



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2025 – 12:37 AM JST

PDB ID : 9JLS / pdb_00009jls
Title : Crystal structure of GH57 family amylopullulanase from Aquifex aeolicus wild type in complex with acarbose
Authors : Zhu, Z.M.; Wang, W.W.; Li, M.J.; Xu, Q.; Zhou, H.; Huang, L.Q.; Wang, Q.S.; Yu, F.
Deposited on : 2024-09-19
Resolution : 3.06 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

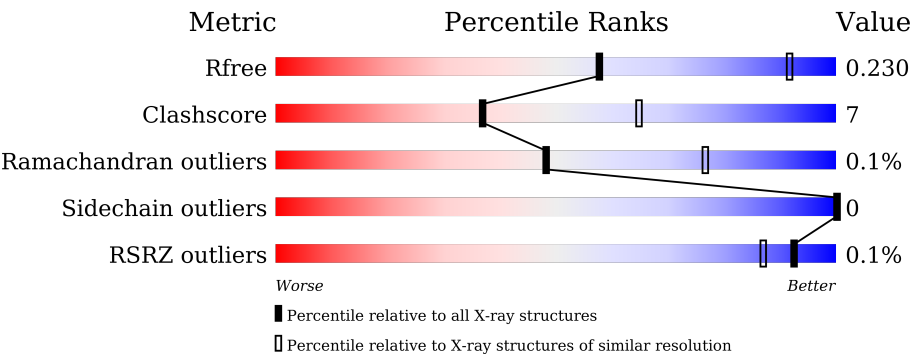
MolProbity : 4-5-2 with Phenix2.0rc1
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

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 3.06 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.




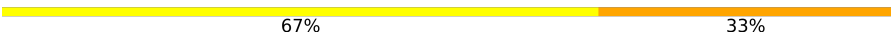
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	2258 (3.10-3.02)
Clashscore	180529	2399 (3.10-3.02)
Ramachandran outliers	177936	2269 (3.10-3.02)
Sidechain outliers	177891	2268 (3.10-3.02)
RSRZ outliers	164620	2258 (3.10-3.02)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	484	<div><div>83%</div><div>16%</div><div>.</div></div>
1	B	484	<div><div>80%</div><div>18%</div><div>.</div></div>
1	C	484	<div><div>79%</div><div>21%</div></div>
1	D	484	<div><div>80%</div><div>20%</div></div>
2	E	3	<div><div>67%</div><div>33%</div></div>
2	F	3	<div><div>33%</div><div>67%</div></div>

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Mol	Chain	Length	Quality of chain
2	G	3	 33%67%
2	H	3	 67%33%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16518 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 Glycoside hydrolase family 57 N-terminal domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	477	Total	C	N	O	S	0	0	0
			4052	2651	648	743	10			
1	B	477	Total	C	N	O	S	0	0	0
			4052	2651	648	743	10			
1	C	484	Total	C	N	O	S	0	0	0
			4119	2692	665	752	10			
1	D	484	Total	C	N	O	S	0	0	0
			4119	2692	665	752	10			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	478	LEU	-	expression tag	UNP O66934
A	479	GLU	-	expression tag	UNP O66934
A	480	HIS	-	expression tag	UNP O66934
A	481	HIS	-	expression tag	UNP O66934
A	482	HIS	-	expression tag	UNP O66934
A	483	HIS	-	expression tag	UNP O66934
A	484	HIS	-	expression tag	UNP O66934
B	478	LEU	-	expression tag	UNP O66934
B	479	GLU	-	expression tag	UNP O66934
B	480	HIS	-	expression tag	UNP O66934
B	481	HIS	-	expression tag	UNP O66934
B	482	HIS	-	expression tag	UNP O66934
B	483	HIS	-	expression tag	UNP O66934
B	484	HIS	-	expression tag	UNP O66934
C	478	LEU	-	expression tag	UNP O66934
C	479	GLU	-	expression tag	UNP O66934
C	480	HIS	-	expression tag	UNP O66934
C	481	HIS	-	expression tag	UNP O66934
C	482	HIS	-	expression tag	UNP O66934
C	483	HIS	-	expression tag	UNP O66934

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Chain	Residue	Modelled	Actual	Comment	Reference
C	484	HIS	-	expression tag	UNP O66934
D	478	LEU	-	expression tag	UNP O66934
D	479	GLU	-	expression tag	UNP O66934
D	480	HIS	-	expression tag	UNP O66934
D	481	HIS	-	expression tag	UNP O66934
D	482	HIS	-	expression tag	UNP O66934
D	483	HIS	-	expression tag	UNP O66934
D	484	HIS	-	expression tag	UNP O66934

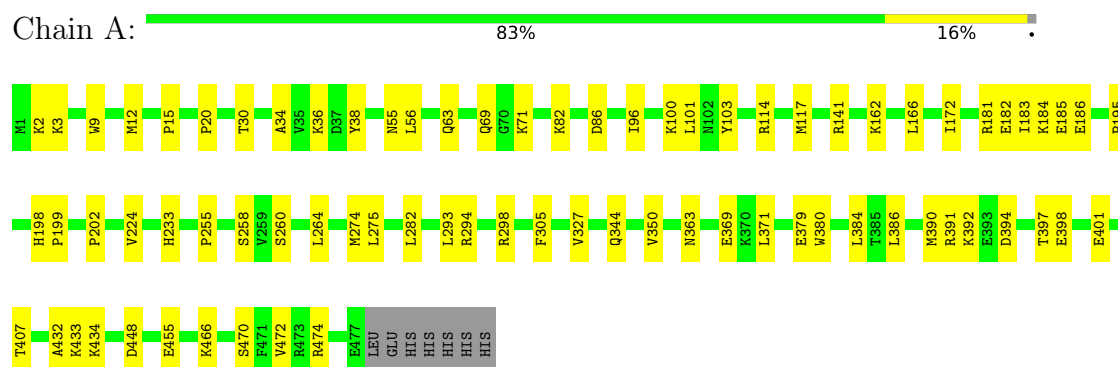
- Molecule 2 is an oligosaccharide called 4,6-dideoxy-4-([(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	0	0	0
			44	25	1	18			
2	F	3	Total	C	N	O	0	0	0
			44	25	1	18			
2	G	3	Total	C	N	O	0	0	0
			44	25	1	18			
2	H	3	Total	C	N	O	0	0	0
			44	25	1	18			

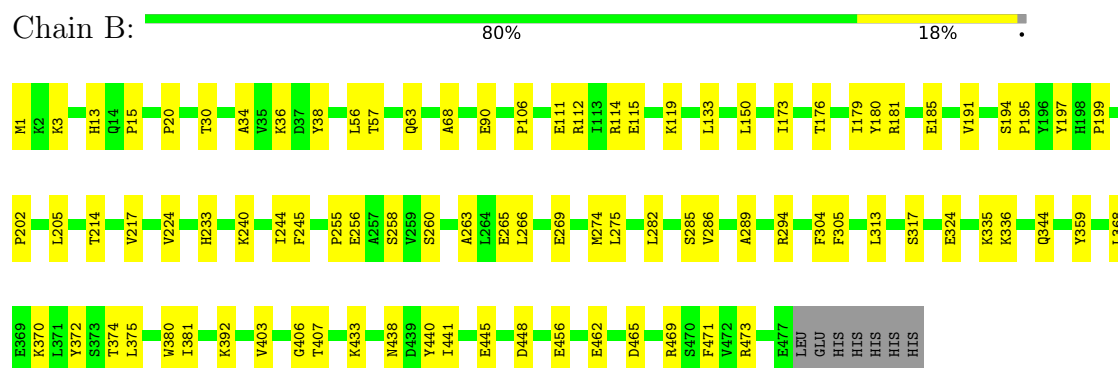
3 Residue-property plots

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

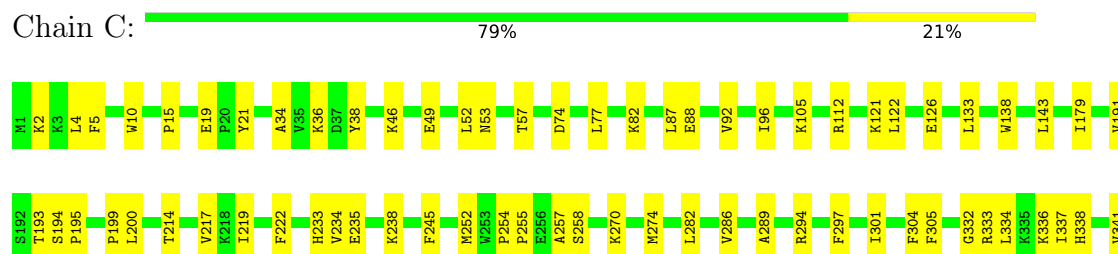
- Molecule 1: Glycoside hydrolase family 57 N-terminal domain-containing protein

