	0.43
1:G:341:THR:HA	1:G:344:LYS:HG2	2.01	0.43
3:H:12:LYS:CD	3:H:18:VAL:HB	2.49	0.43
1:G:343:GLY:O	1:G:346:VAL:HG12	2.18	0.43
4:L:133:VAL:HG12	4:L:177:TYR:CD1	2.54	0.43
1:G:35:TRP:CZ3	2:B:609:PRO:HD3	2.53	0.42
3:H:200:HIS:CD2	3:H:202:PRO:HD2	2.53	0.42
1:G:108:ILE:HA	1:G:111:LEU:HG	2.00	0.42
3:H:119:PRO:CB	3:H:145:TYR:HB3	2.42	0.42
1:G:279:ASN:ND2	1:G:282:LYS:HG2	2.34	0.42
1:G:329:ALA:CB	1:G:420:ILE:HD11	2.49	0.42
3:H:122:PHE:CE2	4:L:124:GLU:HG2	2.55	0.42
3:H:168:ALA:HA	3:H:178:LEU:HB3	2.02	0.42
1:G:62:GLU:O	1:G:63:THR:OG1	2.30	0.42
4:L:26:THR:HG22	4:L:27:HIS:N	2.34	0.42
4:L:119:PRO:HB2	4:L:120:PRO:HD2	2.02	0.42
4:L:6:GLN:HG3	4:L:23:CYS:HB2	2.01	0.42
1:G:231:LYS:HG3	1:G:232:LYS:HZ3	1.84	0.42
1:G:290:THR:OG1	1:G:344:LYS:NZ	2.50	0.42
1:G:309:ILE:HD12	1:G:309:ILE:O	2.19	0.42
3:H:139:GLY:HA2	3:H:154:TRP:HH2	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:140:TYR:HB2	4:L:171:LYS:CD	2.49	0.42
4:L:189:ARG:O	4:L:207:ALA:HA	2.19	0.42
1:G:161:MET:HB3	1:G:170:GLN:HB3	2.01	0.42
1:G:453:ILE:HG21	1:G:472:GLY:HA2	2.01	0.42
2:B:609:PRO:HB2	2:B:610:TRP:H	1.52	0.42
3:H:4:LEU:HD23	3:H:24:ALA:HA	2.01	0.42
3:H:100(L):LEU:O	3:H:103:TRP:NE1	2.52	0.42
4:L:183:GLU:HG2	4:L:184:GLN:N	2.35	0.42
1:G:363:ASN:ND2	5:M:1:NAG:C7	2.83	0.42
3:H:37:PHE:CA	3:H:48:LEU:HD13	2.50	0.42
1:G:72:HIS:CE1	2:B:571:TRP:HE3	2.37	0.41
1:G:101:VAL:HG23	1:G:479:TRP:CB	2.49	0.41
3:H:36:TRP:HE1	3:H:69:MET:HE3	1.85	0.41
3:H:59:TYR:HB2	3:H:64:GLN:HG2	2.02	0.41
4:L:6:GLN:HA	4:L:22:SER:O	2.20	0.41
4:L:9:SER:HB3	4:L:143:ALA:HB1	2.01	0.41
2:B:608:VAL:HG13	2:B:650:GLN:OE1	2.20	0.41
4:L:36:CYS:SG	4:L:37:GLN:N	2.93	0.41
1:G:329:ALA:HB1	1:G:383:PHE:HE1	1.85	0.41
4:L:13:ALA:CA	4:L:106(A):LEU:HG	2.51	0.41
4:L:37:GLN:HB2	4:L:47:LEU:HD11	2.02	0.41
1:G:93:PHE:HB2	1:G:233:PHE:HZ	1.81	0.41
1:G:101:VAL:N	1:G:483:LEU:HD11	2.35	0.41
2:B:639:THR:HA	2:B:642:ILE:HD12	2.02	0.41
3:H:72:ASP:HB3	3:H:75:THR:OG1	2.20	0.41
1:G:288:PHE:HB3	1:G:290:THR:H	1.85	0.41
2:B:598:CYS:CB	2:B:604:CYS:HG	2.20	0.41
3:H:50:TRP:HE1	3:H:58:ASN:HB3	1.85	0.41
1:G:225:ILE:HG21	1:G:486:TYR:CD1	2.54	0.41
3:H:51:MET:CE	3:H:71:ARG:HB3	2.50	0.41
4:L:61:ARG:HH22	4:L:79:GLN:HB2	1.86	0.41
1:G:55:ALA:N	1:G:216:HIS:O	2.54	0.41
1:G:85:HIS:CD2	1:G:86:LEU:N	2.89	0.41
1:G:179:LEU:HD23	1:G:369:LEU:HD22	2.03	0.41
1:G:377:ASN:ND2	1:G:381:GLU:O	2.54	0.41
4:L:148:TRP:CZ3	4:L:193:CYS:HB2	2.56	0.41
1:G:43:PRO:CG	2:B:628:TRP:NE1	2.76	0.41
1:G:72:HIS:CE1	2:B:571:TRP:CE3	3.08	0.41
1:G:93:PHE:CE2	1:G:239:CYS:HB3	2.55	0.41
1:G:129:LEU:HD11	1:G:193:LEU:HD12	2.03	0.41
1:G:192:ARG:HG2	1:G:193:LEU:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:332:ASN:ND2	8:K:1:NAG:O7	2.54	0.41
1:G:415:THR:HG21	8:K:1:NAG:H62	2.01	0.41
3:H:111:VAL:HG13	3:H:111:VAL:O	2.21	0.41
4:L:12:SER:HB2	4:L:106(A):LEU:HD23	2.03	0.41
4:L:16:GLY:HA2	4:L:77:GLY:HA2	2.03	0.41
1:G:233:PHE:O	1:G:235:GLY:N	2.50	0.41
2:B:569:THR:HG23	2:B:572:GLY:HA3	2.03	0.41
2:B:570:VAL:C	2:B:572:GLY:H	2.25	0.41
2:B:627:THR:N	2:B:630:GLN:OE1	2.54	0.41
1:G:96:TRP:C	1:G:98:ASN:H	2.25	0.40
1:G:221:ALA:C	1:G:223:PHE:N	2.74	0.40
4:L:169:ASN:O	4:L:171:LYS:N	2.42	0.40
2:B:634:GLU:HA	9:B:703:NAG:H82	2.02	0.40
1:G:48:ALA:HB3	1:G:490:LYS:HB2	2.03	0.40
1:G:50:THR:HB	1:G:223:PHE:CZ	2.56	0.40
1:G:52:LEU:HG	1:G:219:ALA:HA	2.04	0.40
1:G:359:ILE:HB	1:G:395:TRP:HB2	2.04	0.40
3:H:126:PRO:O	3:H:127:SER:HB2	2.22	0.40
1:G:90:THR:HG22	1:G:240:PRO:CA	2.44	0.40
1:G:229:LYS:HE2	1:G:242:VAL:C	2.42	0.40
1:G:335:LYS:HG2	1:G:414:ILE:HG13	2.03	0.40
1:G:335:LYS:HZ2	1:G:412:ASP:H	1.69	0.40
4:L:185:TRP:HH2	4:L:206:VAL:HG12	1.86	0.40
4:L:189:ARG:HD3	4:L:207:ALA:HB1	2.03	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	G	443/481 (92%)	351 (79%)	61 (14%)	31 (7%)	1 16

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	118/153 (77%)	92 (78%)	21 (18%)	5 (4%)	3 25
3	H	221/229 (96%)	190 (86%)	22 (10%)	9 (4%)	3 25
4	L	198/205 (97%)	163 (82%)	23 (12%)	12 (6%)	1 18
All	All	980/1068 (92%)	796 (81%)	127 (13%)	57 (6%)	1 19

All (57) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	65	LYS
1	G	66	HIS
1	G	268	GLU
1	G	429	ARG
2	B	598	CYS
3	H	53	SER
3	H	191	THR
4	L	51	PHE
4	L	55	PRO
1	G	69	TRP
1	G	70	ALA
1	G	71	THR
1	G	115	SER
1	G	153	GLU
1	G	279	ASN
1	G	314	GLY
1	G	396	ILE
1	G	459	GLY
1	G	460	SER
1	G	463	SER
2	B	609	PRO
2	B	614	TRP
3	H	127	SER
4	L	5	THR
4	L	69	ASN
4	L	77	GLY
4	L	110	LYS
4	L	142	GLY
4	L	189	ARG
1	G	35	TRP
1	G	126	CYS
1	G	258	GLN
1	G	464	THR

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Mol	Chain	Res	Type
3	H	144	ASP
3	H	173	SER
1	G	47	ASP
1	G	114	GLN
1	G	262	ASN
1	G	428	GLN
2	B	599	SER
3	H	27	TYR
3	H	100(F)	GLU
3	H	114	ALA
4	L	151	ASP
1	G	97	LYS
1	G	188	ASN
1	G	234	ASN
1	G	500	ARG
2	B	650	GLN
3	H	117	LYS
4	L	6	GLN
4	L	25	GLY
4	L	170	ASN
1	G	152	GLY
1	G	276	ASN
1	G	184	ILE
1	G	299	PRO

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	G	401/428 (94%)	399 (100%)	2 (0%)	88 93
2	B	105/129 (81%)	105 (100%)	0	100 100
3	H	188/192 (98%)	184 (98%)	4 (2%)	53 72
4	L	168/173 (97%)	166 (99%)	2 (1%)	71 83
All	All	862/922 (94%)	854 (99%)	8 (1%)	78 88

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	377	ASN
1	G	378	CYS
3	H	22	CYS
3	H	100	SER
3	H	100(K)	PHE
3	H	164	HIS
4	L	8	ARG
4	L	23	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	85	HIS
1	G	287	GLN
1	G	377	ASN
1	G	478	ASN
2	B	650	GLN
2	B	651	ASN
3	H	6	GLN
3	H	35	HIS
3	H	58	ASN
3	H	197	ASN
4	L	85	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

