



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:23 PM UTC

PDB ID : 4PQK / pdb_00004pqk
Title : C-Terminal domain of DNA binding protein
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Deposited on : 2014-03-03
Resolution : 3.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

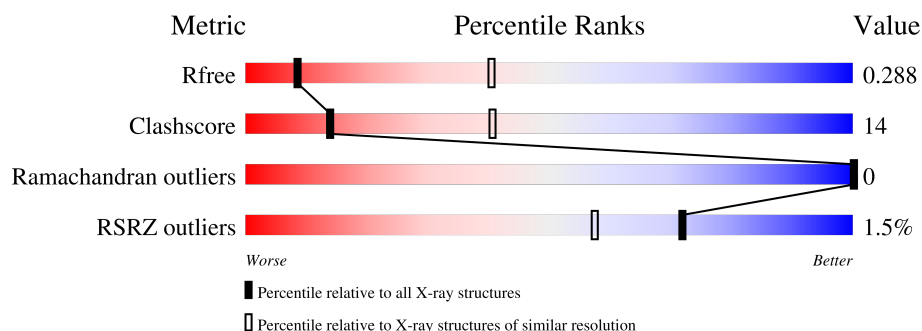
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	487	% 69% 29% .
1	B	487	2% 76% 22% .
1	C	487	% 73% 25% .
1	D	487	% 68% 29% .
2	E	3	33% 67%
2	F	3	100%

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Mol	Chain	Length	Quality of chain
2	G	3	 67%33%
2	H	3	 100%

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	GLC	E	3	-	-	X	-

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 14876 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.

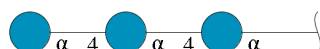
- Molecule 1 is a protein called Maltose ABC transporter periplasmic protein, Truncated replication protein RepA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	478	Total	C	N	O	S	0	0	0
			3700	2371	603	713	13			
1	B	479	Total	C	N	O	S	0	0	0
			3722	2392	615	703	12			
1	C	476	Total	C	N	O	S	0	0	0
			3648	2346	601	690	11			
1	D	472	Total	C	N	O	S	0	0	0
			3670	2351	601	707	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP U6NJU2
B	1	MET	-	expression tag	UNP U6NJU2
C	1	MET	-	expression tag	UNP U6NJU2
D	1	MET	-	expression tag	UNP U6NJU2

- Molecule 2 is an oligosaccharide called 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	O	0	0	0
			34	18	16			
2	F	3	Total	C	O	0	0	0
			34	18	16			
2	G	3	Total	C	O	0	0	0
			34	18	16			

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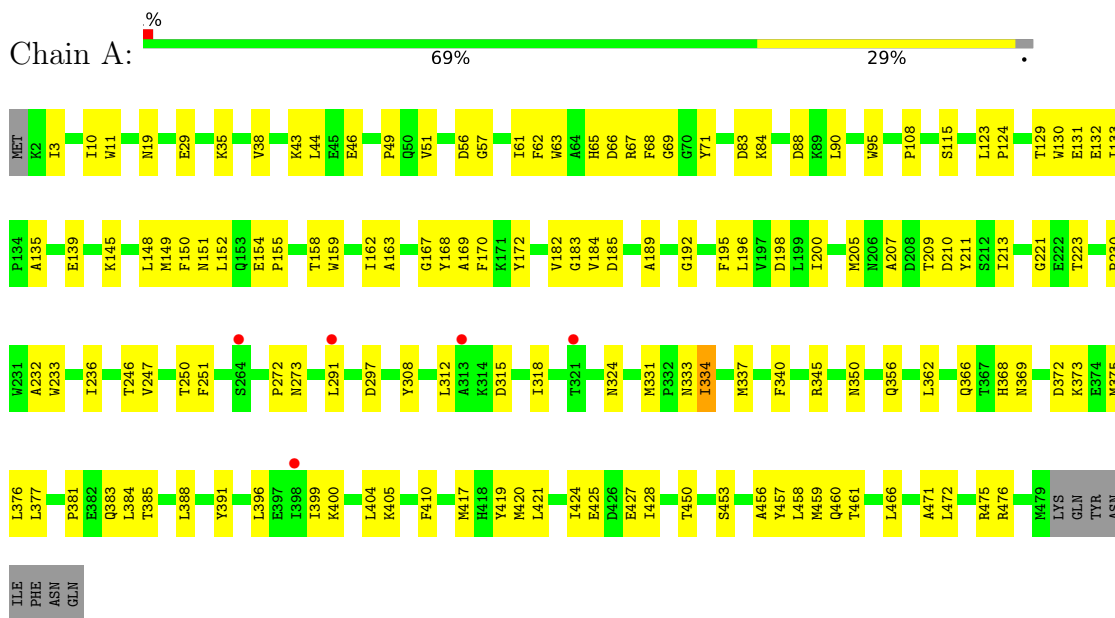
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	H	3	Total	C	O	0	0	0
			34	18	16			