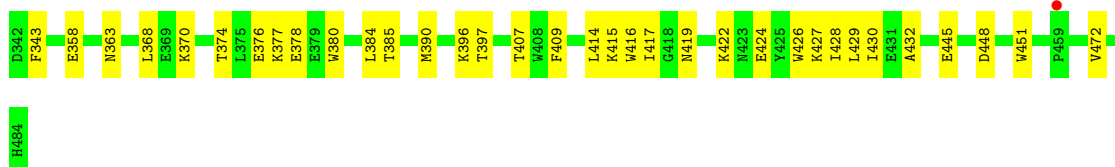


- Molecule 1: Glycoside hydrolase family 57 N-terminal domain-containing protein



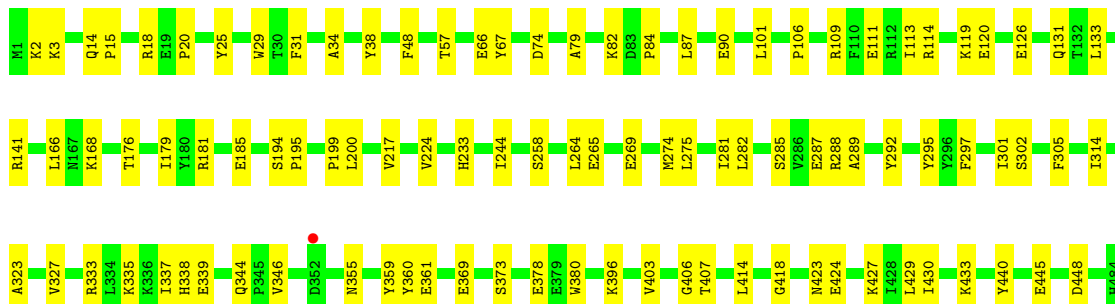
- Molecule 1: Glycoside hydrolase family 57 N-terminal domain-containing protein





- Molecule 1: Glycoside hydrolase family 57 N-terminal domain-containing protein

Chain D: 80% 20%



- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain E: 67% 33%



- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain F: 33% 67%



- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain G: 33% 67%



- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain H: 67% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.43Å 167.41Å 97.98Å 90.00° 109.38° 90.00°	Depositor
Resolution (Å)	66.50 – 3.06 66.50 – 3.06	Depositor EDS
% Data completeness (in resolution range)	99.2 (66.50-3.06) 99.3 (66.50-3.06)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 3.07Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.180 , 0.230 0.180 , 0.230	Depositor DCC
R_{free} test set	2079 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	57.3	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.038 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16518	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GLC, AC1

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	0.11	0/4165	0.27	0/5633
1	B	0.11	0/4165	0.26	0/5633
1	C	0.11	0/4237	0.26	0/5731
1	D	0.11	0/4237	0.28	0/5731
All	All	0.11	0/16804	0.27	0/22728

