Full wwPDB X-ray Structure Validation Report ⓘ

Aug 24, 2021 – 08:10 pm BST

PDB ID	:	7A57
Title	:	La Crosse Virus Envelope Glycoprotein Gc W1066H Mutant Fusion Domains in Postfusion Conformation
Authors	:	Hellert, J.; Guardado-Calvo, P.; Rey, F.A.
Deposited on	:	2020-08-20
Resolution	:	3.15 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 following versions of software and data (see references ①) were used in the production of this report:

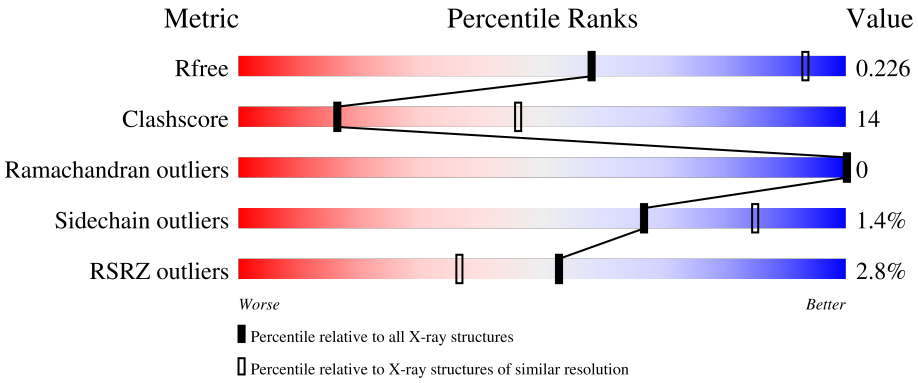
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.1

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 3.15 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	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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	A	458	2% 66% 28% • 5%
1	B	458	2% 70% 27% ••
1	C	458	5% 72% 22% • 5%
2	D	3	67% 33%
2	E	3	33% 67%

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Mol	Chain	Length	Quality of chain
2	F	3	67% 33%

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
2	NAG	F	2	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10385 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 Envelopment polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	437	Total	C	N	O	S	0	0	0
			3397	2152	557	665	23			
1	B	449	Total	C	N	O	S	0	0	0
			3485	2207	573	682	23			
1	C	436	Total	C	N	O	S	0	0	0
			3389	2148	556	662	23			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	907	GLU	-	expression tag	UNP Q8JPR1
A	908	TRP	-	expression tag	UNP Q8JPR1
A	909	SER	-	expression tag	UNP Q8JPR1
A	910	HIS	-	expression tag	UNP Q8JPR1
A	911	PRO	-	expression tag	UNP Q8JPR1
A	912	GLN	-	expression tag	UNP Q8JPR1
A	913	PHE	-	expression tag	UNP Q8JPR1
A	914	GLU	-	expression tag	UNP Q8JPR1
A	915	LYS	-	expression tag	UNP Q8JPR1
A	916	GLY	-	expression tag	UNP Q8JPR1
A	917	GLY	-	expression tag	UNP Q8JPR1
A	1066	HIS	TRP	engineered mutation	UNP Q8JPR1
B	907	GLU	-	expression tag	UNP Q8JPR1
B	908	TRP	-	expression tag	UNP Q8JPR1
B	909	SER	-	expression tag	UNP Q8JPR1
B	910	HIS	-	expression tag	UNP Q8JPR1
B	911	PRO	-	expression tag	UNP Q8JPR1
B	912	GLN	-	expression tag	UNP Q8JPR1
B	913	PHE	-	expression tag	UNP Q8JPR1
B	914	GLU	-	expression tag	UNP Q8JPR1
B	915	LYS	-	expression tag	UNP Q8JPR1
B	916	GLY	-	expression tag	UNP Q8JPR1
B	917	GLY	-	expression tag	UNP Q8JPR1

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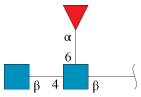
Chain	Residue	Modelled	Actual	Comment	Reference
B	1066	HIS	TRP	engineered mutation	UNP Q8JPR1
C	907	GLU	-	expression tag	UNP Q8JPR1
C	908	TRP	-	expression tag	UNP Q8JPR1
C	909	SER	-	expression tag	UNP Q8JPR1
C	910	HIS	-	expression tag	UNP Q8JPR1
C	911	PRO	-	expression tag	UNP Q8JPR1
C	912	GLN	-	expression tag	UNP Q8JPR1
C	913	PHE	-	expression tag	UNP Q8JPR1
C	914	GLU	-	expression tag	UNP Q8JPR1
C	915	LYS	-	expression tag	UNP Q8JPR1
C	916	GLY	-	expression tag	UNP Q8JPR1
C	917	GLY	-	expression tag	UNP Q8JPR1
C	1066	HIS	TRP	engineered mutation	UNP Q8JPR1

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



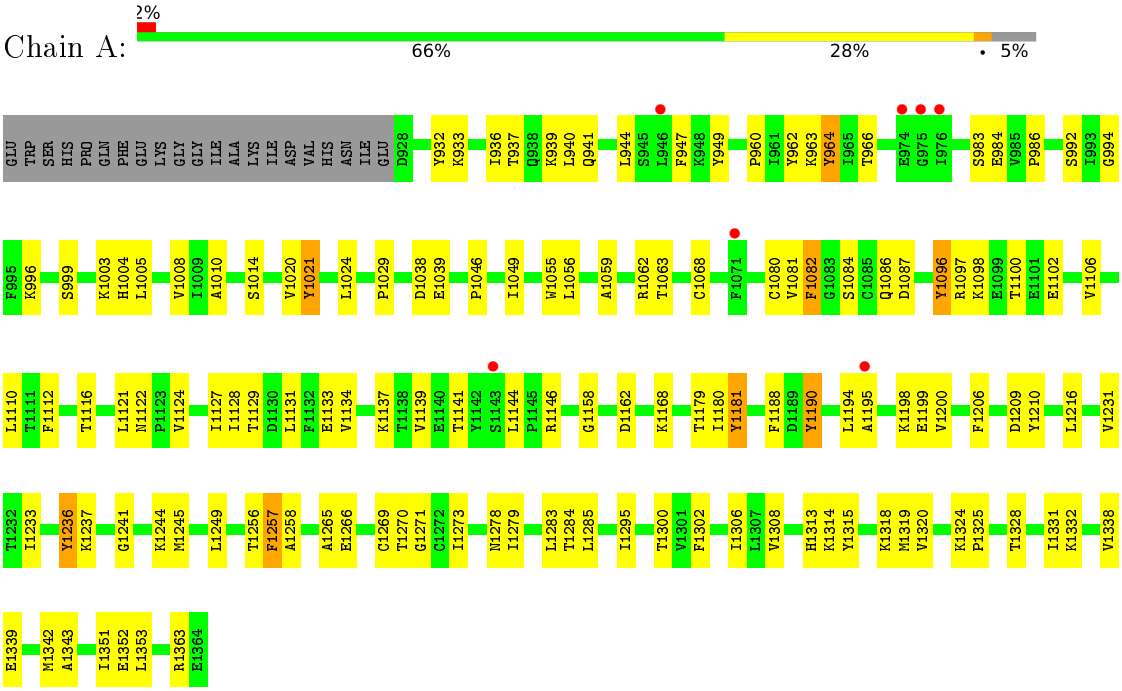
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	E	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	F	3	Total	C	N	O	0	0	0
			38	22	2	14			

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.