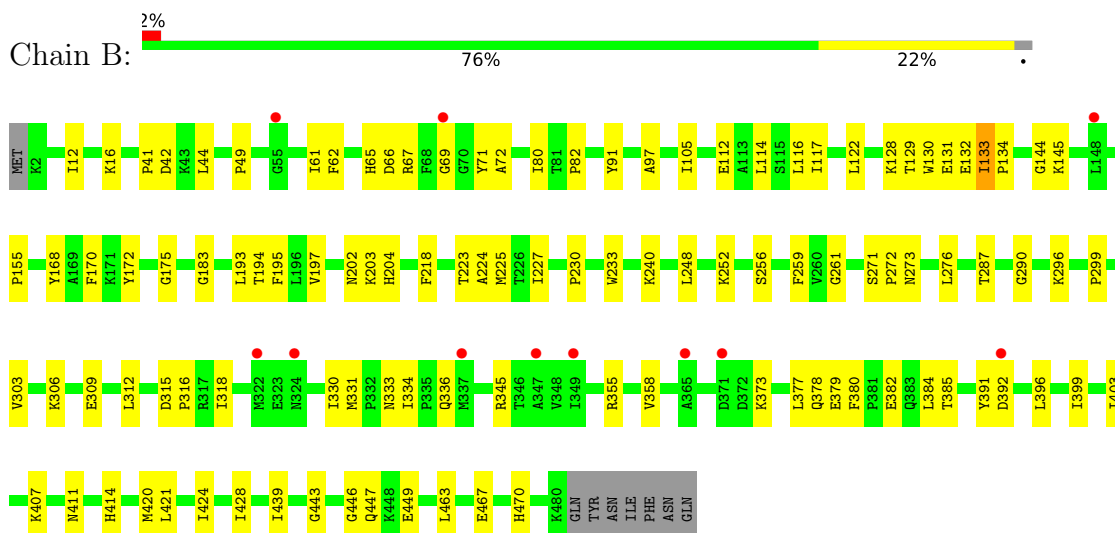
3 Residue-property plots

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.

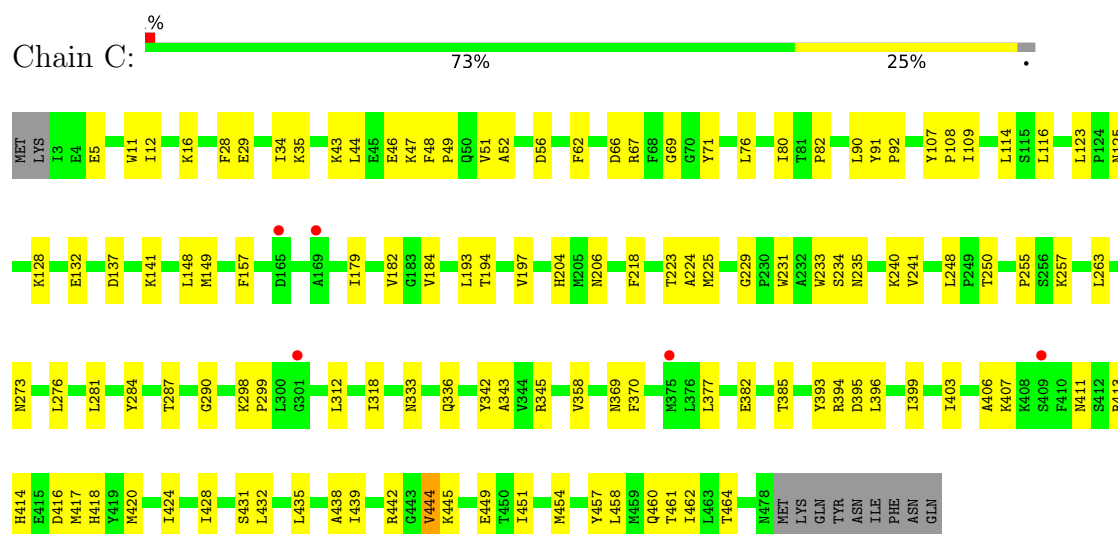
- Molecule 1: Maltose ABC transporter periplasmic protein, Truncated replication protein RepA