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

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	4052	0	3987	47	0
1	B	4052	0	3987	56	0
1	C	4119	0	4039	63	0
1	D	4119	0	4039	62	0
2	E	44	0	30	1	0
2	F	44	0	30	2	0
2	G	44	0	30	0	0
2	H	44	0	30	3	0
All	All	16518	0	16172	229	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (229) 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:C:34:ALA:HA	1:C:38:TYR:HB2	1.66	0.78
1:C:195:PRO:HB3	1:C:233:HIS:HB3	1.65	0.77
1:A:96:ILE:HG23	1:A:117:MET:HE1	1.74	0.70
1:D:258:SER:HB3	1:D:407:THR:HG22	1.76	0.67
1:D:195:PRO:HB3	1:D:233:HIS:HB3	1.79	0.65
1:D:34:ALA:HA	1:D:38:TYR:HB2	1.79	0.64
1:D:141:ARG:NH1	1:D:369:GLU:OE2	2.31	0.63
1:A:390:MET:HE3	1:A:391:ARG:HH12	1.63	0.62
1:B:90:GLU:OE1	1:B:90:GLU:N	2.29	0.61
1:D:101:LEU:HD11	2:H:3:AC1:HCB2	1.83	0.61
1:A:100:LYS:HG3	1:A:117:MET:HE2	1.84	0.60
1:C:258:SER:HB3	1:C:407:THR:HG22	1.83	0.60
1:A:258:SER:HB3	1:A:407:THR:HG22	1.82	0.60
1:D:195:PRO:HB2	1:D:199:PRO:HD2	1.85	0.59
1:C:195:PRO:HB2	1:C:199:PRO:HD2	1.85	0.58
1:C:219:ILE:HG22	1:C:417:ILE:HG12	1.86	0.57
1:A:264:LEU:HD21	1:A:275:LEU:HD11	1.87	0.57
1:A:101:LEU:HD21	2:E:3:AC1:HCB2	1.86	0.56
1:A:141:ARG:NH1	1:A:369:GLU:OE2	2.32	0.56
1:C:426:TRP:O	1:C:430:ILE:HG13	2.06	0.56
1:C:282:LEU:HB2	1:C:305:PHE:CD2	2.39	0.56
1:A:2:LYS:HG3	1:A:344:GLN:HG2	1.87	0.56
1:D:355:ASN:HD21	2:H:2:GLC:H3	1.70	0.56
1:D:111:GLU:HG2	1:D:114:ARG:HH21	1.69	0.56
1:B:181:ARG:O	1:B:185:GLU:HG2	2.06	0.56
1:B:313:LEU:HA	1:B:317:SER:HB3	1.88	0.55
1:C:36:LYS:HD3	1:C:138:TRP:HA	1.87	0.55
1:C:419:ASN:HB3	1:C:422:LYS:HD3	1.89	0.55
1:D:297:PHE:HB3	1:D:301:ILE:HG22	1.89	0.55
1:B:324:GLU:H	1:B:324:GLU:CD	2.15	0.55
1:B:111:GLU:HA	1:B:114:ARG:HH11	1.72	0.54
1:D:113:ILE:HD11	1:D:133:LEU:HB3	1.90	0.54
1:B:286:VAL:HB	1:B:289:ALA:HB2	1.89	0.54
1:D:120:GLU:CD	1:D:120:GLU:H	2.16	0.53
1:C:429:LEU:HD21	1:C:445:GLU:HG2	1.91	0.53
1:B:282:LEU:HB2	1:B:305:PHE:CD1	2.43	0.53
1:D:79:ALA:O	1:D:131:GLN:NE2	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:274:MET:HG3	1:C:390:MET:HE3	1.91	0.53
1:B:258:SER:HB3	1:B:407:THR:HG22	1.90	0.52
1:A:181:ARG:O	1:A:185:GLU:HG2	2.08	0.52
1:C:370:LYS:O	1:C:374:THR:OG1	2.20	0.52
1:B:195:PRO:HB3	1:B:233:HIS:HB3	1.91	0.52
1:C:57:THR:HG22	1:C:194:SER:HB2	1.91	0.52
1:B:456:GLU:N	1:B:456:GLU:OE1	2.44	0.52
1:A:298:ARG:HG2	1:A:401:GLU:HA	1.93	0.51
1:B:260:SER:H	1:B:263:ALA:HB3	1.75	0.51
1:B:335:LYS:HG2	1:B:380:TRP:CZ2	2.45	0.51
1:A:15:PRO:HG3	1:A:448:ASP:CG	2.36	0.51
1:A:293:LEU:HD21	1:A:398:GLU:HB2	1.93	0.51
1:C:112:ARG:HB3	1:C:133:LEU:HD11	1.92	0.51
1:D:57:THR:HG22	1:D:194:SER:HB2	1.93	0.51
1:C:374:THR:O	1:C:378:GLU:HG2	2.10	0.51
1:D:282:LEU:HB2	1:D:305:PHE:CD2	2.46	0.51
1:B:370:LYS:O	1:B:374:THR:OG1	2.24	0.51
1:C:297:PHE:HB3	1:C:301:ILE:HG22	1.93	0.51
1:A:224:VAL:HG11	1:A:433:LYS:HG2	1.93	0.50
1:C:74:ASP:HB3	1:C:77:LEU:HB2	1.92	0.50
1:A:379:GLU:OE1	1:A:379:GLU:N	2.38	0.50
1:C:105:LYS:HD3	1:C:358:GLU:HB3	1.94	0.50
1:D:90:GLU:H	1:D:90:GLU:CD	2.19	0.50
1:A:34:ALA:HA	1:A:38:TYR:HB2	1.94	0.50
1:C:222:PHE:HA	1:C:430:ILE:HD13	1.93	0.50
1:D:265:GLU:O	1:D:269:GLU:HG3	2.11	0.50
1:C:5:PHE:CD1	1:C:384:LEU:HD11	2.47	0.50
1:D:109:ARG:NH1	1:D:361:GLU:OE2	2.45	0.50
1:D:224:VAL:HG21	1:D:433:LYS:HB3	1.93	0.50
1:D:378:GLU:OE1	1:D:380:TRP:NE1	2.44	0.50
1:B:191:VAL:HG11	1:B:245:PHE:CZ	2.47	0.49
1:A:274:MET:HE1	1:A:386:LEU:HD22	1.93	0.49
1:B:195:PRO:HB2	1:B:199:PRO:HD2	1.94	0.49
1:C:49:GLU:N	1:C:376:GLU:OE2	2.32	0.49
1:D:429:LEU:HD21	1:D:445:GLU:HG2	1.94	0.49
1:D:15:PRO:HG3	1:D:448:ASP:CG	2.37	0.49
1:D:84:PRO:HA	1:D:87:LEU:HG	1.95	0.49
1:C:294:ARG:HB3	1:C:397:THR:HG22	1.93	0.49
1:B:465:ASP:O	1:B:469:ARG:HG3	2.11	0.49
1:C:2:LYS:HE2	1:C:338:HIS:CD2	2.46	0.49
1:B:3:LYS:H	1:B:344:GLN:HG2	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:462:GLU:H	1:B:462:GLU:CD	2.21	0.49
1:D:314:ILE:O	1:D:360:TYR:OH	2.18	0.49
1:B:34:ALA:HA	1:B:38:TYR:HB2	1.94	0.49
1:C:121:LYS:HD2	1:C:122:LEU:H	1.77	0.49
1:D:2:LYS:HE2	1:D:338:HIS:CD2	2.48	0.49
1:A:466:LYS:O	1:A:470:SER:OG	2.27	0.48
1:C:304:PHE:HE1	1:C:390:MET:HE2	1.78	0.48
1:D:31:PHE:HB2	1:D:67:TYR:CE1	2.49	0.48
1:A:182:GLU:O	1:A:186:GLU:HG3	2.13	0.48
1:A:20:PRO:HB2	1:A:474:ARG:HD3	1.96	0.48
1:D:418:GLY:O	1:D:423:ASN:ND2	2.47	0.48
1:A:195:PRO:HB2	1:A:199:PRO:HD2	1.95	0.48
1:B:13:HIS:O	2:F:2:GLC:H2	2.14	0.48
1:C:424:GLU:HA	1:C:427:LYS:HD2	1.95	0.48
1:D:380:TRP:H	1:D:380:TRP:CD1	2.32	0.48
1:A:162:LYS:O	1:A:166:LEU:HD12	2.14	0.47
1:C:274:MET:HB2	1:C:390:MET:HE3	1.95	0.47
1:D:323:ALA:O	1:D:327:VAL:HG23	2.14	0.47
1:B:205:LEU:HD21	1:B:266:LEU:HD22	1.95	0.47
1:B:274:MET:HE3	1:B:304:PHE:CD1	2.49	0.47
1:B:336:LYS:HA	1:B:336:LYS:HD3	1.68	0.47
1:B:368:LEU:HD22	1:B:372:TYR:CZ	2.50	0.47
1:B:469:ARG:O	1:B:473:ARG:HG2	2.15	0.47
1:C:15:PRO:HG3	1:C:448:ASP:CG	2.40	0.47
1:D:181:ARG:HD2	1:D:244:ILE:O	2.15	0.47
1:A:82:LYS:HG2	1:A:86:ASP:HB2	1.96	0.47
1:B:438:ASN:ND2	1:B:440:TYR:HB2	2.31	0.47
2:H:2:GLC:H5	2:H:3:AC1:O5	2.14	0.47
1:D:335:LYS:HA	1:D:335:LYS:HD2	1.71	0.46
1:C:122:LEU:HD12	1:C:126:GLU:HB3	1.97	0.46
1:A:3:LYS:N	1:A:344:GLN:OE1	2.47	0.46
1:B:115:GLU:O	1:B:119:LYS:HG3	2.15	0.46
1:C:82:LYS:HE3	1:C:87:LEU:HD23	1.96	0.46
1:A:294:ARG:O	1:A:397:THR:HA	2.16	0.46
1:B:202:PRO:HB3	1:B:260:SER:HB3	1.97	0.46
1:D:67:TYR:OH	1:D:74:ASP:OD2	2.32	0.46
1:B:265:GLU:O	1:B:269:GLU:HG3	2.16	0.46
1:B:57:THR:HG22	1:B:194:SER:HB2	1.98	0.46
1:C:46:LYS:HD3	1:C:179:ILE:HD11	1.98	0.46
1:A:36:LYS:HG2	1:A:363:ASN:HD21	1.80	0.45