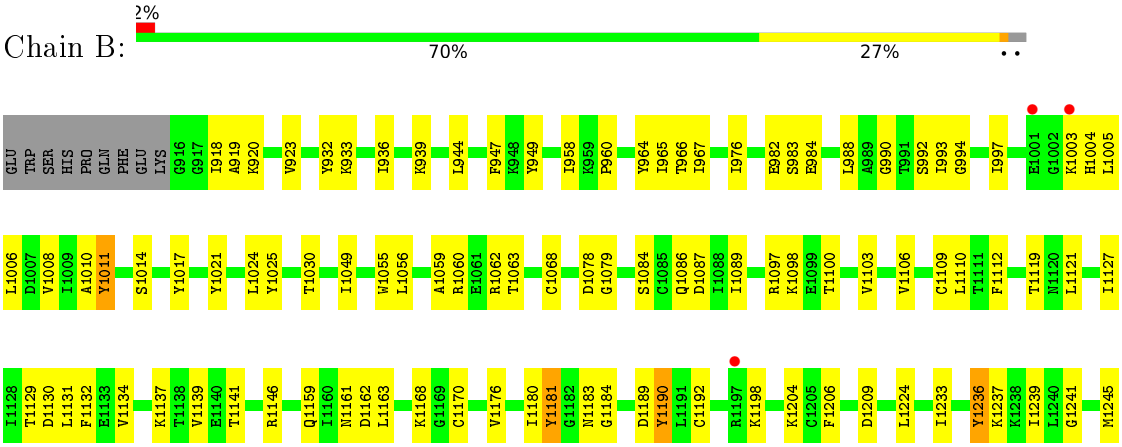
- Molecule 1: Envelopment polyprotein

Chain A:



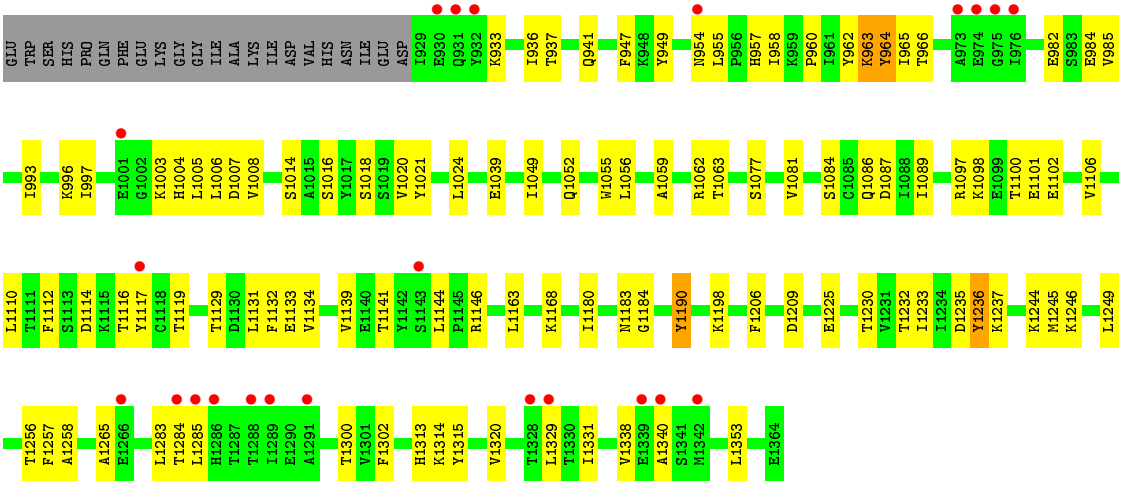
- Molecule 1: Envelopment polyprotein

Chain B:



• Molecule 1: Envelopment polyprotein

Chain C:

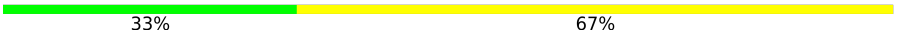


• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:

• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:



• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	170.87Å 170.87Å 379.22Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.06 – 3.15 48.06 – 3.16	Depositor EDS
% Data completeness (in resolution range)	67.0 (48.06-3.15) 67.0 (48.06-3.16)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.10 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.199 , 0.226 0.199 , 0.226	Depositor DCC
R_{free} test set	2298 reflections (6.01%)	wwPDB-VP
Wilson B-factor (Å^2)	99.6	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 70.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	10385	wwPDB-VP
Average B, all atoms (Å^2)	109.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% 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

5.1 Standard geometry

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

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	0.32	0/3460	0.61	0/4688
1	B	0.33	0/3549	0.62	0/4808
1	C	0.32	0/3452	0.59	1/4677 (0.0%)
All	All	0.32	0/10461	0.61	1/14173 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	963	LYS	CD-CE-NZ	6.10	125.73	111.70

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	3397	0	3383	119	0
1	B	3485	0	3472	116	0
1	C	3389	0	3379	97	0
2	D	38	0	34	1	0
2	E	38	0	34	1	0
2	F	38	0	34	0	0
All	All	10385	0	10336	298	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 14.