34 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	$\#\|Z\| > 2$
5	NAG	A	1	5,1	14,14,15	0.40	0	17,19,21	0.37	0
5	NAG	A	2	5	14,14,15	0.49	0	17,19,21	0.68	1 (5%)
5	BMA	A	3	5	11,11,12	0.67	0	15,15,17	0.94	1 (6%)
6	NAG	C	1	1,6	14,14,15	0.25	0	17,19,21	0.45	0
6	NAG	C	2	6	14,14,15	0.35	0	17,19,21	1.39	2 (11%)
7	NAG	D	1	1,7	14,14,15	0.44	0	17,19,21	0.69	0
7	NAG	D	2	7	14,14,15	0.19	0	17,19,21	0.70	0
7	BMA	D	3	7	11,11,12	1.10	1 (9%)	15,15,17	1.11	1 (6%)
7	MAN	D	4	7	11,11,12	1.06	1 (9%)	15,15,17	1.46	3 (20%)
6	NAG	E	1	1,6	14,14,15	0.39	0	17,19,21	0.47	0
6	NAG	E	2	6	14,14,15	0.29	0	17,19,21	0.44	0
5	NAG	F	1	5,1	14,14,15	0.31	0	17,19,21	0.36	0
5	NAG	F	2	5	14,14,15	0.18	0	17,19,21	0.40	0
5	BMA	F	3	5	11,11,12	0.80	0	15,15,17	0.88	0
6	NAG	I	1	1,6	14,14,15	0.20	0	17,19,21	0.72	1 (5%)
6	NAG	I	2	6	14,14,15	0.30	0	17,19,21	0.41	0
6	NAG	J	1	1,6	14,14,15	0.22	0	17,19,21	0.45	0
6	NAG	J	2	6	14,14,15	0.32	0	17,19,21	0.51	0
8	NAG	K	1	8,1	14,14,15	0.46	0	17,19,21	0.50	0
8	NAG	K	2	8	14,14,15	0.26	0	17,19,21	0.59	0
8	BMA	K	3	8	11,11,12	0.70	0	15,15,17	0.94	0
8	MAN	K	4	8	11,11,12	0.72	0	15,15,17	1.31	2 (13%)
8	MAN	K	5	8	11,11,12	0.89	0	15,15,17	1.15	3 (20%)
8	MAN	K	6	8	11,11,12	0.82	0	15,15,17	1.29	2 (13%)
8	MAN	K	7	8	11,11,12	0.82	0	15,15,17	1.12	2 (13%)
5	NAG	M	1	5,1	14,14,15	0.31	0	17,19,21	0.54	0
5	NAG	M	2	5	14,14,15	0.42	0	17,19,21	0.65	0
5	BMA	M	3	5	11,11,12	0.92	1 (9%)	15,15,17	1.25	1 (6%)
6	NAG	N	1	1,6	14,14,15	0.43	0	17,19,21	0.75	0
6	NAG	N	2	6	14,14,15	0.29	0	17,19,21	0.41	0
6	NAG	O	1	1,6	14,14,15	0.36	0	17,19,21	0.68	0
6	NAG	O	2	6	14,14,15	0.27	0	17,19,21	0.49	0
6	NAG	P	1	1,6	14,14,15	0.45	0	17,19,21	0.56	0
6	NAG	P	2	6	14,14,15	0.24	0	17,19,21	0.51	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	A	2	5	-	1/6/23/26	0/1/1/1
5	BMA	A	3	5	-	1/2/19/22	0/1/1/1
6	NAG	C	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	C	2	6	-	5/6/23/26	0/1/1/1
7	NAG	D	1	1,7	-	3/6/23/26	0/1/1/1
7	NAG	D	2	7	-	3/6/23/26	0/1/1/1
7	BMA	D	3	7	-	2/2/19/22	0/1/1/1
7	MAN	D	4	7	-	0/2/19/22	0/1/1/1
6	NAG	E	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	E	2	6	-	3/6/23/26	0/1/1/1
5	NAG	F	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	F	2	5	-	1/6/23/26	0/1/1/1
5	BMA	F	3	5	-	2/2/19/22	0/1/1/1
6	NAG	I	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	I	2	6	-	1/6/23/26	0/1/1/1
6	NAG	J	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	J	2	6	-	2/6/23/26	0/1/1/1
8	NAG	K	1	8,1	-	0/6/23/26	0/1/1/1
8	NAG	K	2	8	-	3/6/23/26	0/1/1/1
8	BMA	K	3	8	-	0/2/19/22	0/1/1/1
8	MAN	K	4	8	-	1/2/19/22	0/1/1/1
8	MAN	K	5	8	-	0/2/19/22	0/1/1/1
8	MAN	K	6	8	-	2/2/19/22	1/1/1/1
8	MAN	K	7	8	-	0/2/19/22	0/1/1/1
5	NAG	M	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	M	2	5	-	3/6/23/26	0/1/1/1
5	BMA	M	3	5	-	0/2/19/22	0/1/1/1
6	NAG	N	1	1,6	-	3/6/23/26	0/1/1/1
6	NAG	N	2	6	-	2/6/23/26	0/1/1/1
6	NAG	O	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	O	2	6	-	2/6/23/26	0/1/1/1
6	NAG	P	1	1,6	-	4/6/23/26	0/1/1/1
6	NAG	P	2	6	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	3	BMA	O5-C1	-2.75	1.39	1.43
5	M	3	BMA	C1-C2	2.17	1.57	1.52
7	D	4	MAN	C2-C3	2.14	1.55	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	C	2	NAG	C2-N2-C7	4.56	129.39	122.90
8	K	6	MAN	C1-O5-C5	3.69	117.20	112.19
8	K	4	MAN	C1-O5-C5	3.04	116.32	112.19
8	K	7	MAN	C1-O5-C5	2.73	115.89	112.19
7	D	4	MAN	C1-O5-C5	2.64	115.76	112.19
7	D	4	MAN	O5-C1-C2	2.63	114.83	110.77
7	D	4	MAN	C1-C2-C3	2.47	112.70	109.67
8	K	7	MAN	O2-C2-C3	-2.44	105.25	110.14
6	C	2	NAG	C1-C2-N2	2.43	114.64	110.49
8	K	5	MAN	C1-O5-C5	2.37	115.40	112.19
8	K	6	MAN	O2-C2-C3	-2.36	105.40	110.14
5	M	3	BMA	O5-C1-C2	2.34	114.38	110.77
6	I	1	NAG	C1-O5-C5	2.29	115.30	112.19
7	D	3	BMA	C3-C4-C5	2.27	114.30	110.24
8	K	5	MAN	O2-C2-C3	-2.15	105.82	110.14
5	A	3	BMA	C1-O5-C5	2.13	115.08	112.19
8	K	4	MAN	O2-C2-C3	-2.10	105.92	110.14
8	K	5	MAN	O5-C1-C2	2.08	113.98	110.77
5	A	2	NAG	C1-O5-C5	2.07	115.00	112.19

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	M	1	NAG	O5-C5-C6-O6
6	E	1	NAG	O5-C5-C6-O6
8	K	2	NAG	O5-C5-C6-O6
5	A	1	NAG	O5-C5-C6-O6
5	F	3	BMA	O5-C5-C6-O6
6	N	1	NAG	O5-C5-C6-O6
6	E	1	NAG	C4-C5-C6-O6
6	C	2	NAG	O5-C5-C6-O6
6	O	2	NAG	O5-C5-C6-O6
6	P	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	D	3	BMA	O5-C5-C6-O6
6	E	2	NAG	O5-C5-C6-O6
5	M	1	NAG	C4-C5-C6-O6
6	N	1	NAG	C4-C5-C6-O6
5	A	1	NAG	C4-C5-C6-O6
8	K	2	NAG	C4-C5-C6-O6
6	P	2	NAG	O5-C5-C6-O6
5	F	1	NAG	C4-C5-C6-O6
6	O	2	NAG	C4-C5-C6-O6
7	D	3	BMA	C4-C5-C6-O6
5	F	3	BMA	C4-C5-C6-O6
6	P	2	NAG	C4-C5-C6-O6
8	K	6	MAN	O5-C5-C6-O6
6	E	2	NAG	C4-C5-C6-O6
6	C	1	NAG	C8-C7-N2-C2
6	C	1	NAG	O7-C7-N2-C2
6	C	2	NAG	C8-C7-N2-C2
6	C	2	NAG	O7-C7-N2-C2
6	P	1	NAG	C8-C7-N2-C2
6	P	1	NAG	O7-C7-N2-C2
7	D	1	NAG	C8-C7-N2-C2
7	D	1	NAG	O7-C7-N2-C2
6	O	1	NAG	C4-C5-C6-O6
6	P	1	NAG	C4-C5-C6-O6
6	C	2	NAG	C4-C5-C6-O6
5	F	1	NAG	O5-C5-C6-O6
5	F	1	NAG	C1-C2-N2-C7
6	N	2	NAG	O5-C5-C6-O6
8	K	6	MAN	C4-C5-C6-O6
6	I	1	NAG	C4-C5-C6-O6
6	I	2	NAG	O5-C5-C6-O6
6	O	1	NAG	O5-C5-C6-O6
6	I	1	NAG	O5-C5-C6-O6
7	D	2	NAG	C3-C2-N2-C7
7	D	2	NAG	C4-C5-C6-O6
6	J	1	NAG	C1-C2-N2-C7
8	K	4	MAN	O5-C5-C6-O6
7	D	2	NAG	O5-C5-C6-O6
5	M	2	NAG	C4-C5-C6-O6
5	M	2	NAG	O5-C5-C6-O6
6	J	2	NAG	C4-C5-C6-O6
7	D	1	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	M	2	NAG	C3-C2-N2-C7
8	K	2	NAG	C3-C2-N2-C7
6	J	2	NAG	O5-C5-C6-O6
5	F	2	NAG	C1-C2-N2-C7
6	E	2	NAG	C1-C2-N2-C7
5	A	2	NAG	O5-C5-C6-O6
6	N	2	NAG	C1-C2-N2-C7
5	F	1	NAG	C3-C2-N2-C7
6	C	2	NAG	C3-C2-N2-C7
6	J	1	NAG	C3-C2-N2-C7
6	N	1	NAG	C3-C2-N2-C7
5	A	3	BMA	C4-C5-C6-O6

All (1) ring outliers are listed below:

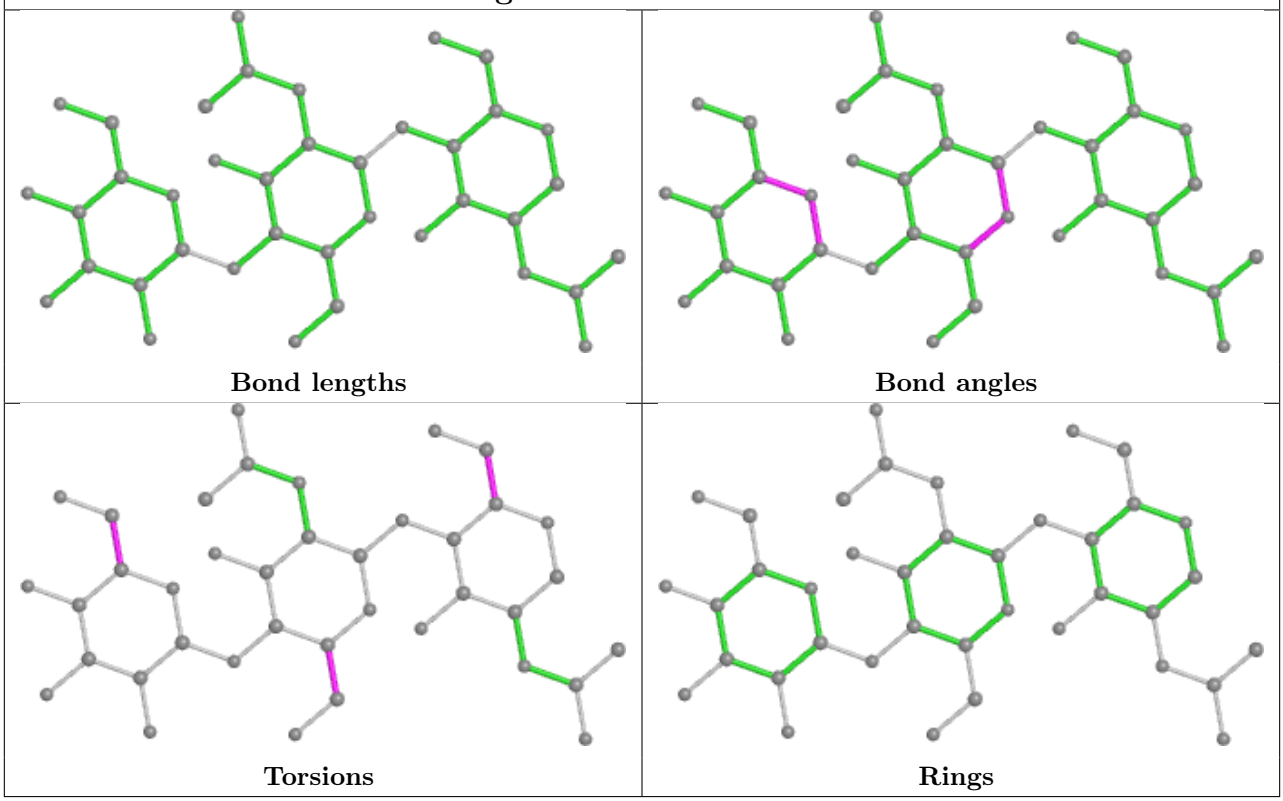
Mol	Chain	Res	Type	Atoms
8	K	6	MAN	C1-C2-C3-C4-C5-O5

5 monomers are involved in 7 short contacts:

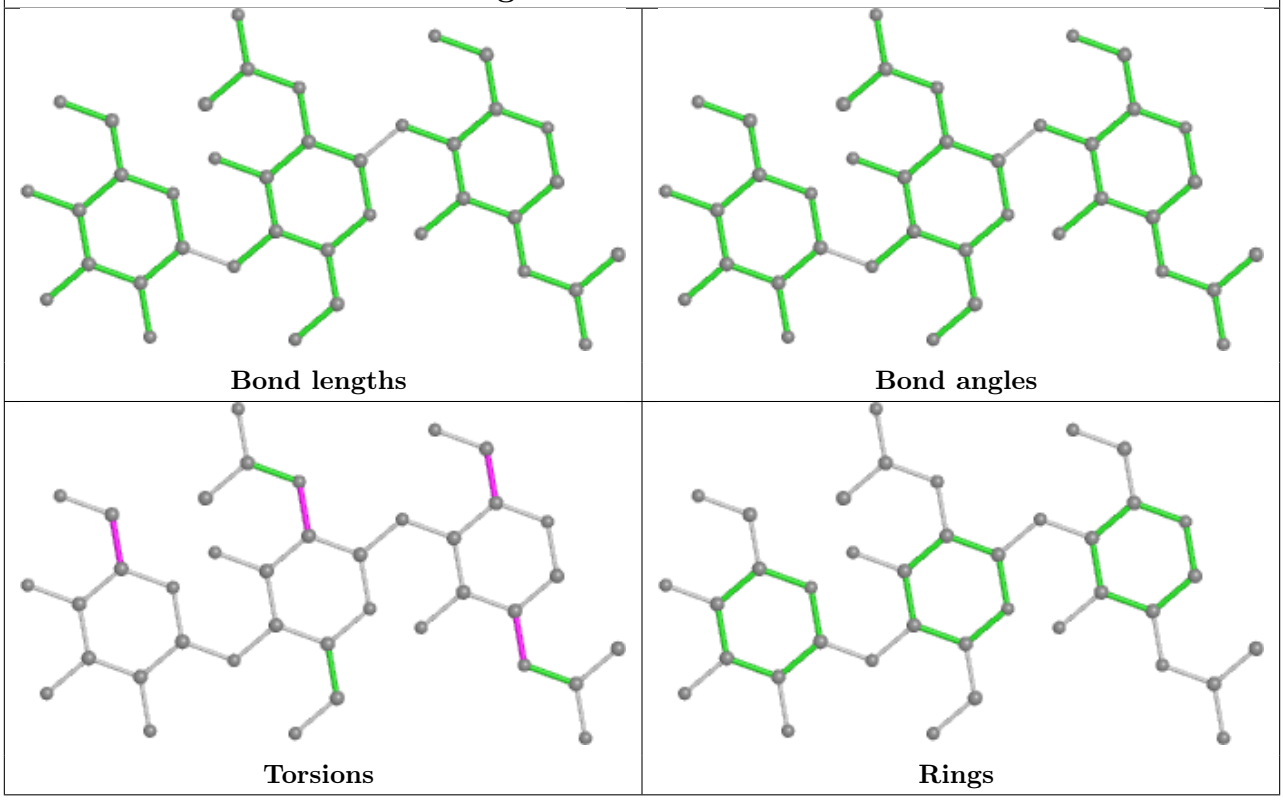
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	2	NAG	1	0
8	K	1	NAG	2	0
5	M	1	NAG	2	0
7	D	1	NAG	1	0
6	C	1	NAG	1	0