1:C:409:PHE:CD2	1:C:415:LYS:HB2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:LEU:HD13	1:B:305:PHE:CE2	2.51	0.45
1:B:195:PRO:HD2	1:B:255:PRO:HD3	1.97	0.45
1:C:4:LEU:HD21	1:C:334:LEU:HD22	1.98	0.45
1:D:292:TYR:HA	1:D:346:VAL:O	2.17	0.45
1:C:377:LYS:HA	1:C:377:LYS:HD3	1.78	0.45
1:B:240:LYS:O	1:B:244:ILE:HG12	2.17	0.45
1:B:15:PRO:HG3	1:B:448:ASP:CG	2.41	0.45
1:D:285:SER:HB3	1:D:403:VAL:HA	1.99	0.45
1:D:287:GLU:OE2	1:D:288:ARG:NE	2.50	0.45
1:A:282:LEU:HB2	1:A:305:PHE:CD2	2.52	0.45
1:C:343:PHE:H	1:C:343:PHE:HD1	1.65	0.45
1:A:432:ALA:HB2	1:A:472:VAL:HG13	1.99	0.45
1:B:176:THR:O	1:B:179:ILE:HG22	2.17	0.45
1:B:392:LYS:HE2	1:B:392:LYS:HB2	1.64	0.45
1:D:258:SER:HA	1:D:406:GLY:O	2.17	0.45
1:A:9:TRP:HE1	1:A:55:ASN:HB2	1.82	0.44
1:C:217:VAL:HG11	1:C:414:LEU:HD21	1.99	0.44
1:D:424:GLU:HA	1:D:427:LYS:HD2	1.99	0.44
1:A:183:ILE:HG13	1:A:184:LYS:N	2.32	0.44
1:A:199:PRO:HG3	1:A:233:HIS:CG	2.51	0.44
1:C:195:PRO:HD2	1:C:255:PRO:HD3	1.99	0.44
1:D:3:LYS:H	1:D:344:GLN:HE21	1.64	0.44
1:C:19:GLU:OE1	1:C:21:TYR:N	2.50	0.44
1:C:254:PRO:HG2	1:C:257:ALA:HA	1.98	0.44
1:D:166:LEU:HD23	1:D:166:LEU:HA	1.88	0.44
1:C:36:LYS:NZ	1:C:363:ASN:OD1	2.51	0.44
1:C:432:ALA:HB2	1:C:472:VAL:HG13	1.98	0.44
1:C:200:LEU:HD23	1:C:200:LEU:HA	1.82	0.44
1:D:181:ARG:HH21	1:D:185:GLU:CD	2.26	0.44
1:D:264:LEU:HD23	1:D:275:LEU:HD21	2.00	0.44
1:B:112:ARG:HB3	1:B:133:LEU:HD11	1.99	0.43
1:D:176:THR:O	1:D:179:ILE:HG22	2.18	0.43
1:D:217:VAL:HG11	1:D:414:LEU:HD21	1.99	0.43
1:A:69:GLN:HB2	1:A:71:LYS:HG3	2.00	0.43
1:A:224:VAL:CG2	1:A:434:LYS:HG3	2.48	0.43
1:D:224:VAL:HG22	1:D:430:ILE:HG23	1.99	0.43
1:A:344:GLN:NE2	1:A:394:ASP:HB3	2.33	0.43
1:D:181:ARG:O	1:D:185:GLU:HG2	2.18	0.43
1:A:9:TRP:HB3	1:A:350:VAL:HG13	2.00	0.43
1:C:2:LYS:NZ	1:C:341:VAL:O	2.50	0.43
1:C:191:VAL:HG11	1:C:245:PHE:CZ	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:ARG:HG2	1:D:18:ARG:HH11	1.83	0.43
1:A:264:LEU:HD23	1:A:275:LEU:HD21	1.99	0.43
1:D:396:LYS:HD2	1:D:396:LYS:HA	1.70	0.43
1:B:370:LYS:HA	1:B:370:LYS:HD3	1.76	0.43
1:A:36:LYS:HG2	1:A:363:ASN:ND2	2.34	0.43
1:A:392:LYS:HD2	1:A:394:ASP:HB2	1.99	0.43
1:C:396:LYS:HD3	1:C:396:LYS:HA	1.90	0.43
1:C:424:GLU:O	1:C:428:ILE:HG13	2.18	0.43
1:B:441:ILE:O	1:B:445:GLU:HG3	2.18	0.43
1:C:53:ASN:HD21	1:C:385:THR:HA	1.84	0.43
1:B:30:THR:HG21	1:B:63:GLN:HB3	2.00	0.42
1:C:92:VAL:O	1:C:96:ILE:HG12	2.19	0.42
1:C:143:LEU:HD12	1:C:143:LEU:HA	1.83	0.42
1:C:414:LEU:HA	1:C:414:LEU:HD23	1.82	0.42
1:D:292:TYR:CZ	1:D:337:ILE:HD13	2.54	0.42
1:C:234:VAL:O	1:C:238:LYS:HG3	2.19	0.42
1:D:48:PHE:HE2	1:D:373:SER:HB3	1.84	0.42
1:A:293:LEU:CD2	1:A:398:GLU:HB2	2.50	0.42
1:B:285:SER:HB3	1:B:403:VAL:HA	2.02	0.42
1:C:286:VAL:HB	1:C:289:ALA:HB2	2.02	0.42
1:D:281:ILE:HD11	1:D:406:GLY:O	2.19	0.42
1:C:88:GLU:O	1:C:92:VAL:HG23	2.19	0.42
1:D:20:PRO:HB3	1:D:440:TYR:CE1	2.55	0.42
1:D:119:LYS:NZ	1:D:126:GLU:OE2	2.48	0.42
1:A:327:VAL:HG22	1:A:371:LEU:HB2	2.02	0.42
1:B:36:LYS:HB3	1:B:36:LYS:HE3	1.81	0.42
1:A:30:THR:HG21	1:A:63:GLN:HB3	2.02	0.42
1:A:198:HIS:CE1	1:A:255:PRO:HB3	2.55	0.42
1:D:199:PRO:HG3	1:D:233:HIS:CG	2.54	0.42
1:A:103:TYR:OH	1:A:114:ARG:HG2	2.20	0.42
1:B:20:PRO:HD3	1:B:471:PHE:CE1	2.55	0.42
1:A:380:TRP:CD1	1:A:380:TRP:H	2.37	0.42
1:B:68:ALA:HB2	1:B:173:ILE:HG22	2.01	0.42
1:B:224:VAL:HG21	1:B:433:LYS:HB3	2.01	0.42
1:B:375:LEU:HD22	1:B:381:ILE:HG21	2.02	0.42
1:C:378:GLU:HB3	1:C:380:TRP:NE1	2.34	0.42
1:A:202:PRO:HB3	1:A:260:SER:HB3	2.03	0.41
1:C:333:ARG:O	1:C:337:ILE:HG13	2.19	0.41
1:C:10:TRP:HE1	1:C:52:LEU:HD22	1.85	0.41
1:D:82:LYS:HD2	1:D:87:LEU:HD23	2.02	0.41
1:D:335:LYS:O	1:D:339:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:HE3	1:B:3:LYS:NZ	2.35	0.41
1:B:106:PRO:HD2	1:B:359:TYR:HA	2.02	0.41
1:B:258:SER:HA	1:B:406:GLY:O	2.21	0.41
1:B:214:THR:HB	1:B:217:VAL:HG13	2.02	0.41
1:B:56:LEU:HD21	1:B:180:TYR:CZ	2.55	0.41
1:D:106:PRO:HD2	1:D:359:TYR:HA	2.02	0.41
1:A:12:MET:HB2	1:A:56:LEU:HD23	2.03	0.41
1:C:416:TRP:NE1	1:C:451:TRP:HD1	2.19	0.41
1:B:274:MET:HE2	1:B:275:LEU:C	2.46	0.41
1:B:294:ARG:NH2	1:B:392:LYS:O	2.46	0.41
1:C:193:THR:HG23	1:C:252:MET:HE3	2.03	0.41
1:C:235:GLU:HG2	1:C:270:LYS:HD2	2.03	0.41
1:C:332:GLY:O	1:C:336:LYS:HG2	2.21	0.41
1:C:368:LEU:HD23	1:C:368:LEU:HA	1.91	0.41
1:D:200:LEU:HD23	1:D:200:LEU:HA	1.79	0.41
1:D:274:MET:HA	1:D:302:SER:O	2.21	0.41
1:B:256:GLU:OE1	2:F:1:GLC:H1	2.21	0.41
1:C:36:LYS:HE2	1:C:138:TRP:CE3	2.56	0.41
1:D:14:GLN:HE22	1:D:29:TRP:HB3	1.86	0.41
1:D:25:TYR:OH	1:D:66:GLU:OE1	2.30	0.41
1:B:197:TYR:HB2	1:B:233:HIS:HE1	1.85	0.40
1:C:214:THR:HB	1:C:217:VAL:HG13	2.03	0.40
1:D:333:ARG:O	1:D:337:ILE:HG13	2.20	0.40
1:B:150:LEU:HD23	1:B:150:LEU:HA	1.82	0.40
1:D:168:LYS:HD2	1:D:168:LYS:HA	1.73	0.40
1:D:289:ALA:HB1	1:D:295:TYR:OH	2.21	0.40
1:A:172:ILE:HD13	1:A:172:ILE:HA	1.90	0.40
1:A:384:LEU:HD11	1:A:392:LYS:HE3	2.04	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	475/484 (98%)	467 (98%)	7 (2%)	1 (0%)	44	71
1	B	475/484 (98%)	468 (98%)	7 (2%)	0	100	100
1	C	482/484 (100%)	474 (98%)	8 (2%)	0	100	100
1	D	482/484 (100%)	470 (98%)	12 (2%)	0	100	100
All	All	1914/1936 (99%)	1879 (98%)	34 (2%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	455	GLU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	437/444 (98%)	437 (100%)	0	100	100
1	B	437/444 (98%)	437 (100%)	0	100	100
1	C	444/444 (100%)	444 (100%)	0	100	100
1	D	444/444 (100%)	444 (100%)	0	100	100
All	All	1762/1776 (99%)	1762 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	A	208	ASN
1	A	355	ASN
1	B	14	GLN
1	B	216	ASN
1	D	208	ASN
1	D	344	GLN