All (298) 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:B:1011:TYR:CE1	1:B:1109:CYS:HB2	2.12	0.84
1:A:1046:PRO:HG2	1:A:1082:PHE:CE2	2.15	0.81
1:A:1055:TRP:NE1	1:A:1086:GLN:OE1	2.15	0.80
1:A:1086:GLN:HE21	1:A:1180:ILE:HB	1.47	0.79
1:B:1055:TRP:NE1	1:B:1086:GLN:OE1	2.17	0.78
1:A:996:LYS:HE3	1:A:1257:PHE:CD2	2.22	0.73
1:B:1086:GLN:HE21	1:B:1180:ILE:HB	1.54	0.73
1:C:1236:TYR:CD1	1:C:1237:LYS:HG3	2.25	0.71
1:B:919:ALA:HB2	1:B:1130:ASP:HB3	1.72	0.71
1:B:1236:TYR:CD1	1:B:1237:LYS:HG3	2.26	0.70
1:A:1139:VAL:HG22	1:B:1139:VAL:HG21	1.73	0.70
1:C:1236:TYR:CE1	1:C:1237:LYS:HG3	2.28	0.69
1:B:1021:TYR:CE2	1:B:1233:ILE:HD12	2.28	0.68
1:B:984:GLU:OE2	1:B:1137:LYS:NZ	2.25	0.68
1:B:1021:TYR:CD1	1:B:1098:LYS:HA	2.29	0.68
1:C:1265:ALA:HB2	1:C:1285:LEU:HD23	1.76	0.68
1:B:939:LYS:HG2	1:B:1131:LEU:HD11	1.76	0.67
1:A:1024:LEU:HD11	1:A:1097:ARG:HB2	1.76	0.67
1:A:1181:TYR:CE2	1:C:1353:LEU:HB3	2.29	0.67
1:A:1306:ILE:HD11	1:A:1315:TYR:HB3	1.78	0.66
1:B:1190:TYR:HD2	1:B:1198:LYS:HD3	1.60	0.66
1:B:1302:PHE:HE1	1:B:1320:VAL:HG23	1.61	0.66
1:C:936:ILE:HD11	1:C:1249:LEU:HD13	1.79	0.65
1:A:984:GLU:OE1	1:B:1137:LYS:NZ	2.29	0.64
1:C:1285:LEU:HD12	1:C:1315:TYR:HD2	1.62	0.64
1:B:1190:TYR:CD2	1:B:1198:LYS:HD3	2.34	0.63
1:A:1038:ASP:HB3	1:A:1082:PHE:CE1	2.34	0.63
1:A:1146:ARG:NH1	1:A:1162:ASP:OD2	2.28	0.63
1:A:964:TYR:HE2	1:A:986:PRO:HD2	1.62	0.63
1:A:994:GLY:HA3	1:A:1257:PHE:CD1	2.33	0.63
1:B:1236:TYR:CE1	1:B:1237:LYS:HG3	2.32	0.63
1:C:1265:ALA:HA	1:C:1284:THR:O	1.98	0.63
1:B:1273:ILE:HD11	1:B:1325:PRO:HD2	1.80	0.62
1:B:923:VAL:O	1:B:932:TYR:HE1	1.82	0.62
1:B:958:ILE:HG12	1:B:1011:TYR:CD2	2.34	0.62
1:B:1025:TYR:HB3	1:B:1176:VAL:HG12	1.81	0.61
1:C:966:THR:HG21	1:C:1256:THR:HG21	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1024:LEU:HD11	1:B:1097:ARG:HB2	1.82	0.61
1:A:944:LEU:HD22	1:A:949:TYR:CE2	2.36	0.61
1:B:1087:ASP:OD2	1:B:1206:PHE:HB2	2.01	0.61
1:A:1137:LYS:NZ	1:C:984:GLU:OE1	2.23	0.61
1:A:1302:PHE:HE1	1:A:1320:VAL:HG23	1.65	0.61
1:A:960:PRO:HG2	1:A:1258:ALA:HA	1.82	0.60
1:B:1265:ALA:HB2	1:B:1285:LEU:HD23	1.83	0.60
1:B:994:GLY:HA3	1:B:1257:PHE:CD1	2.37	0.60
1:B:1056:LEU:O	1:B:1084:SER:HA	2.02	0.60
1:B:1295:ILE:HA	1:B:1332:LYS:O	2.03	0.59
1:C:985:VAL:HG21	1:C:993:ILE:HD13	1.84	0.59
1:C:1021:TYR:CD1	1:C:1098:LYS:HA	2.37	0.59
1:B:1049:ILE:HD11	1:B:1059:ALA:HB2	1.85	0.59
1:C:1302:PHE:HE1	1:C:1320:VAL:HG23	1.67	0.58
1:A:962:TYR:HB2	1:A:964:TYR:HE1	1.67	0.58
1:A:994:GLY:HA3	1:A:1257:PHE:CE1	2.38	0.58
1:C:1087:ASP:OD2	1:C:1206:PHE:HB2	2.03	0.58
1:B:964:TYR:CE2	1:B:993:ILE:HG22	2.38	0.57
1:B:994:GLY:HA3	1:B:1257:PHE:CE1	2.39	0.57
1:C:1283:LEU:HD21	1:C:1331:ILE:HD12	1.87	0.57
1:A:1100:THR:HG22	1:C:1302:PHE:CD1	2.39	0.57
1:B:1055:TRP:CD1	1:B:1086:GLN:HB2	2.39	0.57
1:C:1086:GLN:HE21	1:C:1180:ILE:HB	1.70	0.56
1:A:1131:LEU:HB3	1:A:1249:LEU:HD23	1.86	0.56
1:B:988:LEU:HB2	1:B:1239:ILE:HG12	1.87	0.56
1:C:963:LYS:C	1:C:964:TYR:HD1	2.09	0.56
1:A:1302:PHE:CD1	1:B:1100:THR:HG22	2.41	0.56
1:A:1353:LEU:HB3	1:B:1181:TYR:CE2	2.41	0.56
1:C:1102:GLU:HG2	1:C:1144:LEU:HD12	1.89	0.55
1:B:947:PHE:CD2	1:B:1129:THR:HA	2.41	0.55
1:A:1056:LEU:O	1:A:1084:SER:HA	2.06	0.55
1:A:1190:TYR:CE1	1:B:1062:ARG:CZ	2.89	0.55
1:B:1302:PHE:CD1	1:C:1100:THR:HG22	2.42	0.55
1:B:1302:PHE:HD1	1:C:1100:THR:HG22	1.71	0.55
1:C:1024:LEU:HD11	1:C:1097:ARG:HB2	1.88	0.55
1:A:1270:THR:HG22	1:A:1343:ALA:HB3	1.87	0.55
1:B:1078:ASP:OD1	1:B:1079:GLY:N	2.39	0.55
1:A:963:LYS:C	1:A:964:TYR:HD1	2.10	0.55
1:C:962:TYR:HB2	1:C:964:TYR:HE1	1.70	0.55
1:A:1039:GLU:OE2	1:A:1198:LYS:NZ	2.23	0.55
1:B:997:ILE:HG12	1:B:1249:LEU:HD12	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1100:THR:HG22	1:C:1302:PHE:HD1	1.71	0.54
1:C:1014:SER:O	1:C:1106:VAL:HA	2.07	0.54
1:A:1295:ILE:HA	1:A:1332:LYS:O	2.08	0.54
1:C:933:LYS:HG2	1:C:1005:LEU:HG	1.89	0.54
1:C:955:LEU:HB3	1:C:957:HIS:CE1	2.43	0.54
1:B:966:THR:HG22	1:B:983:SER:HB2	1.90	0.54
1:B:1303:HIS:ND1	1:C:1101:GLU:OE1	2.36	0.54
1:A:1082:PHE:HD1	1:A:1082:PHE:H	1.54	0.54
1:A:1273:ILE:HD11	1:A:1324:LYS:HA	1.90	0.54
1:A:1008:VAL:HG12	1:A:1112:PHE:HB3	1.90	0.53
1:C:936:ILE:HG21	1:C:1005:LEU:HD21	1.91	0.53
1:B:960:PRO:HG2	1:B:1258:ALA:HA	1.90	0.53
1:A:1236:TYR:CD1	1:A:1237:LYS:HG3	2.42	0.53
1:C:996:LYS:HG2	1:C:1007:ASP:OD1	2.08	0.53
1:B:918:ILE:HG22	1:B:1130:ASP:OD2	2.09	0.53
1:B:944:LEU:HD22	1:B:949:TYR:CE2	2.44	0.53
1:B:992:SER:HA	1:B:1010:ALA:O	2.08	0.53
1:A:1302:PHE:HD1	1:B:1100:THR:HG22	1.74	0.53
1:A:1110:LEU:O	1:A:1116:THR:HA	2.09	0.53
1:B:1241:GLY:HA2	1:C:1139:VAL:HG12	1.90	0.53
1:A:1096:TYR:CE2	1:A:1231:VAL:HG23	2.44	0.52
1:A:1063:THR:HG21	1:A:1081:VAL:HG23	1.91	0.52
1:A:966:THR:HG22	1:A:983:SER:HB2	1.91	0.52
1:A:1265:ALA:HB2	1:A:1285:LEU:HD23	1.92	0.52
1:A:1351:ILE:HG22	1:A:1352:GLU:HG3	1.92	0.52
1:A:1055:TRP:CD1	1:A:1086:GLN:HB2	2.45	0.52
1:A:999:SER:HB2	1:A:1005:LEU:HD11	1.92	0.51
1:B:1190:TYR:CE1	1:C:1062:ARG:CZ	2.92	0.51
1:C:960:PRO:HG2	1:C:1258:ALA:HA	1.92	0.51
1:A:1049:ILE:HD11	1:A:1059:ALA:HB2	1.92	0.51
1:A:1241:GLY:HA2	1:B:1139:VAL:HG12	1.93	0.51
1:A:966:THR:HG21	1:A:1256:THR:HG21	1.92	0.51
1:C:1049:ILE:HD11	1:C:1059:ALA:HB2	1.92	0.51
1:C:1063:THR:HB	1:C:1190:TYR:HE2	1.75	0.51
1:A:1181:TYR:HE2	1:C:1353:LEU:HB3	1.74	0.51
1:B:1269:CYS:O	1:B:1342:MET:HA	2.10	0.51
1:B:1285:LEU:HD12	1:B:1315:TYR:HD2	1.75	0.51
1:B:1008:VAL:HG12	1:B:1112:PHE:HB3	1.91	0.51
1:C:1024:LEU:HD22	1:C:1163:LEU:HB2	1.93	0.51
1:C:949:TYR:CD1	1:C:1119:THR:HB	2.45	0.51
1:C:1055:TRP:NE1	1:C:1086:GLN:OE1	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1236:TYR:CE1	1:A:1237:LYS:HG3	2.45	0.50
1:B:1265:ALA:HA	1:B:1284:THR:O	2.10	0.50
1:C:937:THR:HG23	1:C:1112:PHE:CD1	2.46	0.50
1:B:949:TYR:CD1	1:B:1119:THR:HB	2.46	0.50
1:B:1063:THR:HB	1:B:1190:TYR:HE2	1.76	0.50
1:C:1008:VAL:HG12	1:C:1112:PHE:HB3	1.94	0.50
1:B:1063:THR:HB	1:B:1190:TYR:CE2	2.46	0.50
1:C:1021:TYR:CE2	1:C:1233:ILE:HD12	2.47	0.50
1:B:1121:LEU:HD23	1:B:1127:ILE:HD12	1.94	0.49
1:A:1068:CYS:SG	1:A:1198:LYS:HD3	2.52	0.49
1:C:996:LYS:HG3	1:C:1257:PHE:CD2	2.47	0.49
1:A:1285:LEU:HD12	1:A:1315:TYR:HD2	1.77	0.49
1:A:1134:VAL:HG12	1:A:1245:MET:HG2	1.94	0.49
1:C:1005:LEU:HD23	1:C:1006:LEU:HB2	1.94	0.49
1:A:1190:TYR:HD1	1:A:1190:TYR:H	1.61	0.49
1:B:1134:VAL:HG12	1:B:1245:MET:HG2	1.93	0.49
1:A:1062:ARG:CZ	1:C:1190:TYR:CE1	2.96	0.49
1:A:1141:THR:O	1:B:1141:THR:HG21	2.13	0.49
1:A:1188:PHE:HE1	1:A:1200:VAL:HG13	1.78	0.49
1:C:937:THR:O	1:C:941:GLN:HB2	2.12	0.49
1:A:1190:TYR:CD2	1:A:1198:LYS:HD2	2.48	0.49
1:B:958:ILE:HG21	1:B:1011:TYR:CE2	2.48	0.49
1:B:1224:LEU:HD23	1:B:1233:ILE:HG12	1.94	0.49
1:B:1204:LYS:HD2	1:B:1357:ASP:O	2.13	0.49
1:C:1063:THR:HB	1:C:1198:LYS:HE3	1.94	0.48
1:C:1190:TYR:CD2	1:C:1198:LYS:HD2	2.48	0.48
1:C:1052:GLN:HB2	1:C:1055:TRP:CD1	2.48	0.48
1:C:1110:LEU:HD11	1:C:1132:PHE:CE2	2.47	0.48
1:C:1056:LEU:O	1:C:1084:SER:HA	2.13	0.48
1:C:1183:ASN:OD1	1:C:1184:GLY:N	2.47	0.48
1:A:1188:PHE:CE1	1:A:1200:VAL:HG13	2.49	0.48
1:B:1269:CYS:SG	1:B:1329:LEU:HD23	2.54	0.48
1:A:1265:ALA:HA	1:A:1284:THR:O	2.13	0.48
1:A:962:TYR:HB2	1:A:964:TYR:CE1	2.48	0.48