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

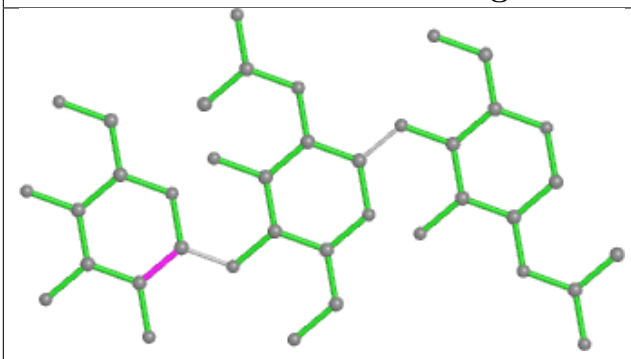
Oligosaccharide Chain A



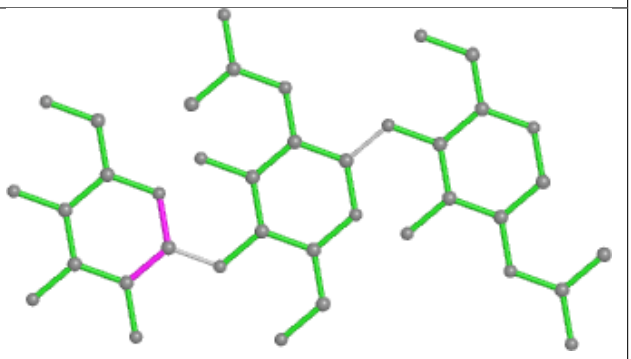
Oligosaccharide Chain F



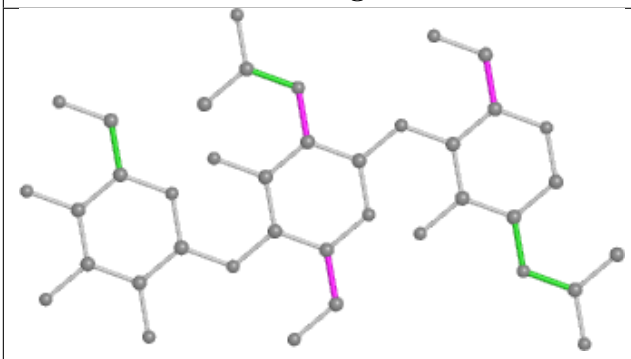
Oligosaccharide Chain M



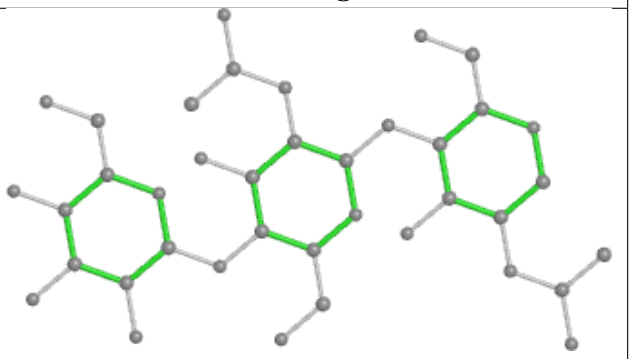
Bond lengths



Bond angles

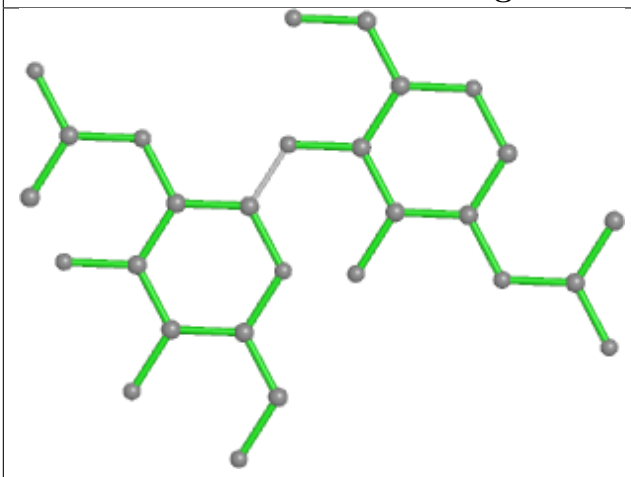


Torsions

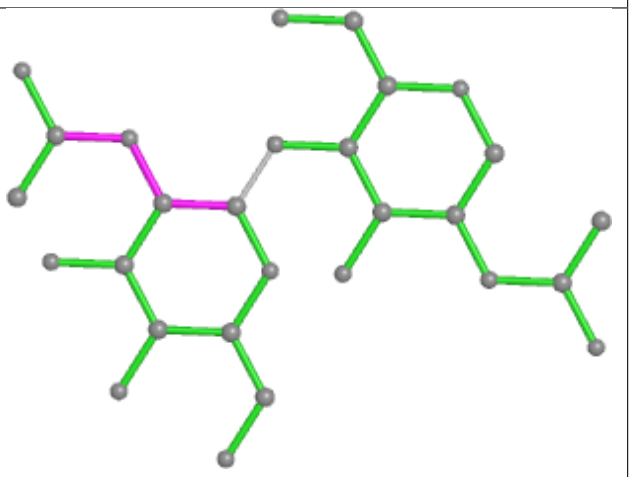


Rings

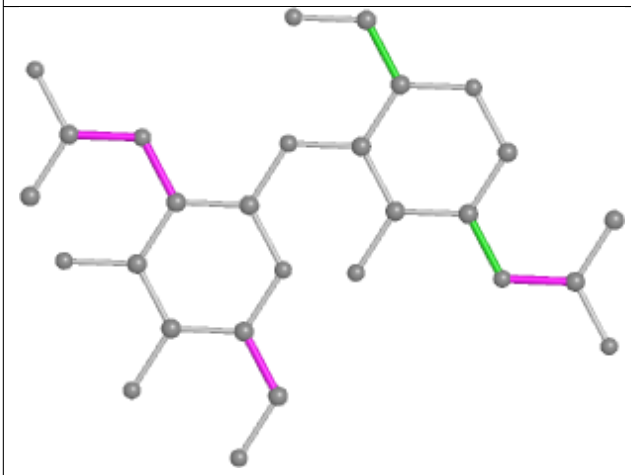
Oligosaccharide Chain C



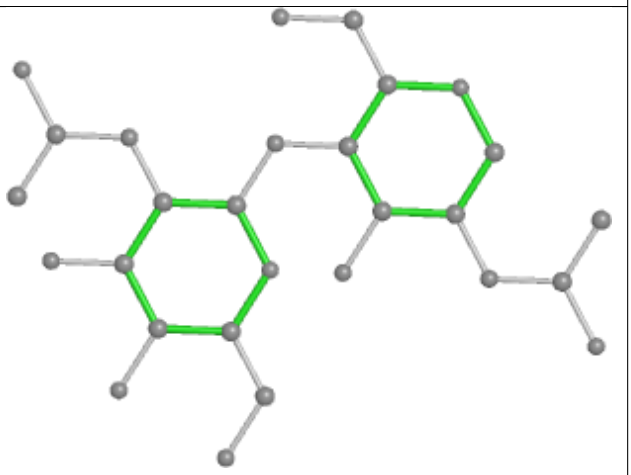
Bond lengths



Bond angles

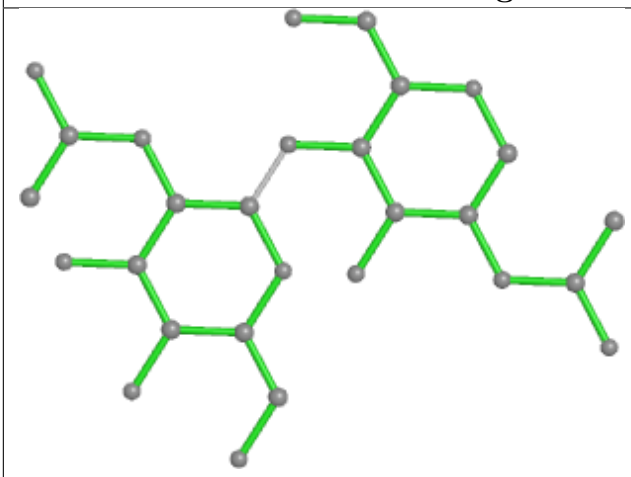


Torsions

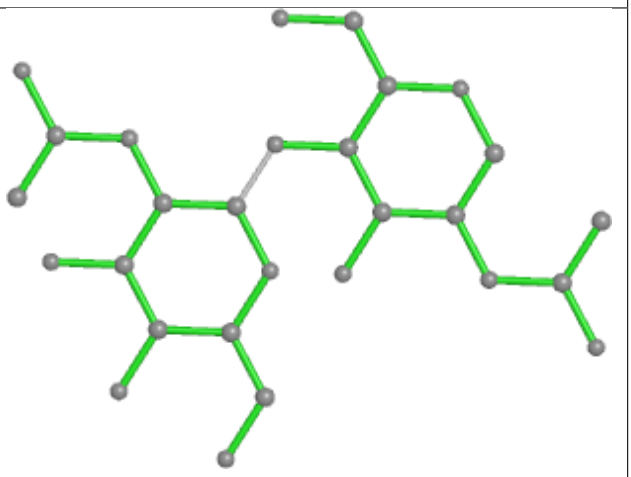


Rings

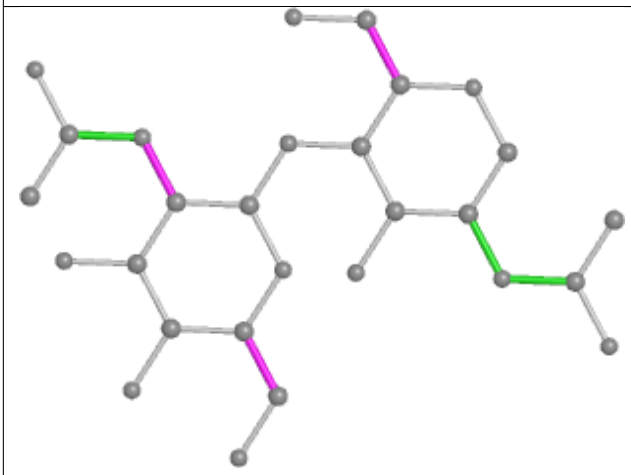
Oligosaccharide Chain E



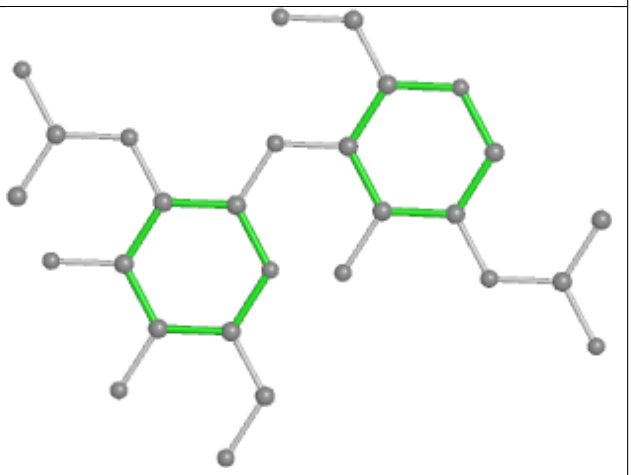
Bond lengths



Bond angles

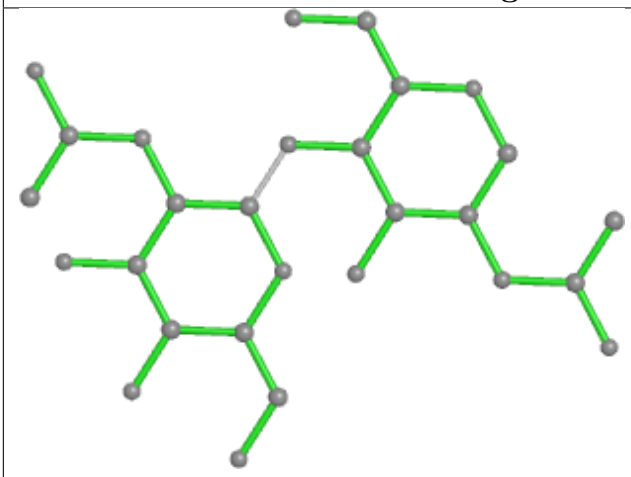


Torsions

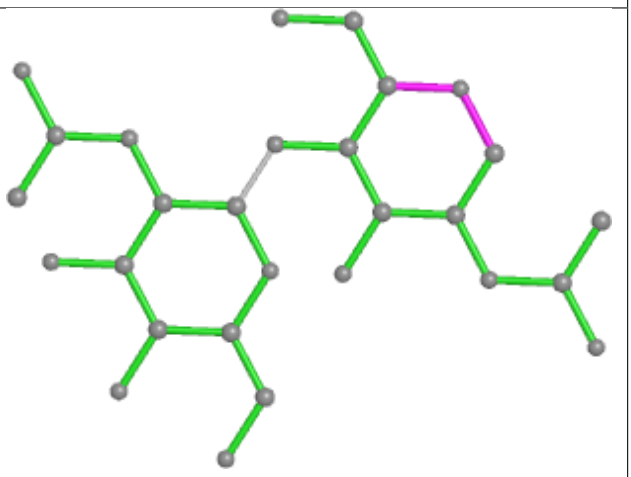


Rings

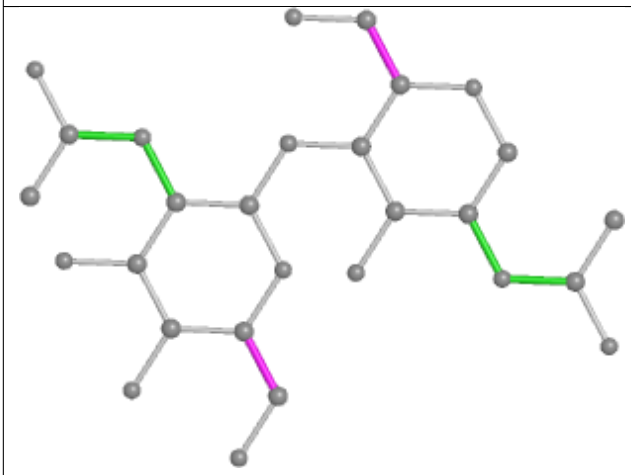
Oligosaccharide Chain I



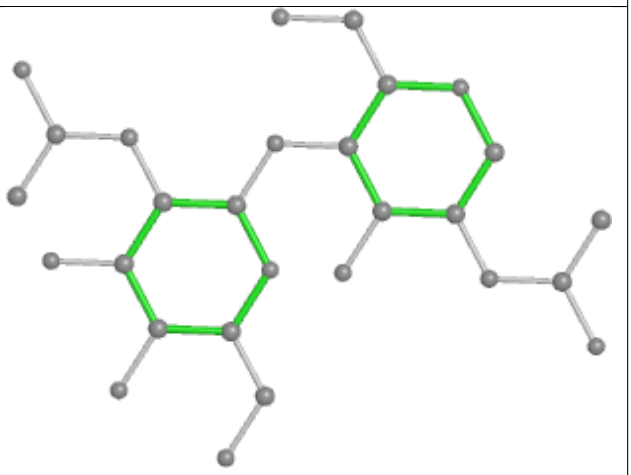
Bond lengths



Bond angles

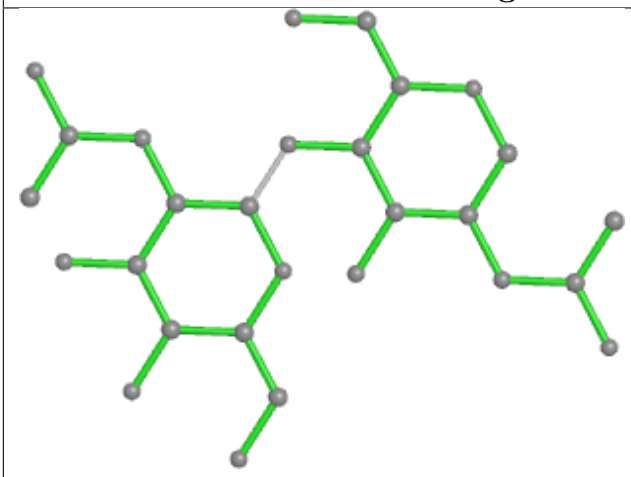


Torsions

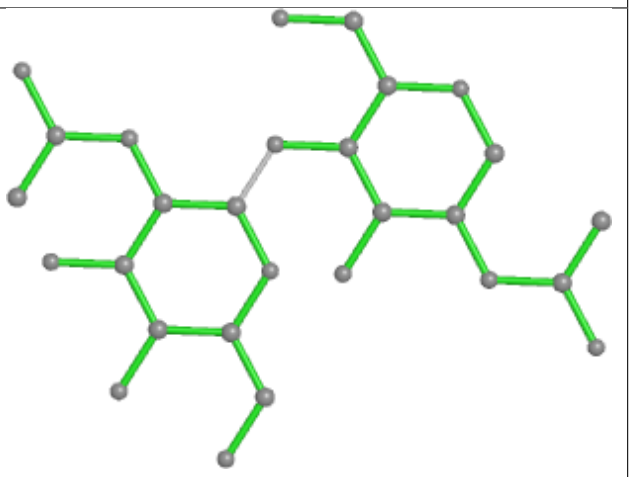


Rings

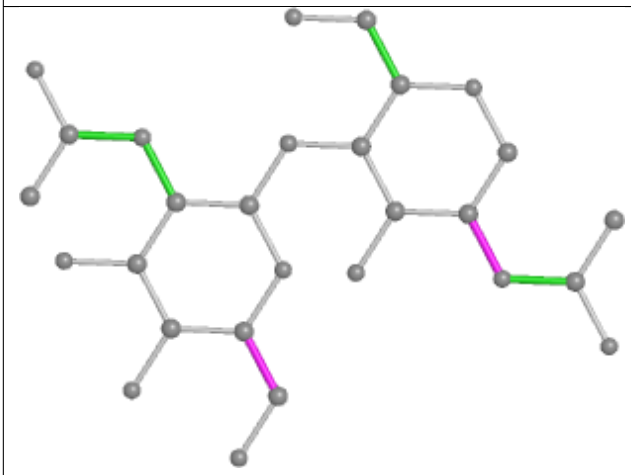
Oligosaccharide Chain J



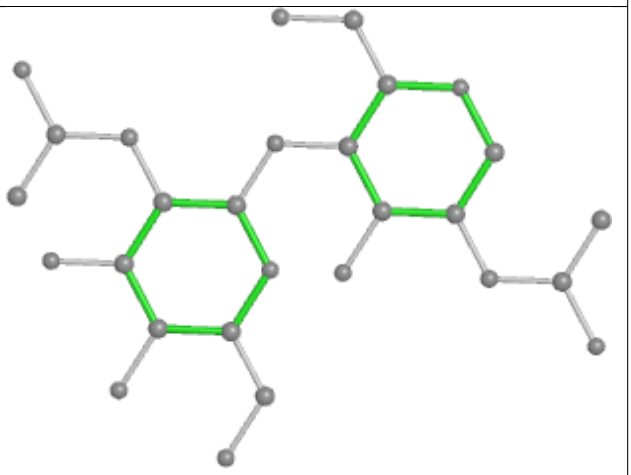
Bond lengths



Bond angles

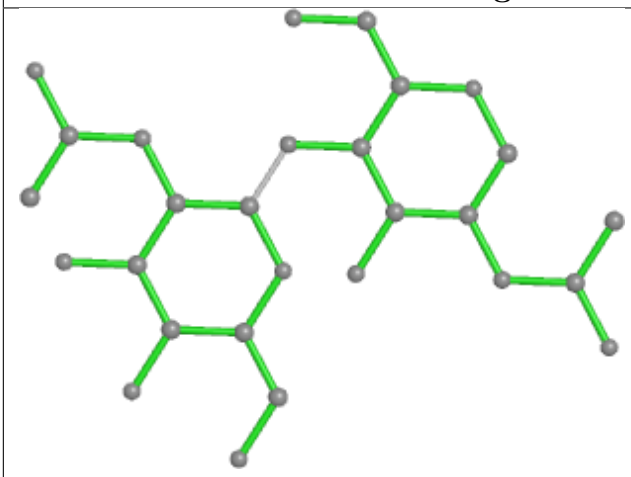


Torsions

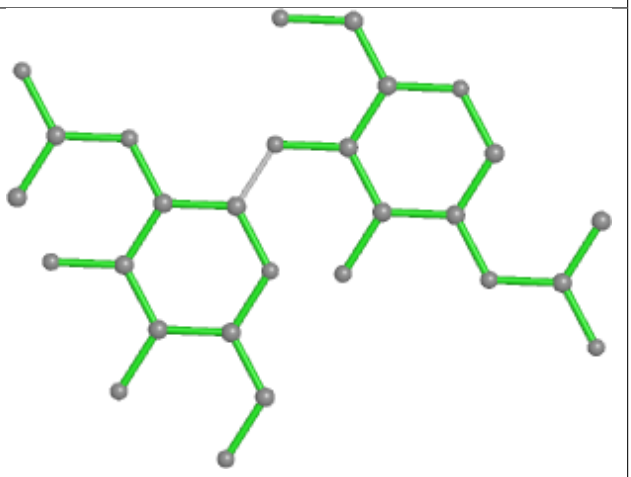


Rings

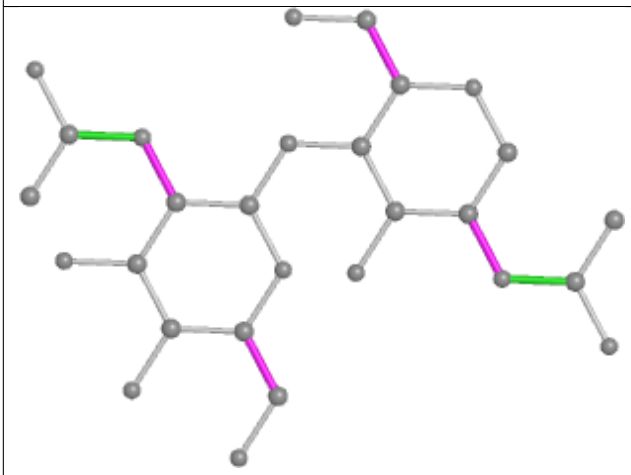
Oligosaccharide Chain N



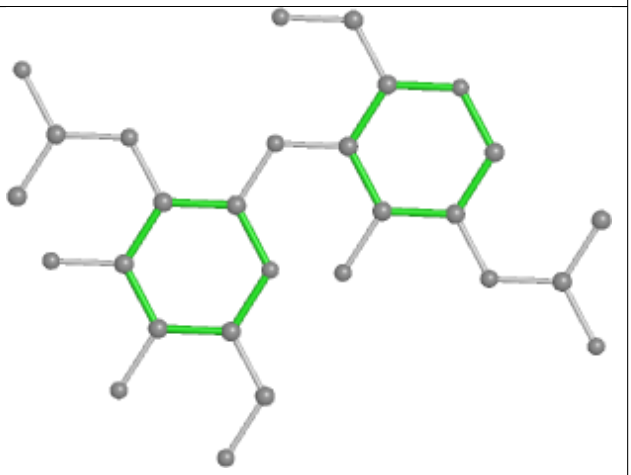
Bond lengths



Bond angles

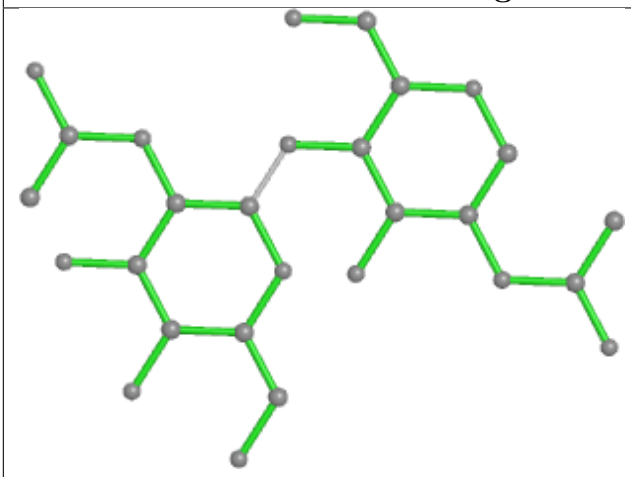


Torsions

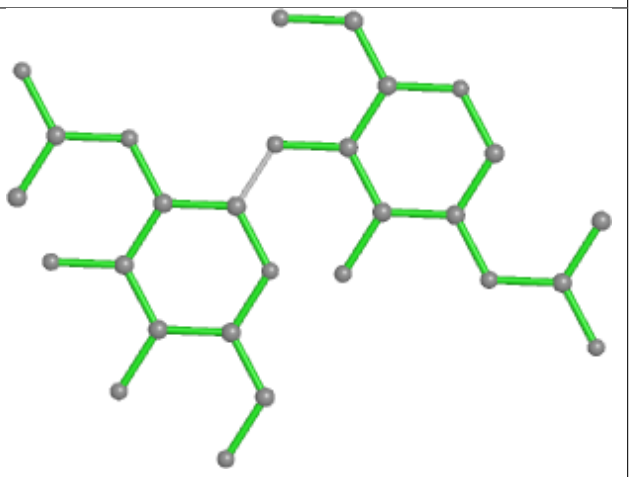


Rings

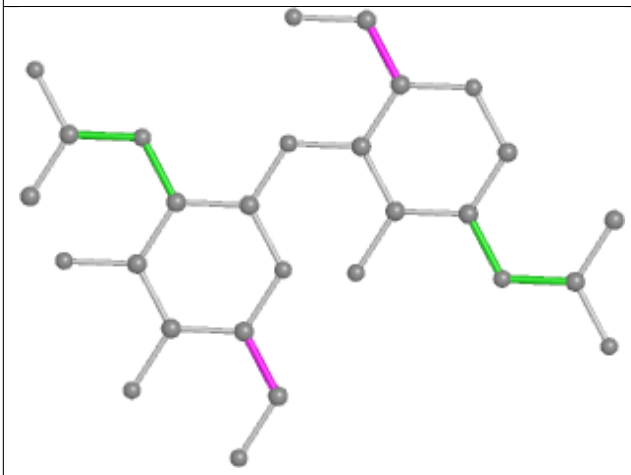
Oligosaccharide Chain O



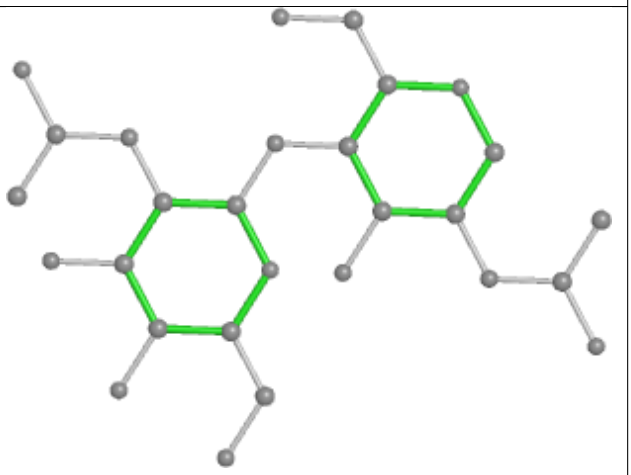
Bond lengths



Bond angles

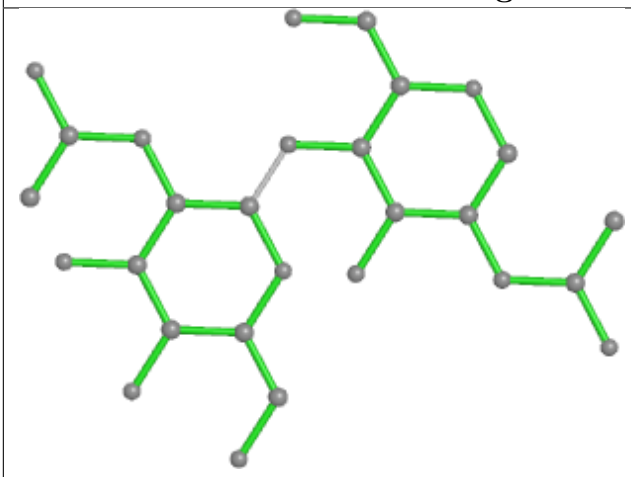


Torsions

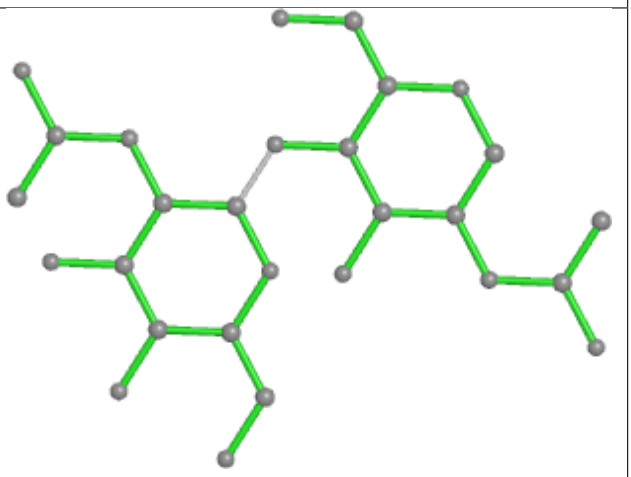


Rings

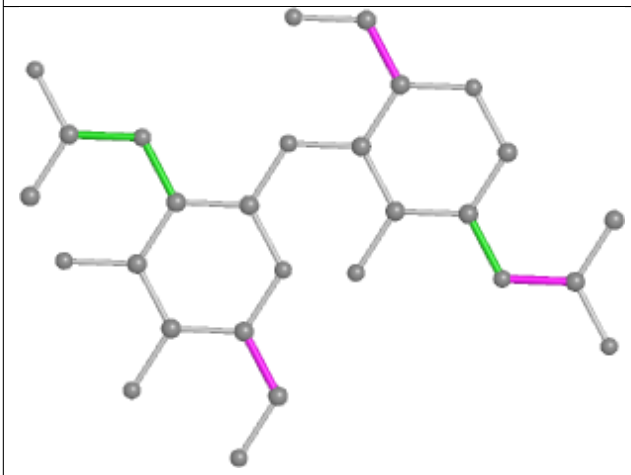
Oligosaccharide Chain P



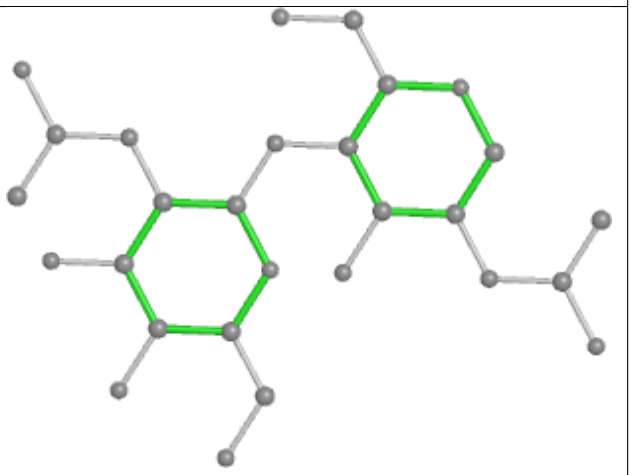
Bond lengths



Bond angles

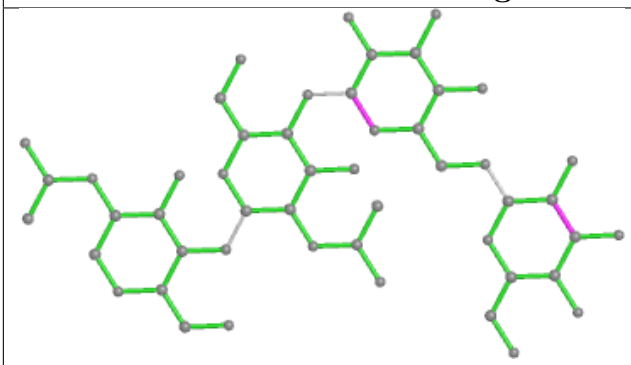


Torsions

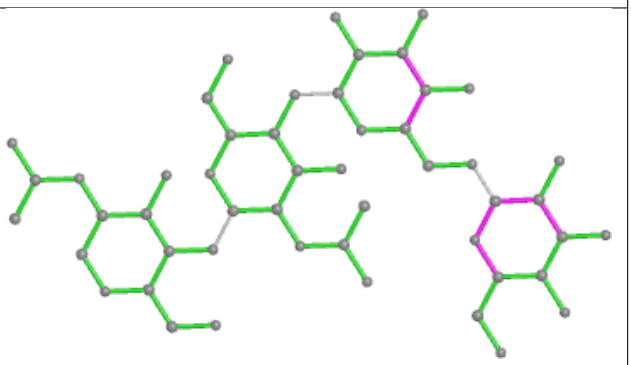


Rings

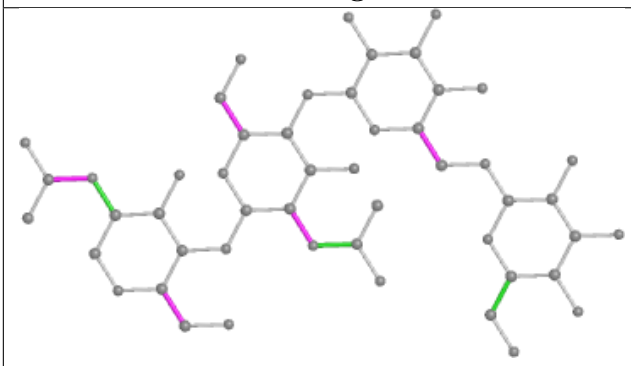
Oligosaccharide Chain D