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Mol	Chain	Res	Type
1	D	453	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 ⓘ

12 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GLC	E	1	2	12,12,12	0.41	0	17,17,17	0.76	0
2	GLC	E	2	2	11,11,12	0.61	0	15,15,17	1.02	0
2	AC1	E	3	2	21,22,23	0.29	0	22,32,34	0.71	0
2	GLC	F	1	2	12,12,12	0.46	0	17,17,17	1.13	2 (11%)
2	GLC	F	2	2	11,11,12	0.62	0	15,15,17	0.93	1 (6%)
2	AC1	F	3	2	21,22,23	0.25	0	22,32,34	0.52	0
2	GLC	G	1	2	12,12,12	0.42	0	17,17,17	1.15	2 (11%)
2	GLC	G	2	2	11,11,12	0.81	0	15,15,17	1.53	3 (20%)
2	AC1	G	3	2	21,22,23	0.28	0	22,32,34	0.81	0
2	GLC	H	1	2	12,12,12	0.40	0	17,17,17	1.08	2 (11%)
2	GLC	H	2	2	11,11,12	0.62	0	15,15,17	1.34	2 (13%)
2	AC1	H	3	2	21,22,23	0.26	0	22,32,34	0.52	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	E	1	2	-	2/2/22/22	0/1/1/1
2	GLC	E	2	2	-	0/2/19/22	0/1/1/1
2	AC1	E	3	2	-	2/6/43/46	0/2/2/2
2	GLC	F	1	2	-	2/2/22/22	0/1/1/1
2	GLC	F	2	2	-	1/2/19/22	0/1/1/1
2	AC1	F	3	2	-	3/6/43/46	0/2/2/2
2	GLC	G	1	2	-	2/2/22/22	0/1/1/1
2	GLC	G	2	2	-	0/2/19/22	0/1/1/1
2	AC1	G	3	2	-	1/6/43/46	0/2/2/2
2	GLC	H	1	2	-	2/2/22/22	0/1/1/1
2	GLC	H	2	2	-	0/2/19/22	0/1/1/1
2	AC1	H	3	2	-	3/6/43/46	0/2/2/2

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	2	GLC	C1-O5-C5	4.19	117.87	112.19
2	G	2	GLC	C1-O5-C5	3.82	117.37	112.19
2	G	2	GLC	O4-C4-C3	-3.51	102.22	110.35
2	G	1	GLC	O4-C4-C3	-3.02	103.36	110.35
2	H	2	GLC	O5-C5-C6	2.63	111.33	107.20
2	F	2	GLC	C1-C2-C3	2.61	112.87	109.67
2	H	1	GLC	O4-C4-C3	-2.37	104.86	110.35
2	F	1	GLC	C3-C4-C5	2.30	114.35	110.24
2	H	1	GLC	C3-C4-C5	2.22	114.21	110.24
2	F	1	GLC	O4-C4-C5	-2.12	104.05	109.30
2	G	2	GLC	O2-C2-C3	-2.09	105.95	110.14
2	G	1	GLC	C3-C4-C5	2.08	113.95	110.24

There are no chirality outliers.

All (18) torsion outliers are listed below:

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

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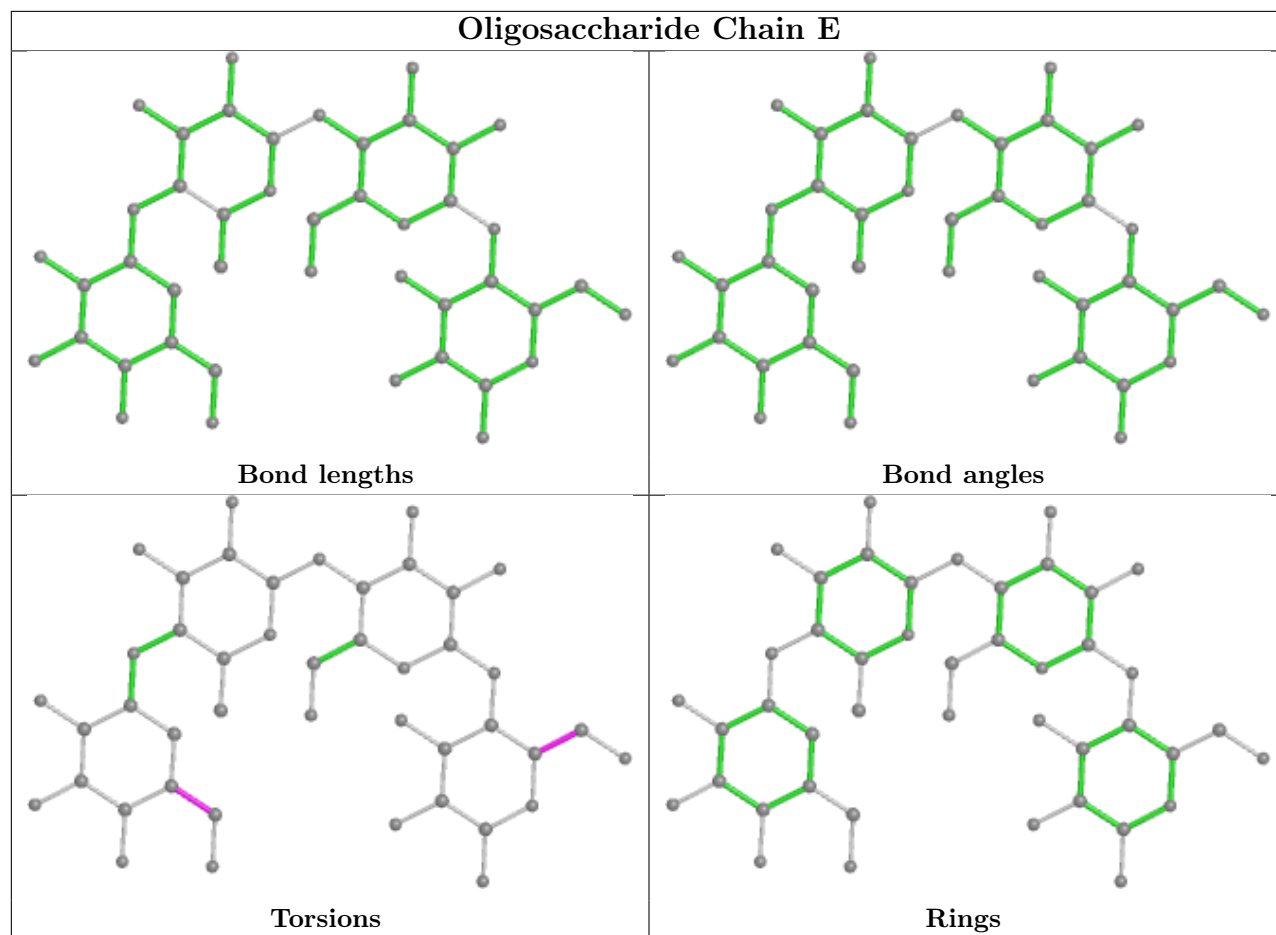
Mol	Chain	Res	Type	Atoms
2	G	1	GLC	O5-C5-C6-O6
2	F	1	GLC	C4-C5-C6-O6
2	F	2	GLC	O5-C5-C6-O6
2	F	1	GLC	O5-C5-C6-O6
2	H	3	AC1	C2B-C1B-N4A-C4
2	F	3	AC1	C5-C4-N4A-C1B
2	E	3	AC1	C4A-C5B-C6B-O6B
2	H	3	AC1	C4A-C5B-C6B-O6B
2	F	3	AC1	C7B-C1B-N4A-C4
2	E	3	AC1	C7B-C5B-C6B-O6B
2	F	3	AC1	C7B-C5B-C6B-O6B
2	G	3	AC1	C7B-C5B-C6B-O6B
2	H	3	AC1	C7B-C5B-C6B-O6B