1:A:1266:GLU:HG2	1:A:1284:THR:HB	1.95	0.48
1:B:997:ILE:O	1:B:1005:LEU:HB2	2.13	0.48
1:A:1062:ARG:CZ	1:C:1190:TYR:HE1	2.26	0.48
1:B:1285:LEU:HD12	1:B:1315:TYR:CD2	2.50	0.47
1:B:1086:GLN:NE2	1:B:1180:ILE:HB	2.27	0.47
1:C:1020:VAL:HG12	1:C:1232:THR:OG1	2.14	0.47
1:A:1046:PRO:HG2	1:A:1082:PHE:HE2	1.76	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:2:NAG:H61	2:E:3:FUC:H61	1.97	0.47
1:B:933:LYS:HA	1:B:936:ILE:HG22	1.97	0.47
1:B:1328:THR:OG1	1:B:1339:GLU:OE1	2.33	0.47
1:C:1089:ILE:HG12	1:C:1180:ILE:HD13	1.95	0.47
1:B:990:GLY:HA3	1:B:1307:LEU:HD11	1.96	0.47
1:B:1302:PHE:CD2	1:B:1318:LYS:HB2	2.50	0.47
1:A:1122:ASN:ND2	1:A:1124:VAL:HB	2.30	0.47
1:B:966:THR:HG21	1:B:1256:THR:HG21	1.97	0.47
1:C:947:PHE:CD2	1:C:1129:THR:HG22	2.50	0.47
1:B:1168:LYS:HB3	1:B:1209:ASP:OD2	2.15	0.47
1:C:1063:THR:HG21	1:C:1081:VAL:HG23	1.97	0.46
1:A:937:THR:O	1:A:941:GLN:HB2	2.16	0.46
1:B:1159:GLN:HG3	1:C:1146:ARG:HH22	1.81	0.46
1:A:1082:PHE:CD1	1:A:1082:PHE:N	2.81	0.46
1:A:1096:TYR:CD1	1:A:1096:TYR:N	2.84	0.46
1:B:932:TYR:CE2	1:B:1005:LEU:HD21	2.50	0.46
1:C:1313:HIS:CD2	1:C:1314:LYS:HG3	2.50	0.46
1:A:940:LEU:HA	1:A:1131:LEU:HD12	1.97	0.46
1:B:1300:THR:HG22	1:C:1100:THR:HG21	1.97	0.46
1:C:964:TYR:CD1	1:C:964:TYR:N	2.84	0.46
1:C:1003:LYS:HA	1:C:1004:HIS:HA	1.53	0.46
1:B:1030:THR:HA	1:B:1089:ILE:HG22	1.98	0.46
1:B:1110:LEU:HD11	1:B:1132:PHE:CE2	2.51	0.46
1:A:1179:THR:HG1	1:A:1181:TYR:HE1	1.63	0.45
1:A:1278:ASN:HB3	1:A:1319:MET:O	2.17	0.45
1:B:1190:TYR:HD1	1:B:1190:TYR:H	1.63	0.45
1:B:967:ILE:HD11	1:C:982:GLU:HG3	1.99	0.45
1:A:1285:LEU:HB2	1:A:1315:TYR:HD2	1.82	0.45
1:A:1003:LYS:HA	1:A:1004:HIS:HA	1.59	0.45
1:A:1096:TYR:N	1:A:1096:TYR:HD1	2.14	0.45
1:C:1018:SER:HA	1:C:1233:ILE:O	2.16	0.45
1:B:1277:GLU:OE2	1:B:1348:LYS:HD2	2.16	0.45
1:A:1098:LYS:NZ	1:A:1102:GLU:OE2	2.33	0.45
1:A:1102:GLU:HG3	1:A:1144:LEU:HD12	1.97	0.45
1:A:1285:LEU:HB2	1:A:1315:TYR:CD2	2.52	0.45
1:C:1039:GLU:OE2	1:C:1198:LYS:NZ	2.28	0.45
1:A:1131:LEU:HD22	1:A:1249:LEU:HA	1.99	0.45
1:A:1273:ILE:CD1	1:A:1324:LYS:HA	2.47	0.45
1:B:964:TYR:CD2	1:B:993:ILE:HG22	2.52	0.45
1:C:1331:ILE:HG13	1:C:1338:VAL:HB	1.98	0.45
1:B:947:PHE:HD2	1:B:1129:THR:HA	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:997:ILE:O	1:C:1005:LEU:HB3	2.16	0.44
1:A:1141:THR:HG21	1:C:1141:THR:O	2.17	0.44
1:B:1183:ASN:OD1	1:B:1184:GLY:N	2.49	0.44
1:C:1131:LEU:HD22	1:C:1249:LEU:HA	1.98	0.44
1:A:1194:LEU:HD12	1:A:1195:ALA:N	2.33	0.44
1:B:1131:LEU:HB3	1:B:1249:LEU:HD23	1.98	0.44
1:C:1225:GLU:HB2	1:C:1232:THR:HB	1.99	0.44
1:A:933:LYS:HA	1:A:936:ILE:HG22	1.99	0.44
1:A:1121:LEU:HD23	1:A:1127:ILE:HD12	2.00	0.44
1:B:1236:TYR:CE1	1:B:1237:LYS:HE3	2.53	0.44
1:A:947:PHE:CE2	1:A:1129:THR:HA	2.53	0.44
1:A:1087:ASP:OD2	1:A:1206:PHE:HB2	2.17	0.44
1:A:1244:LYS:HE3	1:C:965:ILE:HB	2.00	0.44
1:A:1158:GLY:HA3	1:A:1216:LEU:HD22	2.00	0.44
1:A:1265:ALA:HB3	1:A:1338:VAL:HG22	2.00	0.44
1:B:1014:SER:O	1:B:1106:VAL:HA	2.18	0.44
1:B:1024:LEU:HD22	1:B:1163:LEU:HB2	1.99	0.44
1:A:1181:TYR:HD1	1:A:1181:TYR:N	2.15	0.43
1:C:937:THR:HG23	1:C:1112:PHE:HD1	1.83	0.43
1:A:1144:LEU:HD23	1:A:1144:LEU:HA	1.86	0.43
1:A:1190:TYR:HE1	1:B:1062:ARG:CZ	2.31	0.43
1:B:1006:LEU:HD11	1:B:1112:PHE:HB2	1.99	0.43
1:A:992:SER:HA	1:A:1010:ALA:O	2.18	0.43
1:A:1181:TYR:N	1:A:1181:TYR:CD1	2.86	0.43
1:C:1190:TYR:HD1	1:C:1190:TYR:H	1.65	0.43
1:A:1020:VAL:C	1:A:1021:TYR:HD1	2.22	0.43
1:A:1128:ILE:HG23	1:A:1133:GLU:HG2	2.00	0.43
1:A:1302:PHE:CD2	1:A:1318:LYS:HB2	2.53	0.43
1:A:1021:TYR:CE2	1:A:1233:ILE:HD12	2.53	0.43
1:A:1100:THR:HG21	1:C:1300:THR:HG22	1.99	0.43
1:C:1016:SER:HA	1:C:1235:ASP:O	2.18	0.43
1:B:1011:TYR:CD1	1:B:1109:CYS:HB2	2.51	0.43
1:A:949:TYR:HE1	1:A:1127:ILE:HG21	1.83	0.43
1:A:1300:THR:HG22	1:B:1100:THR:HG21	2.00	0.43
1:A:1353:LEU:HD23	1:B:1181:TYR:CE2	2.54	0.43
1:A:964:TYR:CD1	1:A:964:TYR:N	2.87	0.43
1:B:992:SER:HB3	1:B:1011:TYR:HB3	2.01	0.43
1:B:1146:ARG:NH1	1:B:1162:ASP:OD2	2.45	0.43
1:C:1168:LYS:HB3	1:C:1209:ASP:OD2	2.19	0.43
1:B:1003:LYS:HA	1:B:1004:HIS:HA	1.42	0.43
1:B:1192:CYS:SG	1:C:1077:SER:HB3	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1236:TYR:CE1	1:C:1237:LYS:HE3	2.54	0.43
1:B:947:PHE:CE2	1:B:1129:THR:HA	2.54	0.42
1:C:1110:LEU:O	1:C:1116:THR:HA	2.19	0.42
1:A:1062:ARG:NE	1:C:1190:TYR:CE1	2.87	0.42
1:A:1190:TYR:HE1	1:B:1062:ARG:NH2	2.16	0.42
1:B:949:TYR:HE1	1:B:1127:ILE:HG21	1.85	0.42
1:B:1181:TYR:N	1:B:1181:TYR:HD1	2.17	0.42
1:C:1190:TYR:HD2	1:C:1198:LYS:HD2	1.84	0.42
1:B:982:GLU:HA	1:B:1245:MET:O	2.19	0.42
1:C:1129:THR:OG1	1:C:1132:PHE:N	2.50	0.42
1:A:1273:ILE:HD11	1:A:1325:PRO:HD2	2.01	0.42
1:C:1144:LEU:HD23	1:C:1144:LEU:HA	1.85	0.42
1:C:954:ASN:ND2	1:C:1114:ASP:HA	2.35	0.42
1:A:1199:GLU:OE1	1:A:1363:ARG:NE	2.42	0.42
1:C:1329:LEU:O	1:C:1340:ALA:HB3	2.20	0.42
1:A:939:LYS:HG2	1:A:1131:LEU:HD11	2.00	0.42
1:B:1303:HIS:HD2	1:B:1305:ARG:O	2.03	0.42
1:B:1350:ILE:HG21	1:B:1353:LEU:HD22	2.01	0.42
1:C:982:GLU:HA	1:C:1245:MET:O	2.20	0.42
1:A:1014:SER:O	1:A:1106:VAL:HA	2.20	0.41
1:B:1060:ARG:NH1	1:B:1189:ASP:OD1	2.52	0.41
1:A:1328:THR:OG1	1:A:1339:GLU:OE1	2.38	0.41
1:A:1283:LEU:HD21	1:A:1331:ILE:HD11	2.02	0.41
1:B:932:TYR:HE2	1:B:1005:LEU:HD21	1.83	0.41
2:D:1:NAG:O3	2:D:2:NAG:N2	2.53	0.41
1:A:964:TYR:HD1	1:A:964:TYR:N	2.18	0.41
1:B:1181:TYR:N	1:B:1181:TYR:CD1	2.88	0.41
1:B:920:LYS:O	1:B:976:ILE:HD11	2.20	0.41
1:B:1159:GLN:HG3	1:C:1146:ARG:NH2	2.35	0.41
1:A:1168:LYS:HB3	1:A:1209:ASP:OD2	2.20	0.41
1:B:1017:TYR:HA	1:B:1103:VAL:O	2.21	0.41
1:B:1141:THR:O	1:C:1141:THR:HG21	2.20	0.41
1:A:1269:CYS:O	1:A:1342:MET:HA	2.20	0.41
1:C:964:TYR:HD1	1:C:964:TYR:N	2.18	0.41
1:B:1011:TYR:HE1	1:B:1109:CYS:HB2	1.79	0.41
1:A:1271:GLY:HA3	1:A:1279:ILE:HA	2.03	0.41
1:B:1190:TYR:HE1	1:C:1062:ARG:CZ	2.34	0.41
1:C:933:LYS:HE2	1:C:1005:LEU:HA	2.02	0.41
1:C:1133:GLU:OE1	1:C:1246:LYS:HD3	2.21	0.41
1:C:1134:VAL:HG12	1:C:1245:MET:HG2	2.03	0.41
1:A:1029:PRO:HD3	1:A:1210:TYR:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1190:TYR:N	1:A:1190:TYR:CD1	2.88	0.41
1:A:1313:HIS:CD2	1:A:1314:LYS:HG3	2.55	0.41
1:B:1161:ASN:ND2	1:B:1170:CYS:O	2.54	0.41
1:C:955:LEU:HD22	1:C:958:ILE:HD11	2.04	0.40
1:B:947:PHE:HD2	1:B:1129:THR:HG22	1.85	0.40
1:B:965:ILE:HB	1:C:1244:LYS:HE3	2.03	0.40
1:A:1080:CYS:SG	1:A:1081:VAL:N	2.94	0.40
1:A:1285:LEU:HD12	1:A:1315:TYR:CD2	2.56	0.40
1:C:941:GLN:HG3	1:C:1117:TYR:CE2	2.56	0.40
1:C:1102:GLU:OE2	1:C:1144:LEU:HB2	2.21	0.40
1:A:932:TYR:HE2	1:A:1005:LEU:HD21	1.87	0.40
1:A:1308:VAL:HG12	1:A:1315:TYR:CE2	2.57	0.40
1:B:1011:TYR:H	1:B:1011:TYR:HD1	1.69	0.40
1:B:1068:CYS:O	1:B:1198:LYS:HE3	2.22	0.40
1:C:1021:TYR:O	1:C:1230:THR:HA	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	435/458 (95%)	427 (98%)	8 (2%)	0	100	100
1	B	447/458 (98%)	438 (98%)	9 (2%)	0	100	100
1	C	434/458 (95%)	425 (98%)	9 (2%)	0	100	100
All	All	1316/1374 (96%)	1290 (98%)	26 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	386/404 (96%)	378 (98%)	8 (2%)	53 78
1	B	395/404 (98%)	390 (99%)	5 (1%)	69 86
1	C	385/404 (95%)	382 (99%)	3 (1%)	81 92
All	All	1166/1212 (96%)	1150 (99%)	16 (1%)	67 85