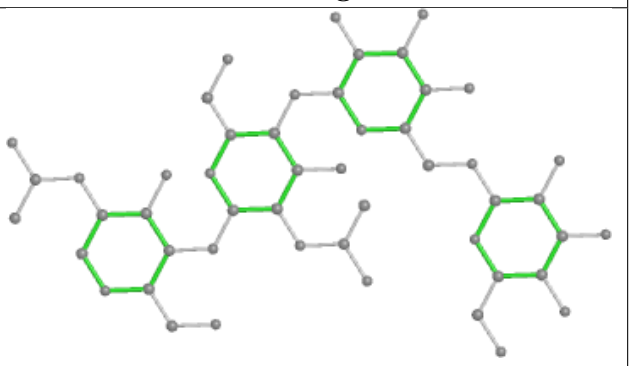
Bond lengths



Bond angles

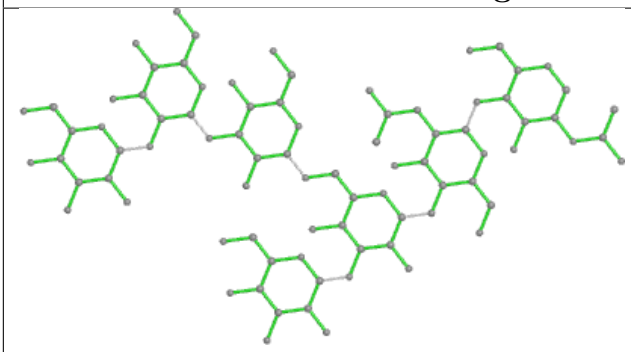


Torsions

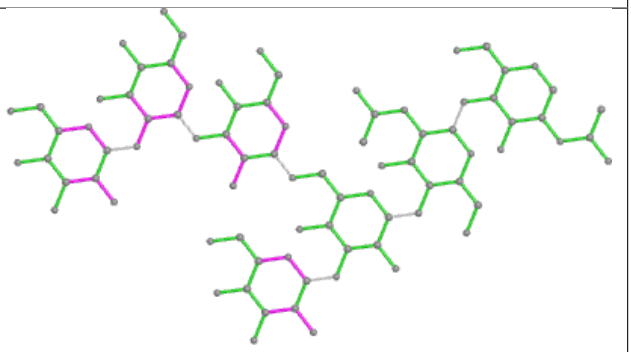


Rings

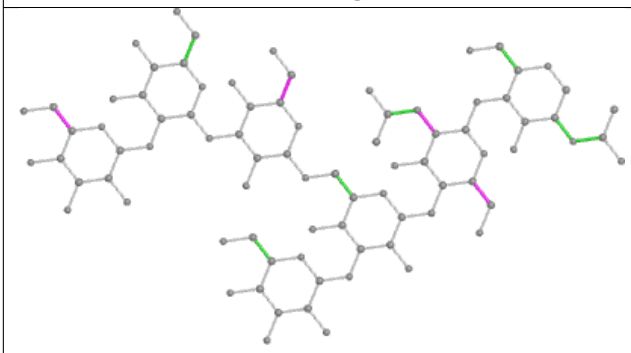
Oligosaccharide Chain K



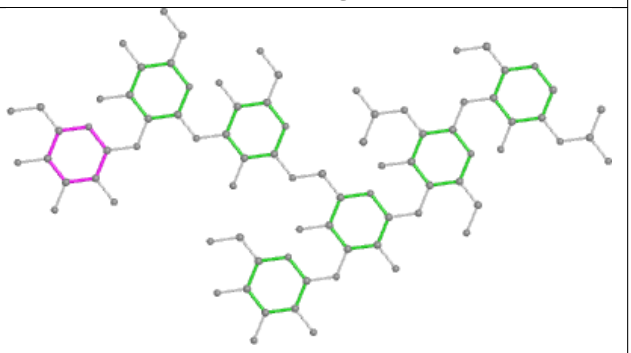
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

7 ligands 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	$\#\|Z\| > 2$
9	NAG	G	603	1	14,14,15	0.21	0	17,19,21	0.53	0
9	NAG	B	702	2	14,14,15	0.38	0	17,19,21	0.42	0
9	NAG	G	601	1	14,14,15	0.40	0	17,19,21	0.40	0
9	NAG	G	602	1	14,14,15	0.51	0	17,19,21	1.39	2 (11%)
9	NAG	G	604	1	14,14,15	0.33	0	17,19,21	0.34	0
9	NAG	B	703	2	14,14,15	0.67	0	17,19,21	0.39	0
9	NAG	B	701	2	14,14,15	0.24	0	17,19,21	0.33	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	G	603	1	-	4/6/23/26	0/1/1/1
9	NAG	B	702	2	-	0/6/23/26	0/1/1/1
9	NAG	G	601	1	-	2/6/23/26	0/1/1/1
9	NAG	G	602	1	-	5/6/23/26	0/1/1/1
9	NAG	G	604	1	-	2/6/23/26	0/1/1/1
9	NAG	B	703	2	-	0/6/23/26	0/1/1/1
9	NAG	B	701	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	G	602	NAG	C2-N2-C7	4.67	129.56	122.90
9	G	602	NAG	C1-C2-N2	2.23	114.30	110.49

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	G	601	NAG	C4-C5-C6-O6
9	G	601	NAG	O5-C5-C6-O6
9	G	603	NAG	O5-C5-C6-O6
9	G	604	NAG	O5-C5-C6-O6
9	B	701	NAG	O5-C5-C6-O6
9	G	602	NAG	O5-C5-C6-O6
9	G	602	NAG	C4-C5-C6-O6
9	G	604	NAG	C4-C5-C6-O6
9	B	701	NAG	C4-C5-C6-O6
9	G	602	NAG	C8-C7-N2-C2
9	G	602	NAG	O7-C7-N2-C2
9	G	603	NAG	C8-C7-N2-C2
9	G	603	NAG	O7-C7-N2-C2
9	G	603	NAG	C4-C5-C6-O6
9	G	602	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B	702	NAG	1	0
9	G	602	NAG	4	0
9	B	703	NAG	1	0

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 (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	G	449/481 (93%)	-0.06	10 (2%)	62	52	154, 213, 266, 325	0
2	B	122/153 (79%)	0.07	8 (6%)	18	14	215, 289, 343, 400	0
3	H	225/229 (98%)	-0.12	5 (2%)	62	52	149, 190, 243, 257	0
4	L	200/205 (97%)	0.03	10 (5%)	28	24	168, 212, 245, 265	0
All	All	996/1068 (93%)	-0.04	33 (3%)	46	37	149, 213, 298, 400	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	496	VAL	5.1
1	G	498	PRO	5.0
1	G	497	ALA	4.8
2	B	541	ALA	4.5
2	B	544	LEU	4.4
2	B	537	LEU	4.3
1	G	40	TYR	3.5
4	L	35	TRP	3.4