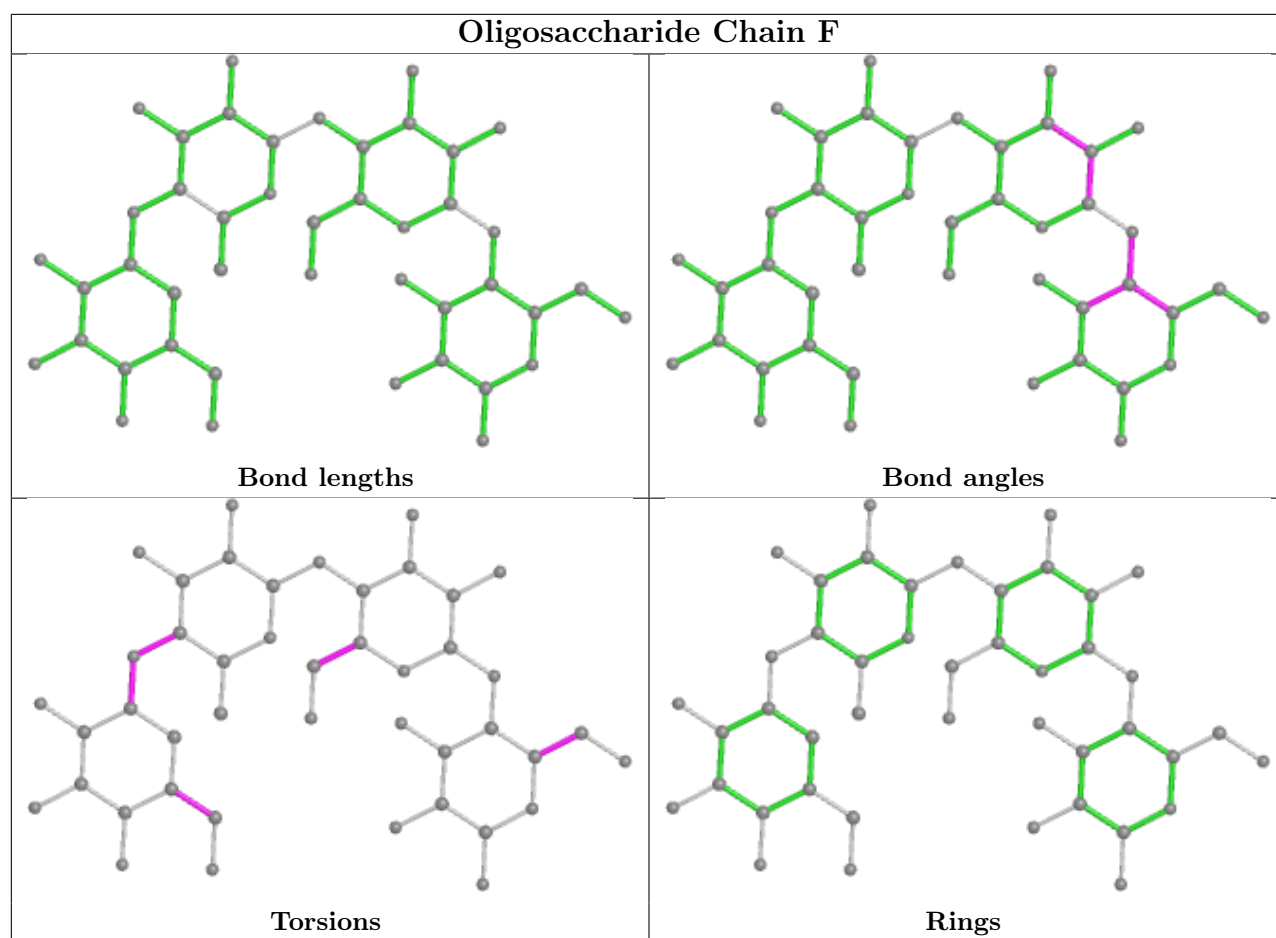
There are no ring outliers.

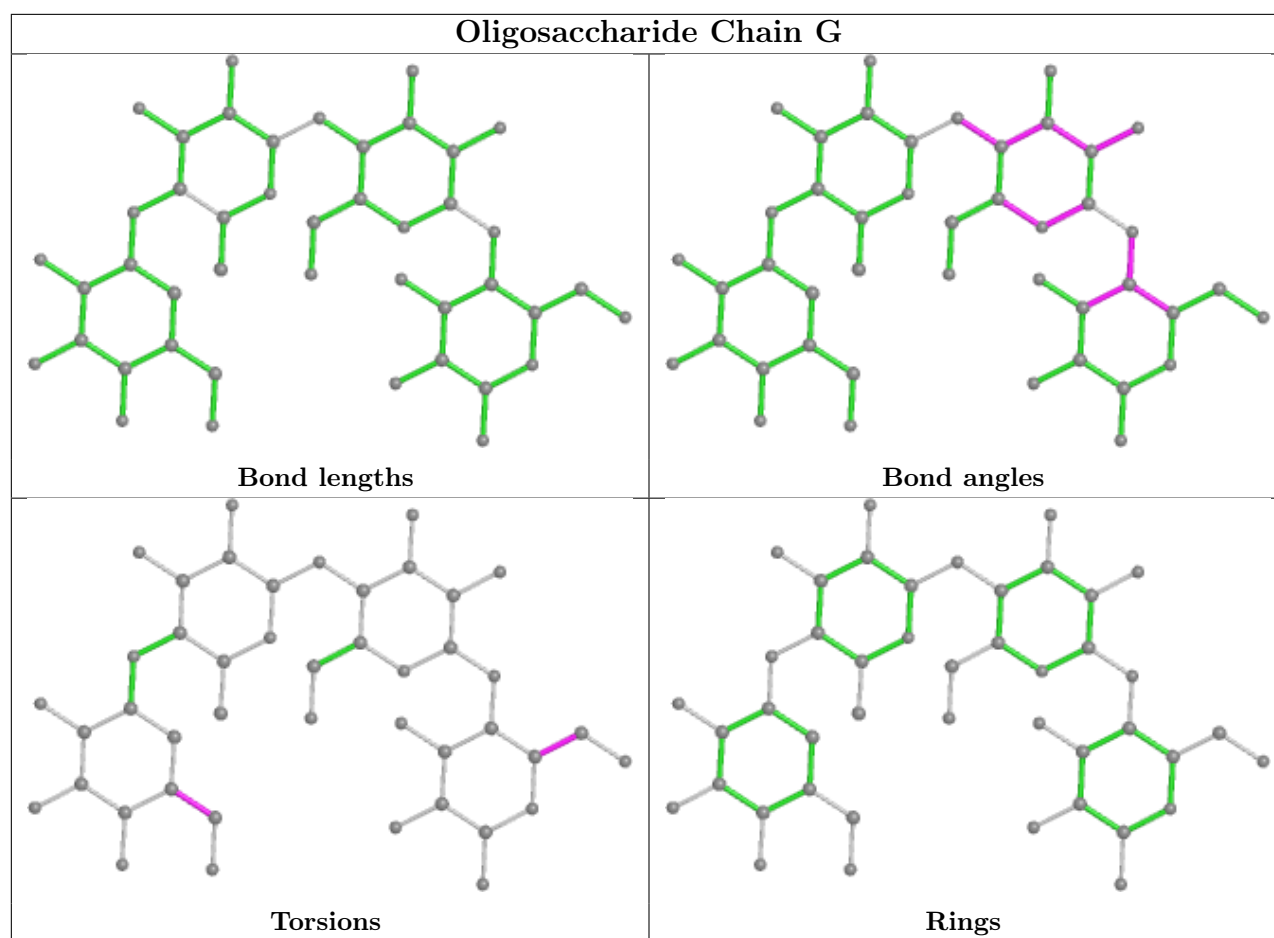
5 monomers are involved in 6 short contacts:

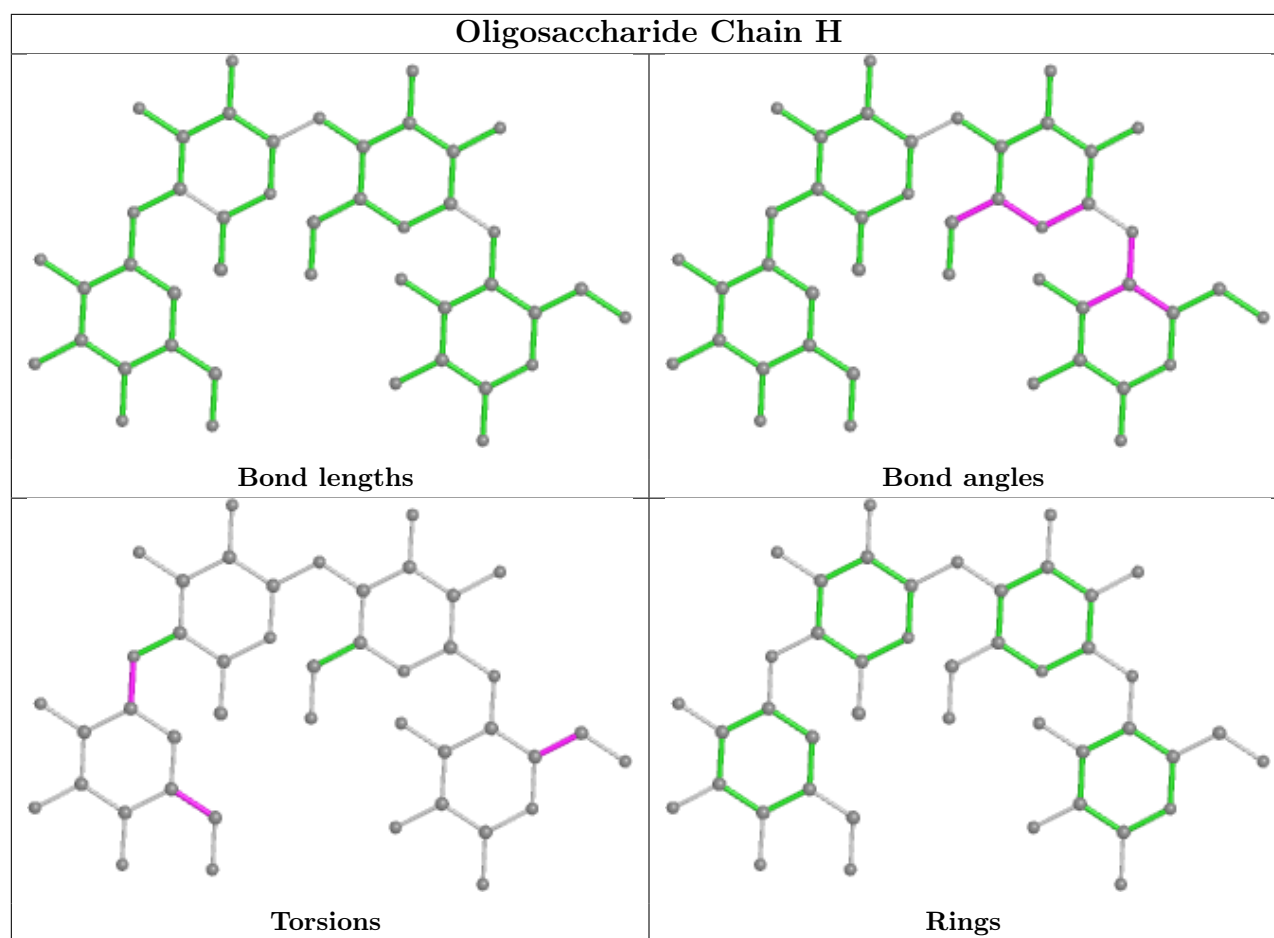
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	2	GLC	2	0
2	F	2	GLC	1	0
2	F	1	GLC	1	0
2	E	3	AC1	1	0
2	H	3	AC1	2	0

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









5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	477/484 (98%)	-0.49	0 100 100	35, 49, 75, 121	0
1	B	477/484 (98%)	-0.47	0 100 100	35, 52, 77, 134	0
1	C	484/484 (100%)	-0.29	1 (0%) 92 85	40, 63, 100, 156	0
1	D	484/484 (100%)	-0.31	1 (0%) 92 85	42, 63, 98, 166	0
All	All	1922/1936 (99%)	-0.39	2 (0%) 92 87	35, 57, 92, 166	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	459	PRO	2.7
1	D	352	ASP	2.1

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

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

6.3 Carbohydrates [i](#)

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

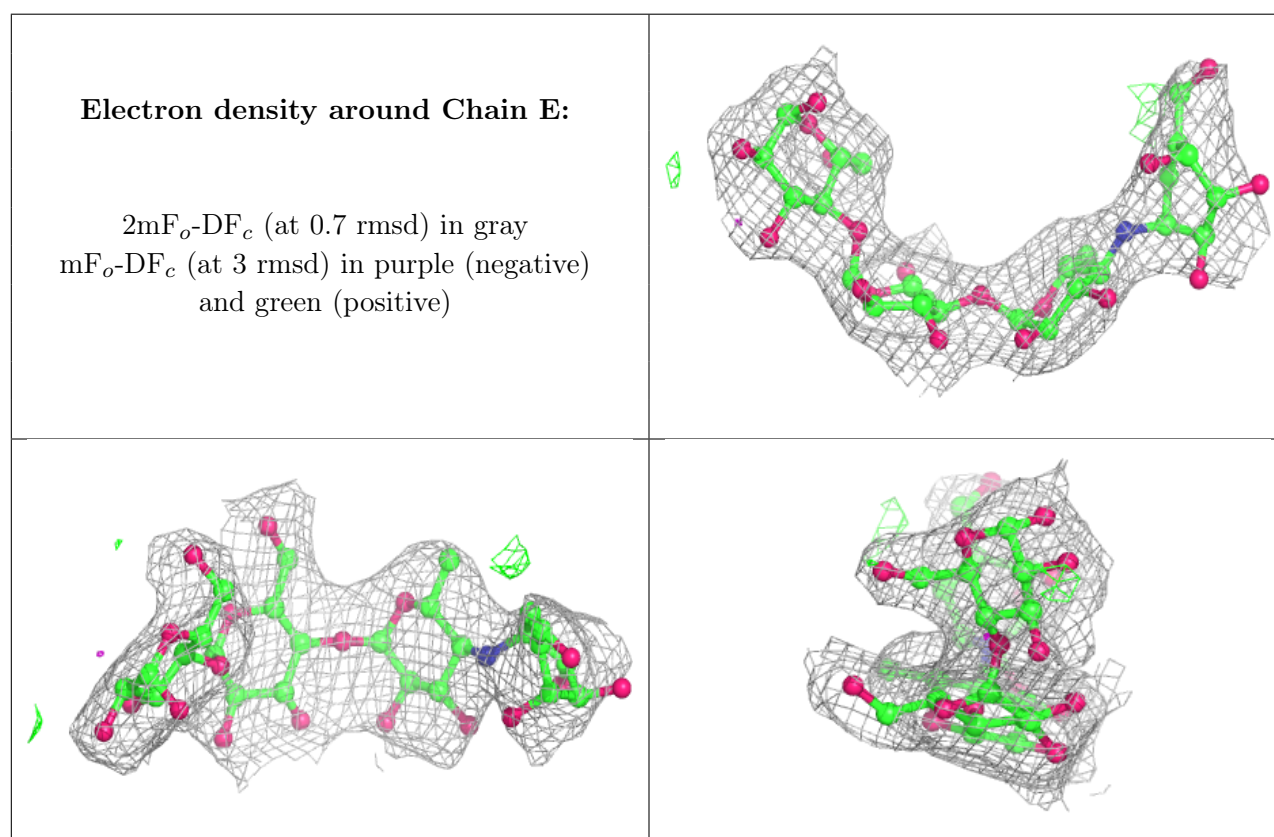
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	H	1	12/12	0.82	0.17	51,60,66,66	0
2	GLC	H	2	11/12	0.87	0.12	49,56,69,83	0
2	AC1	G	3	21/22	0.88	0.11	50,70,84,94	0
2	AC1	E	3	21/22	0.88	0.12	40,54,69,78	0
2	GLC	G	1	12/12	0.88	0.11	49,51,57,60	0