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	964	TYR
1	A	1021	TYR
1	A	1082	PHE
1	A	1096	TYR
1	A	1181	TYR
1	A	1190	TYR
1	A	1236	TYR
1	A	1257	PHE
1	B	1011	TYR
1	B	1181	TYR
1	B	1190	TYR
1	B	1236	TYR
1	B	1257	PHE
1	C	964	TYR
1	C	1190	TYR
1	C	1236	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1086	GLN
1	C	1086	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

9 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	NAG	D	1	2,1	14,14,15	0.63	1 (7%)	17,19,21	0.52	0
2	NAG	D	2	2	14,14,15	0.51	0	17,19,21	0.57	0
2	FUC	D	3	2	10,10,11	1.23	2 (20%)	14,14,16	1.34	3 (21%)
2	NAG	E	1	2,1	14,14,15	0.32	0	17,19,21	0.48	0
2	NAG	E	2	2	14,14,15	0.57	0	17,19,21	0.37	0
2	FUC	E	3	2	10,10,11	1.01	0	14,14,16	1.01	0
2	NAG	F	1	2,1	14,14,15	0.32	0	17,19,21	0.55	0
2	NAG	F	2	2	14,14,15	0.56	0	17,19,21	0.56	0
2	FUC	F	3	2	10,10,11	1.00	0	14,14,16	1.08	1 (7%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	D	1	2,1	-	1/6/23/26	0/1/1/1
2	NAG	D	2	2	-	4/6/23/26	0/1/1/1
2	FUC	D	3	2	-	-	0/1/1/1
2	NAG	E	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	NAG	O5-C1	-2.25	1.40	1.43
2	D	3	FUC	C1-C2	2.21	1.57	1.52
2	D	3	FUC	C2-C3	2.03	1.55	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	3	FUC	C1-C2-C3	2.72	113.01	109.67
2	D	3	FUC	O5-C5-C4	2.40	113.82	109.52
2	D	3	FUC	C1-O5-C5	2.26	117.91	112.78
2	F	3	FUC	C1-O5-C5	2.11	117.56	112.78

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	2	NAG	O5-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C1-C2-N2-C7
2	D	2	NAG	C3-C2-N2-C7
2	F	2	NAG	C3-C2-N2-C7
2	D	2	NAG	C1-C2-N2-C7
2	D	1	NAG	O5-C5-C6-O6