2	B	602	LEU	3.2
2	B	545	LEU	3.0
2	B	540	GLN	2.9
1	G	444	ARG	2.8
3	H	121	VAL	2.7
4	L	21	ILE	2.6
4	L	86	TYR	2.5
2	B	614	TRP	2.5
1	G	291	PRO	2.5
4	L	37	GLN	2.4
1	G	495	GLY	2.4
4	L	20	THR	2.3
3	H	133	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
4	L	47	LEU	2.3
3	H	209	LYS	2.2
1	G	252	LYS	2.2
4	L	82	ASP	2.1
4	L	48	ILE	2.1
4	L	19	VAL	2.1
1	G	443	ILE	2.1
4	L	179	SER	2.1
3	H	100(B)	ARG	2.1
1	G	153	GLU	2.1
3	H	210	ARG	2.0
2	B	539	VAL	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
5	BMA	A	3	11/12	0.63	0.52	336,340,341,342	0
8	MAN	K	4	11/12	0.65	0.24	389,393,397,398	0
5	BMA	M	3	11/12	0.66	0.40	327,331,336,342	0
8	MAN	K	7	11/12	0.69	0.36	372,380,398,404	0
8	BMA	K	3	11/12	0.70	0.19	371,384,392,397	0
6	NAG	O	2	14/15	0.72	0.33	344,351,355,359	0
6	NAG	I	1	14/15	0.72	0.50	304,313,328,340	0
7	MAN	D	4	11/12	0.75	0.75	321,323,326,327	0
6	NAG	I	2	14/15	0.77	0.27	350,360,365,365	0
7	BMA	D	3	11/12	0.78	0.48	305,308,316,318	0
5	BMA	F	3	11/12	0.79	0.28	333,336,338,338	0
6	NAG	C	2	14/15	0.79	0.41	302,310,313,314	0
6	NAG	N	2	14/15	0.79	0.16	323,332,342,345	0
6	NAG	E	2	14/15	0.79	1.34	357,364,370,372	0
6	NAG	P	2	14/15	0.79	0.27	390,405,416,432	0
6	NAG	J	2	14/15	0.82	0.32	336,346,350,351	0

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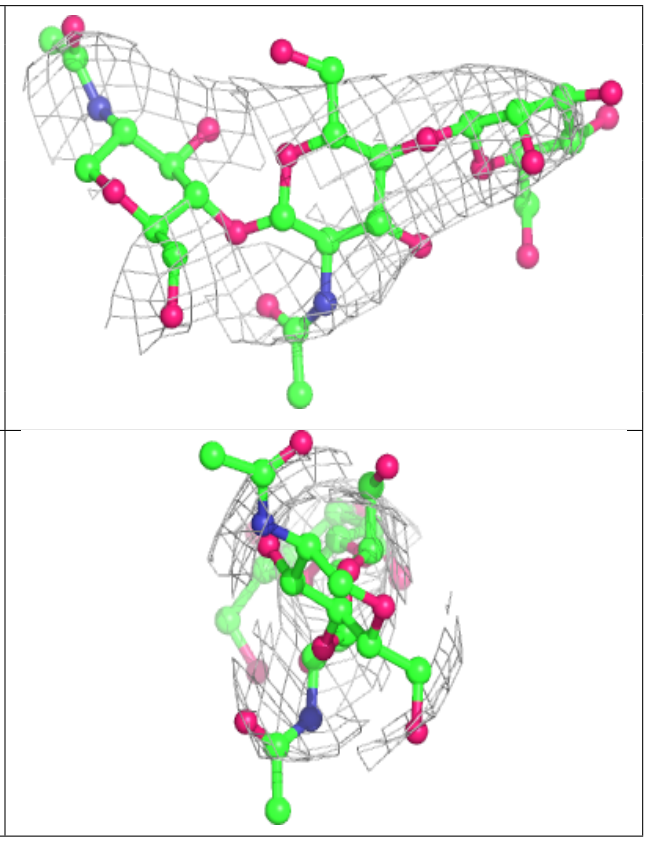
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	NAG	M	2	14/15	0.83	0.23	322,327,334,338	0
6	NAG	O	1	14/15	0.84	0.35	300,315,324,335	0
6	NAG	C	1	14/15	0.85	0.38	267,275,284,298	0
8	MAN	K	6	11/12	0.86	0.15	378,386,389,389	0
8	NAG	K	1	14/15	0.86	0.68	298,311,319,334	0
8	MAN	K	5	11/12	0.88	0.21	390,397,401,402	0
5	NAG	A	2	14/15	0.88	0.22	306,314,321,328	0
6	NAG	P	1	14/15	0.88	0.33	321,332,353,373	0
8	NAG	K	2	14/15	0.91	0.35	340,348,353,363	0
7	NAG	D	2	14/15	0.91	0.29	291,296,304,307	0
6	NAG	E	1	14/15	0.91	0.50	312,326,341,347	0
5	NAG	A	1	14/15	0.92	0.20	270,282,289,296	0
5	NAG	M	1	14/15	0.93	0.26	279,288,299,310	0
6	NAG	J	1	14/15	0.93	0.20	293,308,315,329	0
7	NAG	D	1	14/15	0.94	0.24	253,270,277,284	0
5	NAG	F	2	14/15	0.95	0.22	318,327,333,334	0
5	NAG	F	1	14/15	0.96	0.32	281,290,303,311	0
6	NAG	N	1	14/15	0.96	0.22	279,291,313,318	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.