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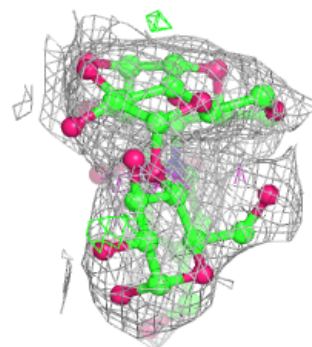
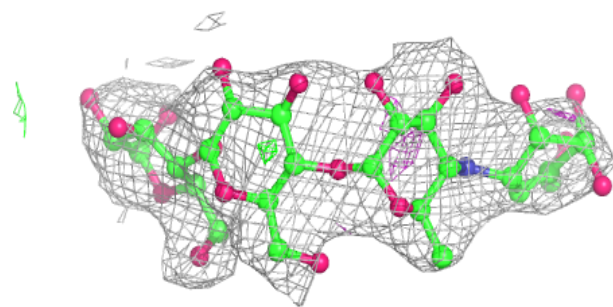
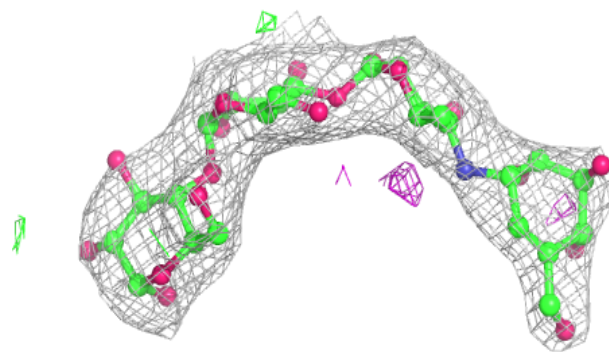
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GLC	G	2	11/12	0.90	0.10	45,49,56,60	0
2	AC1	F	3	21/22	0.91	0.10	44,52,76,89	0
2	GLC	F	2	11/12	0.91	0.08	41,44,57,57	0
2	GLC	F	1	12/12	0.92	0.09	32,49,60,61	0
2	AC1	H	3	21/22	0.92	0.10	49,62,87,99	0
2	GLC	E	1	12/12	0.94	0.07	34,42,56,61	0
2	GLC	E	2	11/12	0.96	0.06	33,38,43,46	0

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

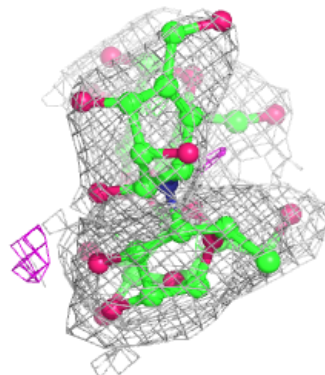
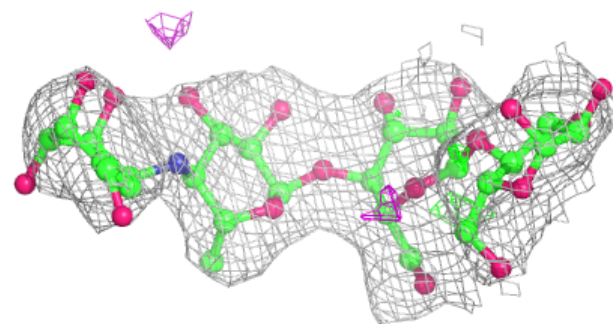
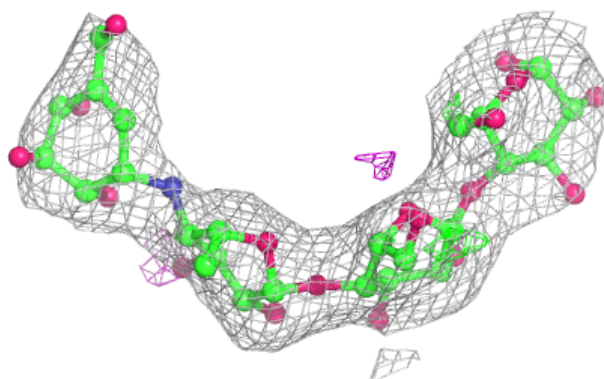


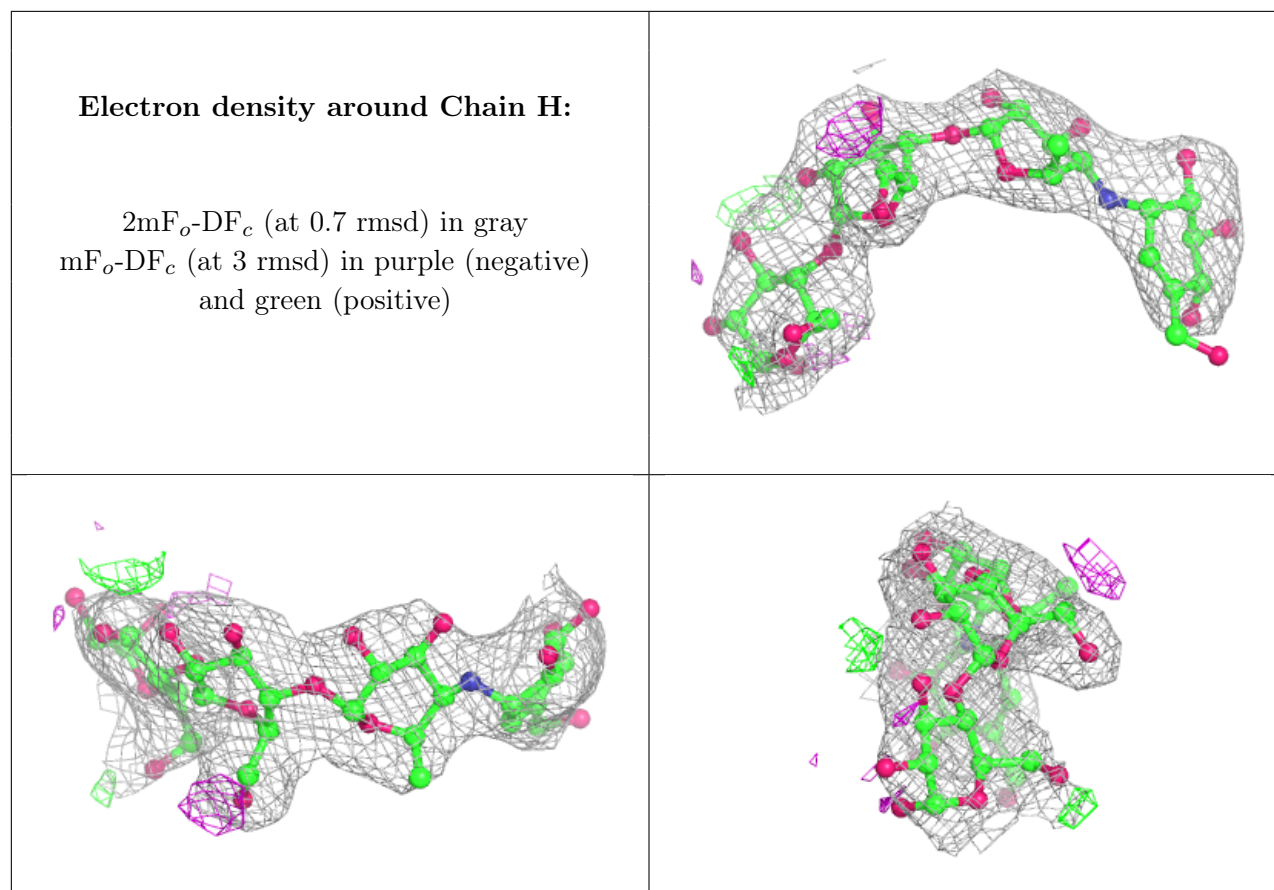
Electron density around Chain F:

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

**Electron density around Chain G:**

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





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.