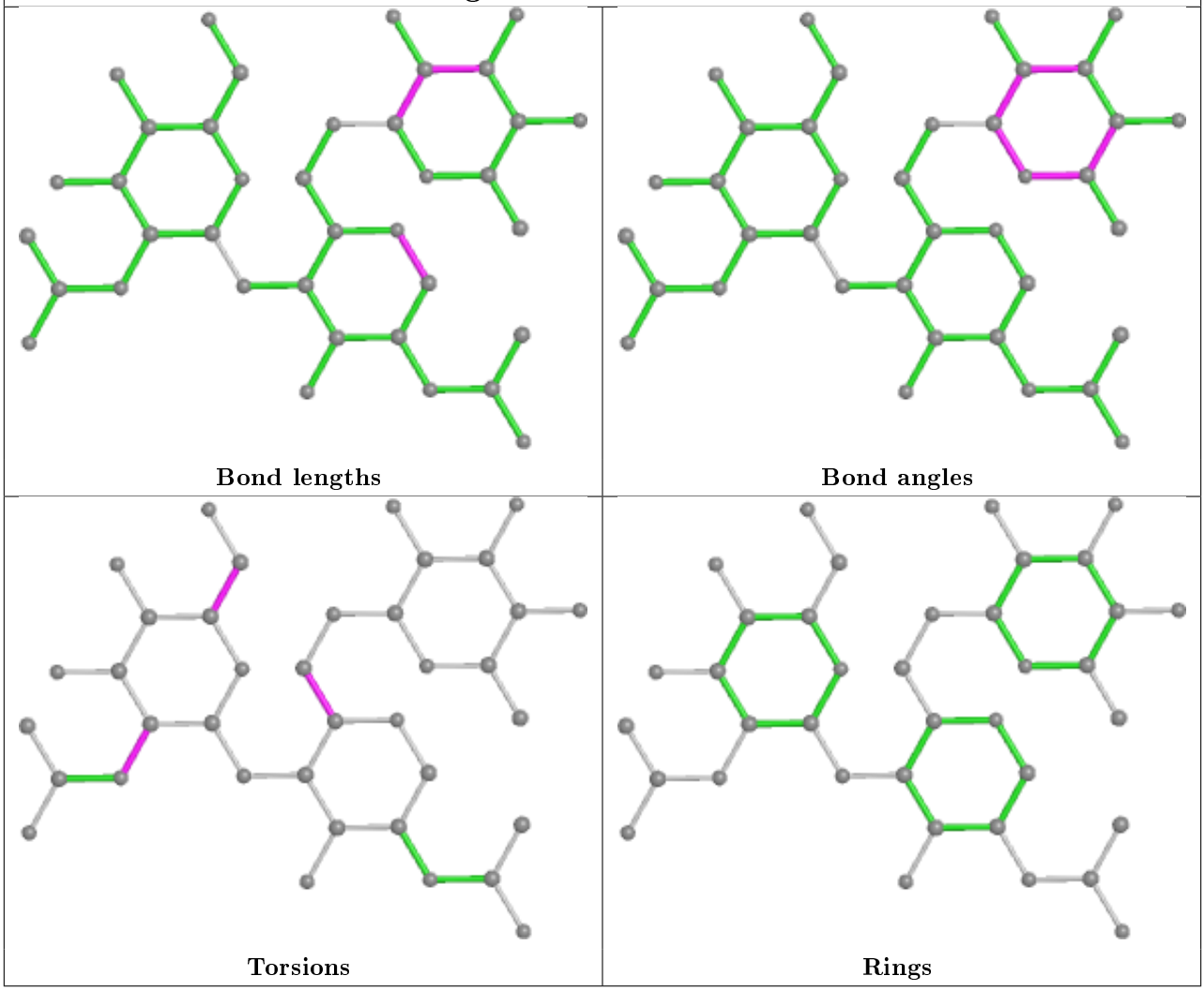
There are no ring outliers.

4 monomers are involved in 2 short contacts:

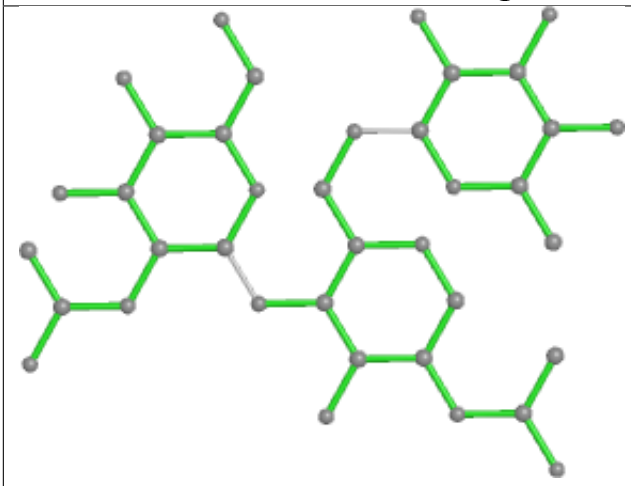
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	NAG	1	0
2	E	2	NAG	1	0
2	E	3	FUC	1	0
2	D	2	NAG	1	0