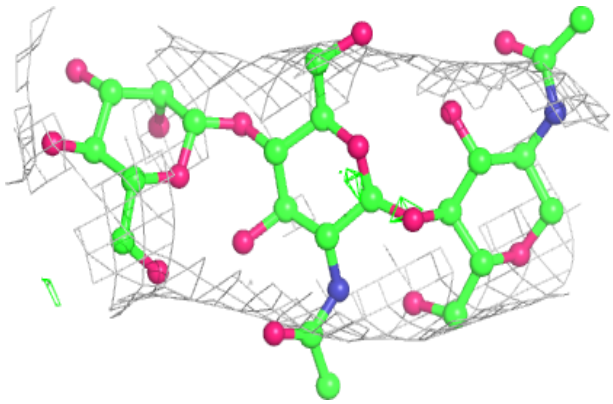
Electron density around Chain A:

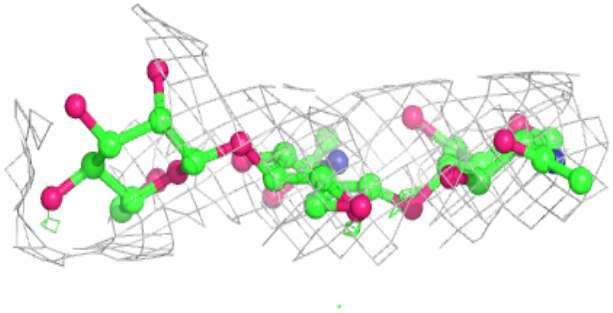
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

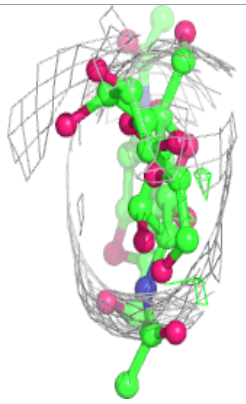


Electron density around Chain F:

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

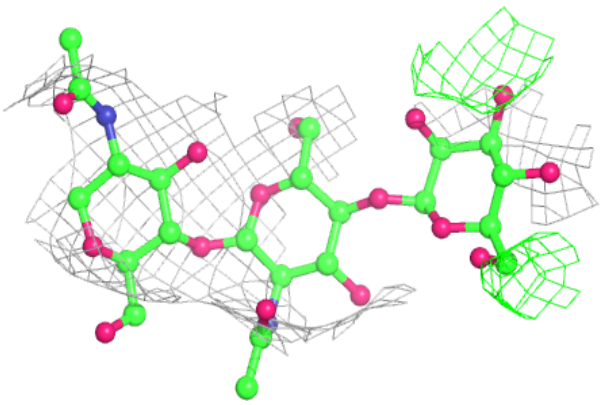


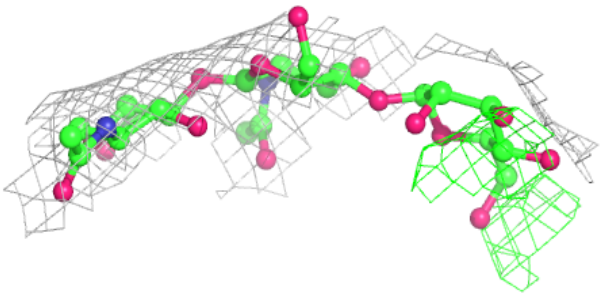


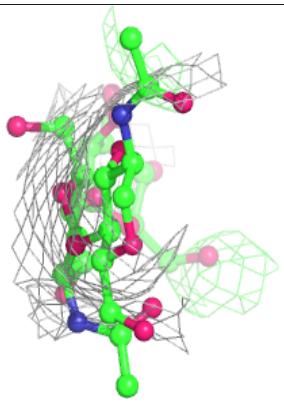


Electron density around Chain M:

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

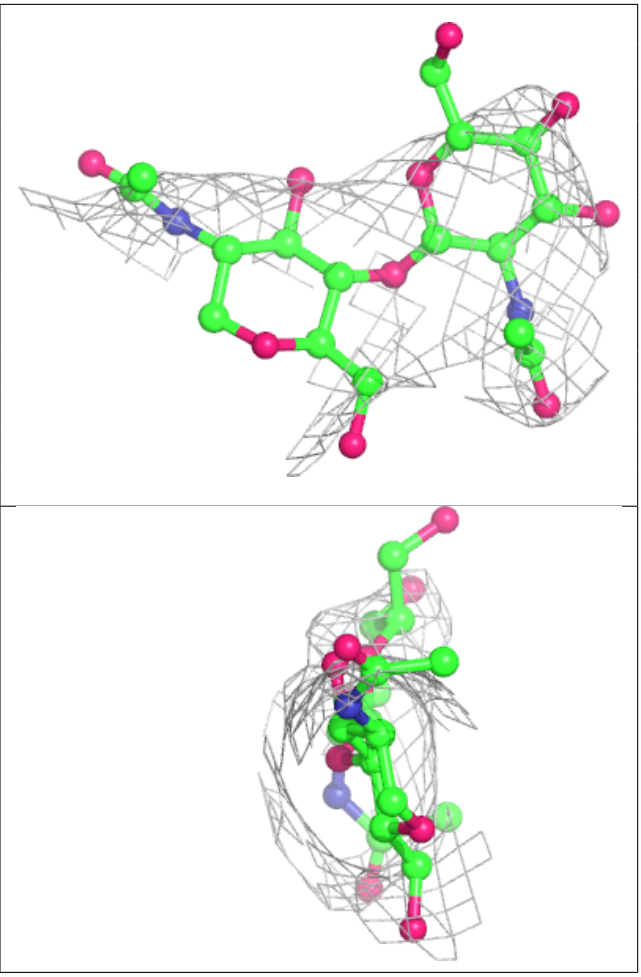






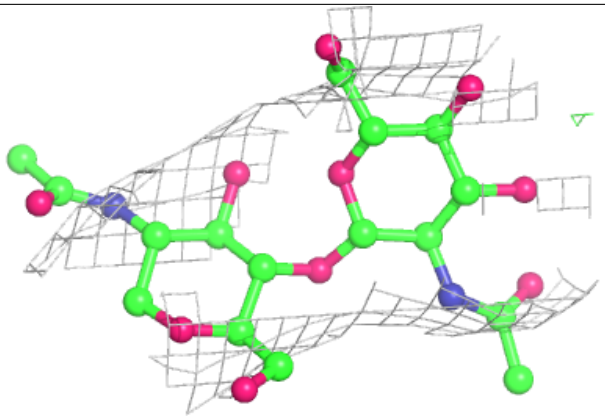
Electron density around Chain C:

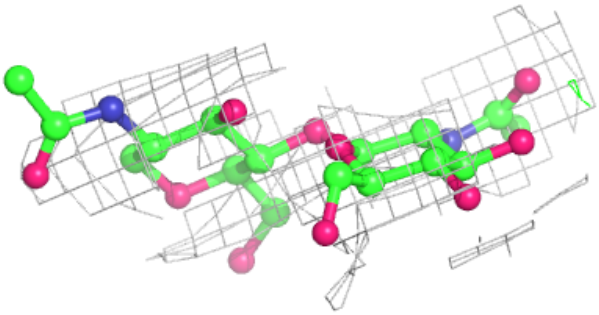
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

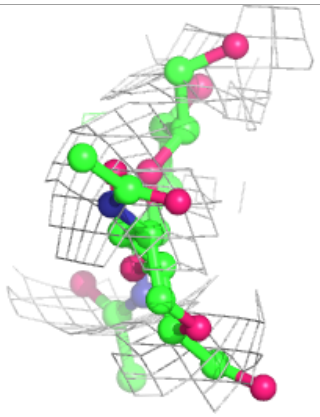


Electron density around Chain E:

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

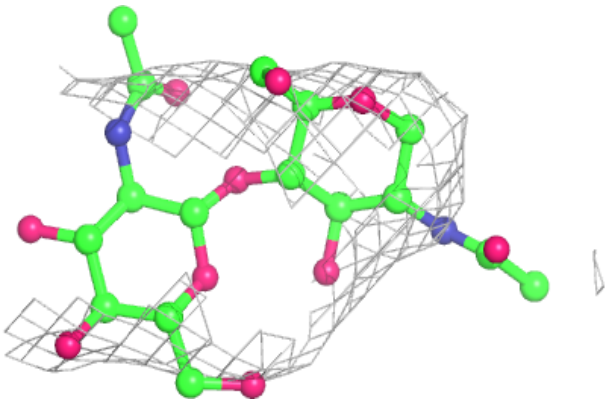


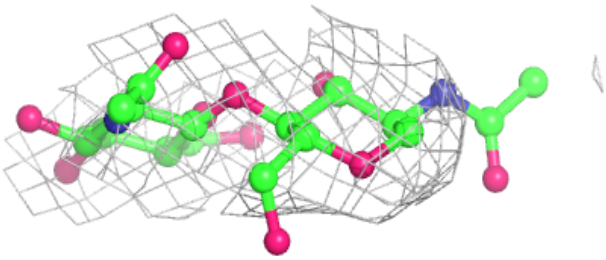


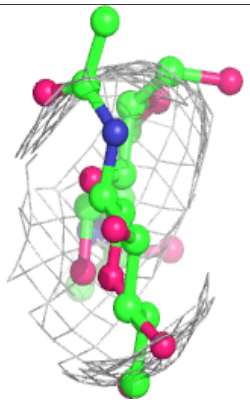


Electron density around Chain I:

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

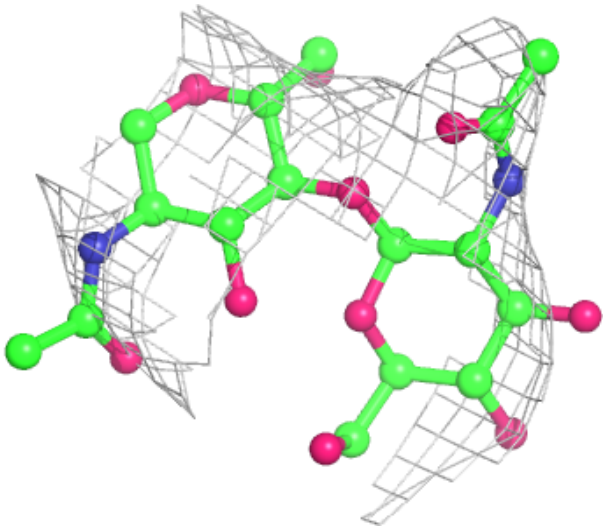


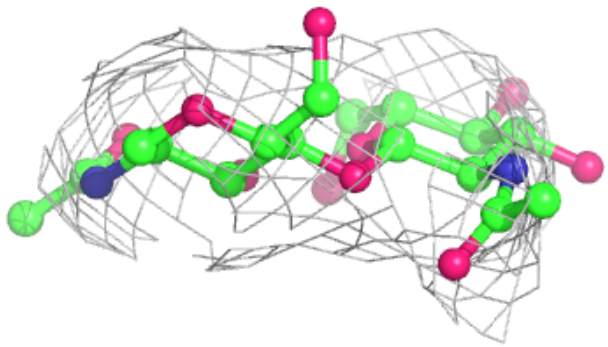


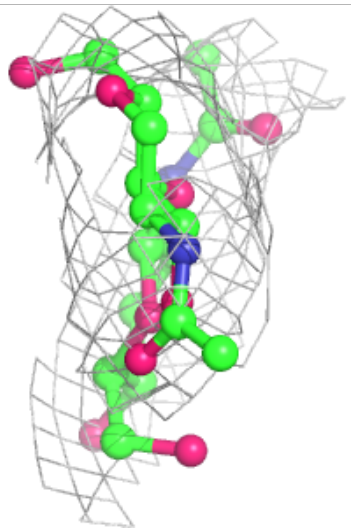


Electron density around Chain J:

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

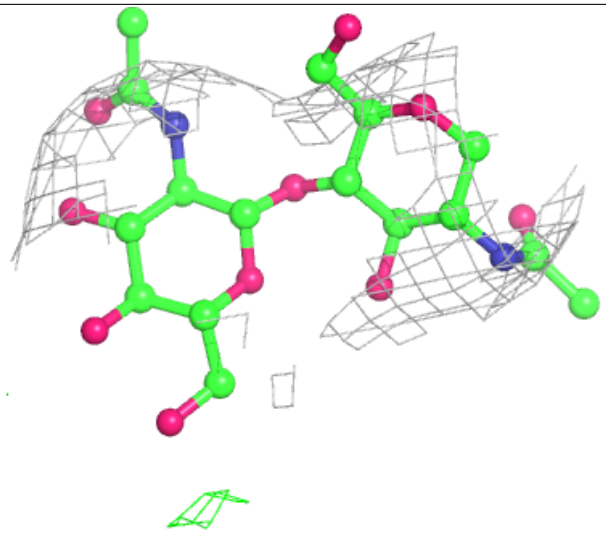


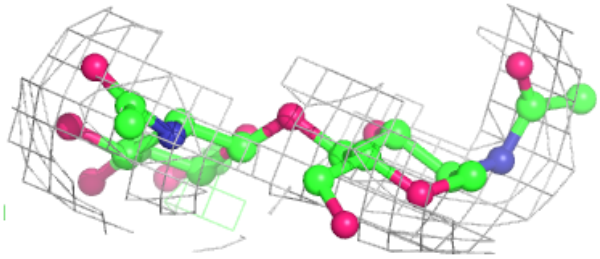


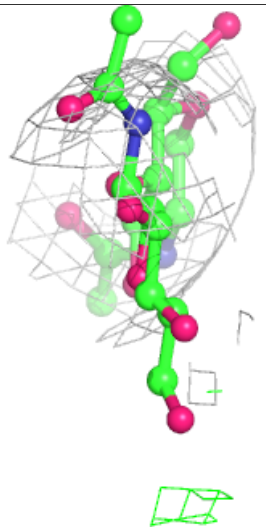


Electron density around Chain N:

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

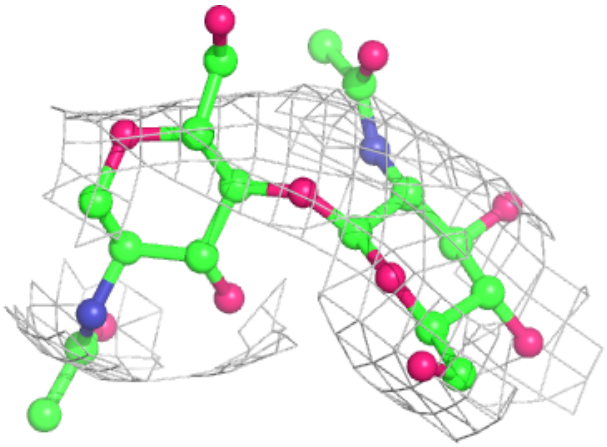


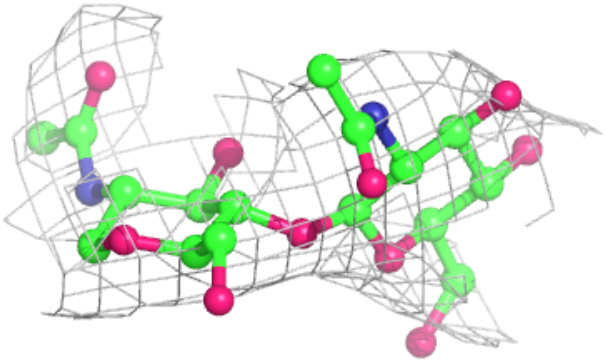


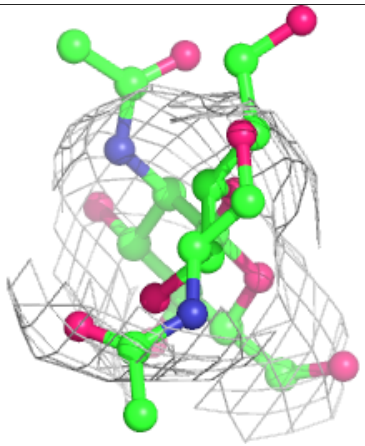


Electron density around Chain O:

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

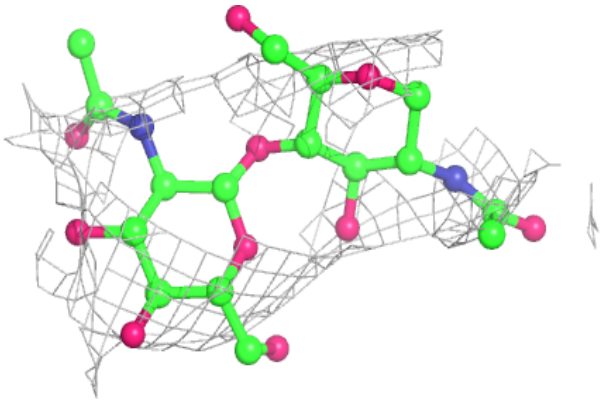


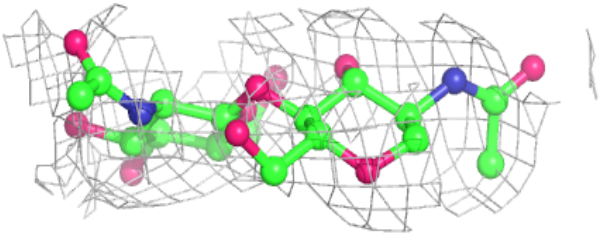


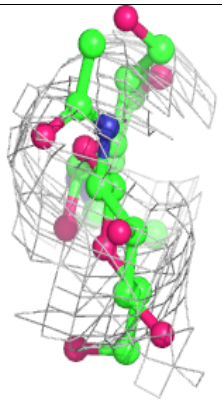


Electron density around Chain P:

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

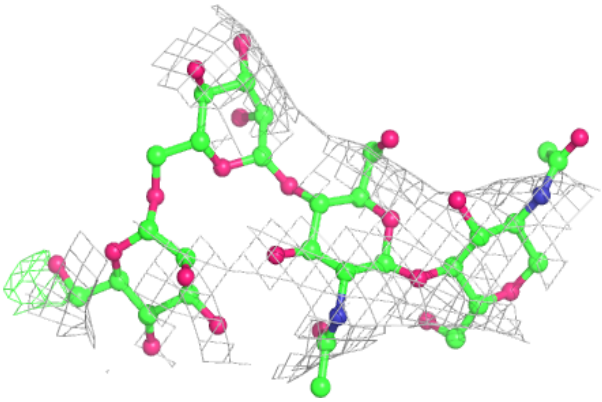


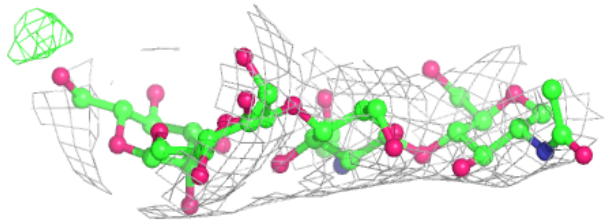


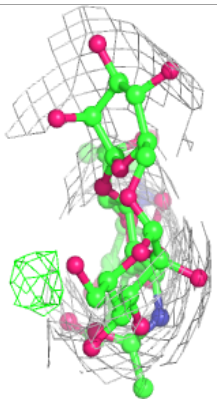


Electron density around Chain D:

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

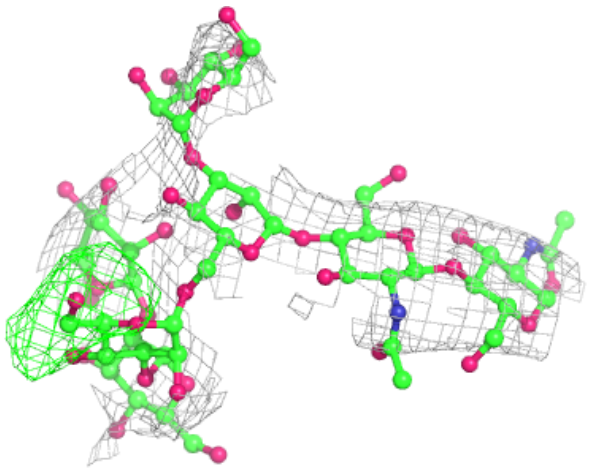


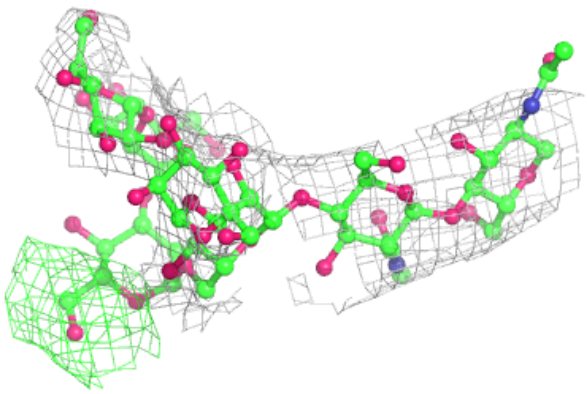


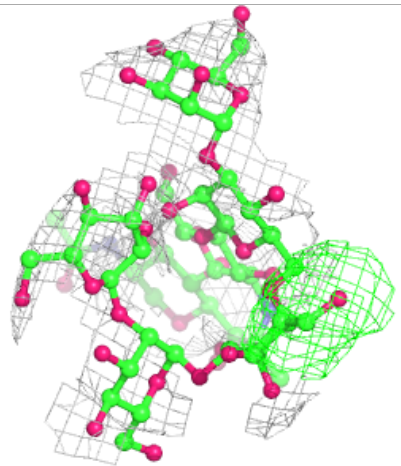


Electron density around Chain K:

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







6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
9	NAG	G	603	14/15	0.80	0.53	341,352,358,360	0
9	NAG	B	703	14/15	0.82	0.18	200,230,258,262	0
9	NAG	B	702	14/15	0.83	0.20	221,236,260,261	0
9	NAG	G	604	14/15	0.83	0.26	321,334,340,341	0
9	NAG	G	601	14/15	0.85	0.30	262,271,275,276	0
9	NAG	G	602	14/15	0.86	0.36	293,305,313,314	0
9	NAG	B	701	14/15	0.86	0.47	212,234,263,273	0

6.5 Other polymers (i)

There are no such residues in this entry.