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

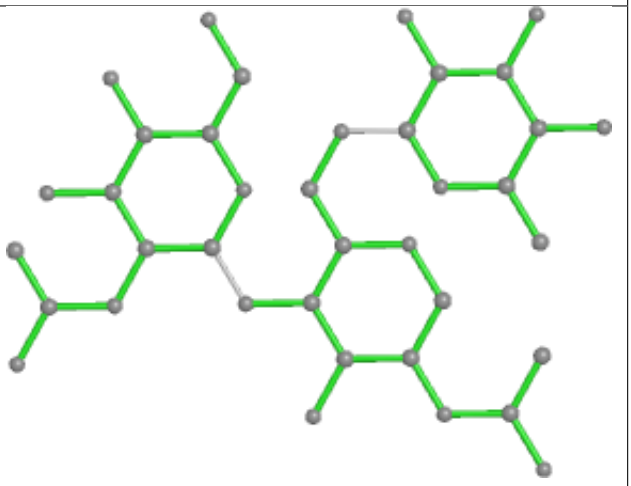
Oligosaccharide Chain D



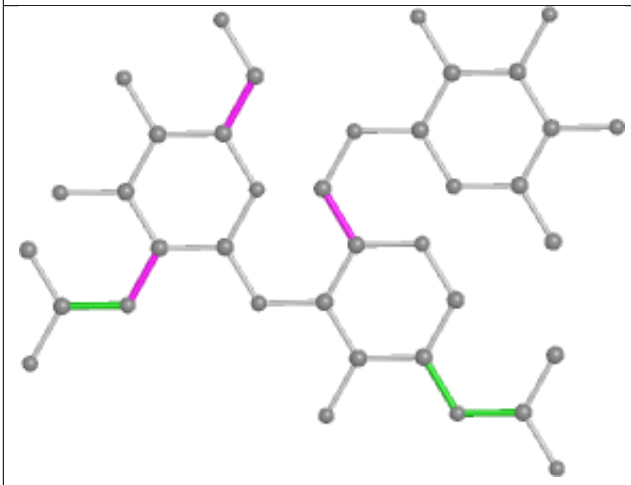
Oligosaccharide Chain E



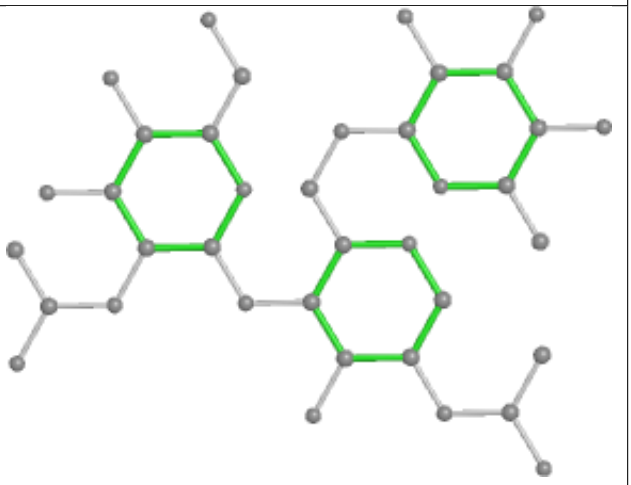
Bond lengths



Bond angles

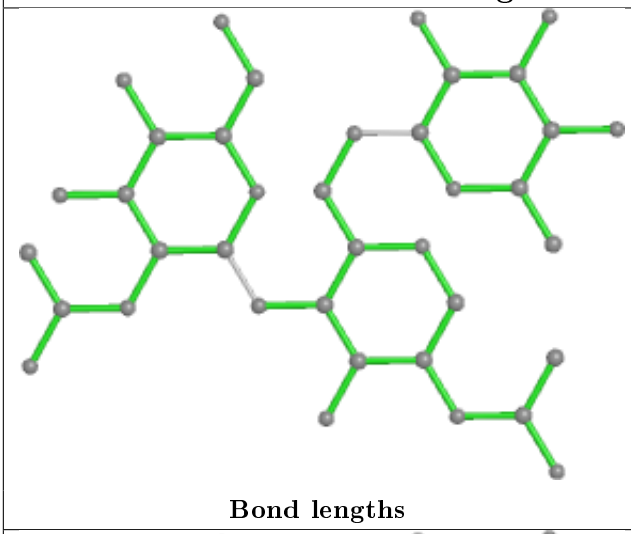


Torsions

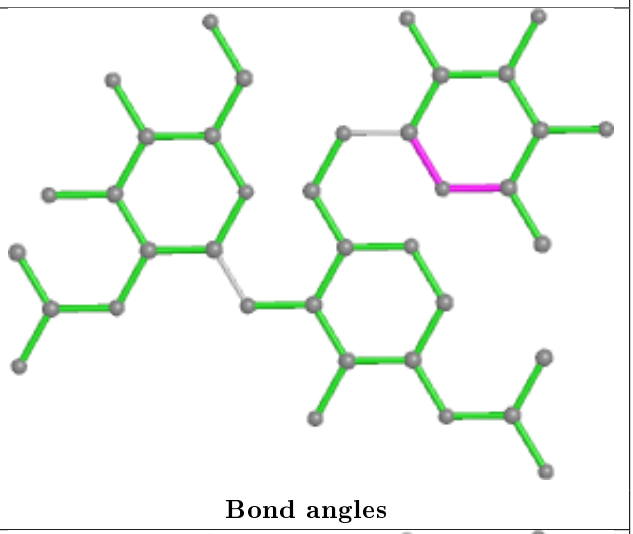


Rings

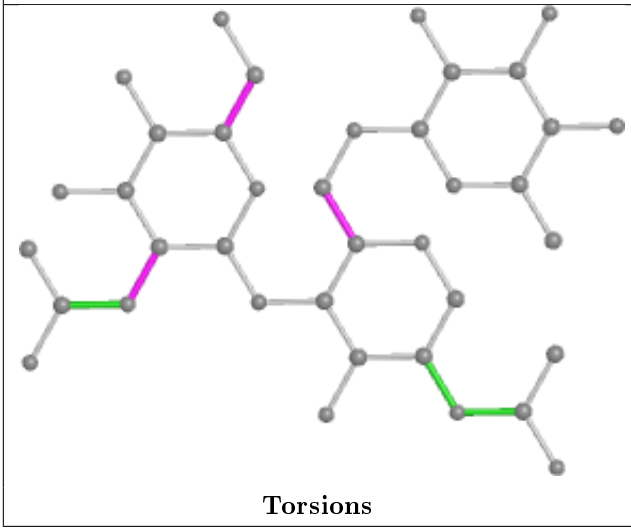
Oligosaccharide Chain F



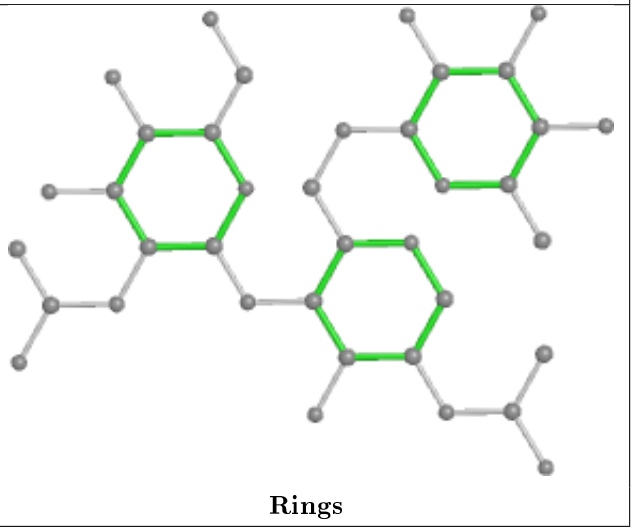
Bond lengths



Bond angles



Torsions



Rings

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, 95^{th} 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	437/458 (95%)	-0.06	7 (1%)	72	59	57, 92, 154, 208	0
1	B	449/458 (98%)	-0.03	7 (1%)	72	59	58, 102, 151, 209	0
1	C	436/458 (95%)	0.13	23 (5%)	26	13	69, 115, 191, 234	0
All	All	1322/1374 (96%)	0.01	37 (2%)	53	36	57, 103, 170, 234	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1195	ALA	4.0
1	C	975	GLY	3.9
1	A	975	GLY	3.7
1	B	1285	LEU	3.6
1	C	974	GLU	3.6
1	B	1331	ILE	3.4
1	B	1001	GLU	3.3
1	A	974	GLU	3.2
1	B	1329	LEU	3.1
1	C	954	ASN	3.0
1	C	1266	GLU	2.9
1	A	946	LEU	2.9
1	B	1003	LYS	2.9
1	C	930	GLU	2.5
1	C	931	GLN	2.5
1	C	1001	GLU	2.5
1	C	1285	LEU	2.4
1	C	1143	SER	2.4
1	C	1284	THR	2.4
1	A	976	ILE	2.4
1	C	1286	HIS	2.3
1	C	932	TYR	2.3
1	C	1339	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	1342	MET	2.3
1	B	1197	ARG	2.2
1	C	1328	THR	2.2
1	A	1071	PHE	2.2
1	C	1291	ALA	2.2
1	C	1117	TYR	2.1
1	B	1286	HIS	2.1
1	A	1143	SER	2.1
1	C	976	ILE	2.1
1	C	1289	ILE	2.1
1	C	1340	ALA	2.0
1	C	1288	THR	2.0
1	C	973	ALA	2.0
1	C	1329	LEU	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)

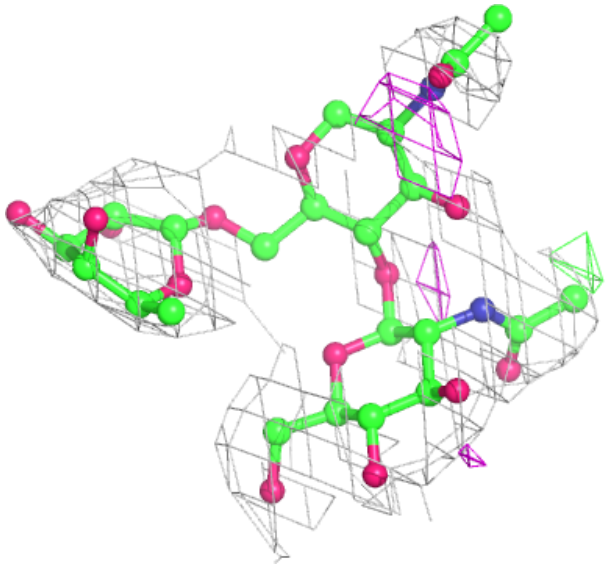
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

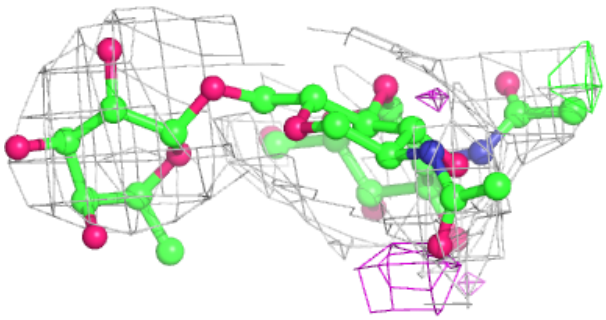
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	NAG	F	2	14/15	0.79	0.46	162,172,182,182	0
2	NAG	E	2	14/15	0.84	0.32	153,174,183,184	0
2	FUC	F	3	10/11	0.88	0.59	147,158,163,164	0
2	NAG	D	2	14/15	0.89	0.32	132,151,154,156	0
2	NAG	F	1	14/15	0.92	0.25	101,130,157,168	0
2	FUC	D	3	10/11	0.92	0.51	152,157,165,167	0
2	FUC	E	3	10/11	0.92	0.50	143,162,165,169	0
2	NAG	E	1	14/15	0.94	0.15	89,98,152,153	0
2	NAG	D	1	14/15	0.94	0.21	98,110,139,147	0

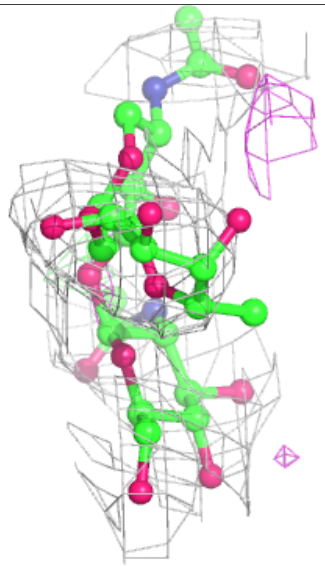
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain D:

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

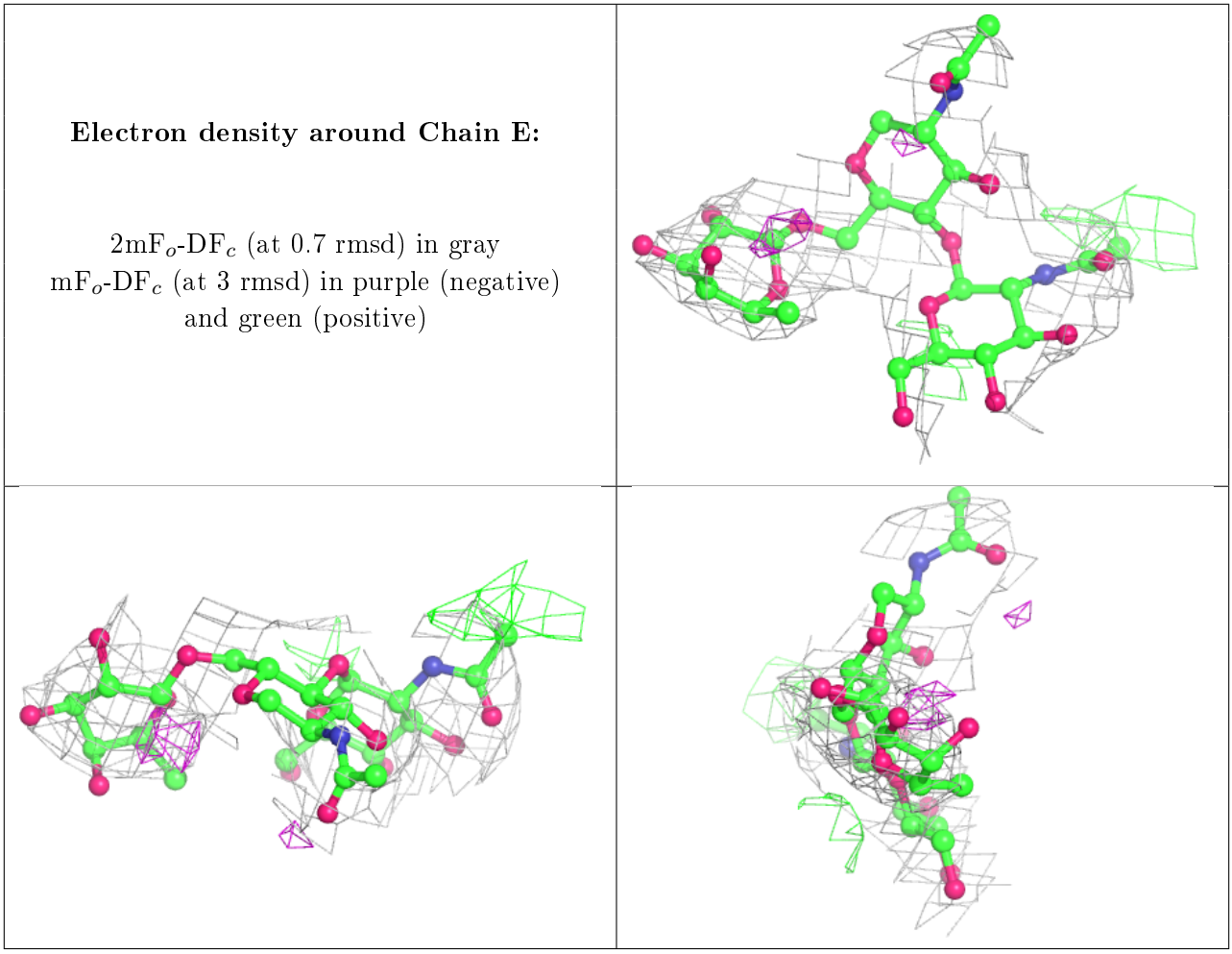






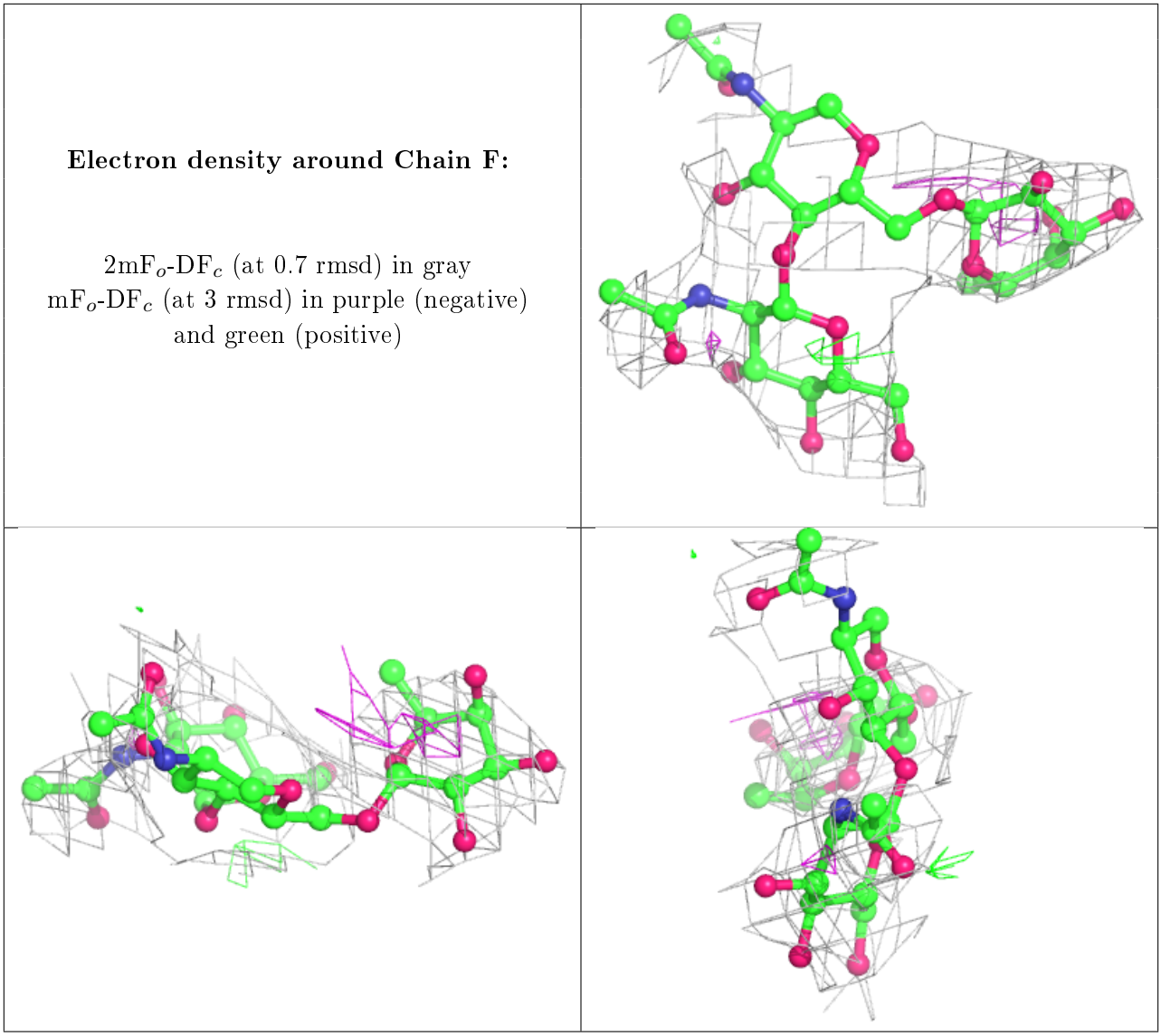
Electron density around Chain E:

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



Electron density around Chain F:

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



6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.