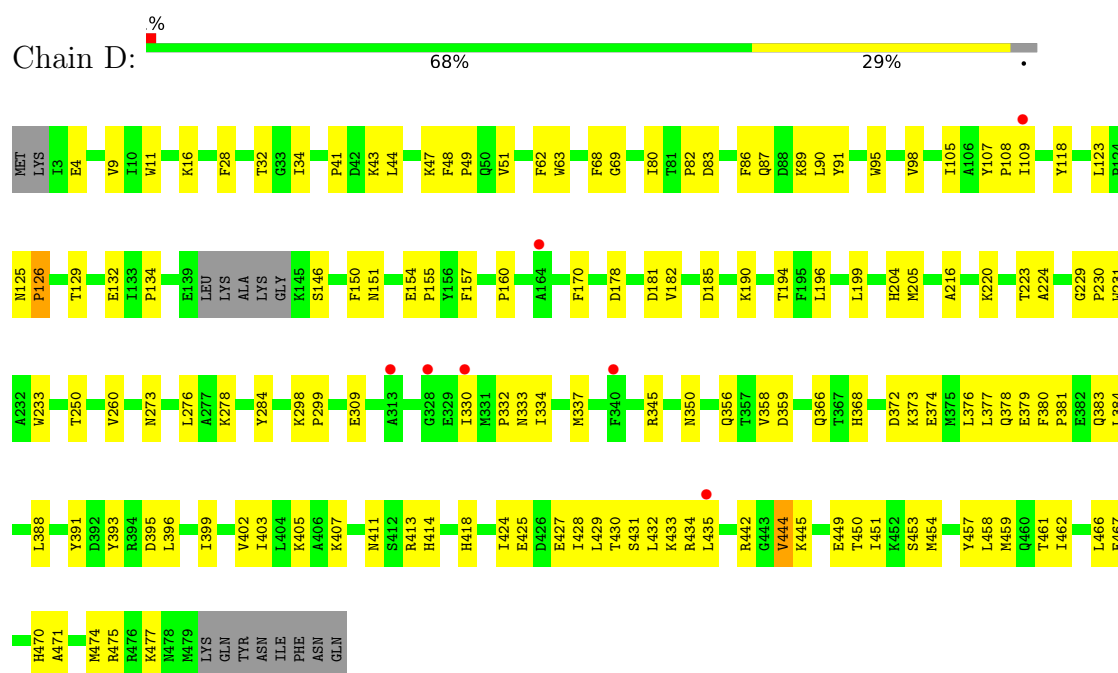
- Molecule 1: Maltose ABC transporter periplasmic protein, Truncated replication protein RepA



- Molecule 1: Maltose ABC transporter periplasmic protein, Truncated replication protein RepA



- Molecule 1: Maltose ABC transporter periplasmic protein, Truncated replication protein RepA



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



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



GLC1
GLC2
GLC3

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

Chain G:  67% 33%

GLC1
GLC2
GLC3

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

Chain H:  100%

GLC1
GLC2
GLC3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	64.22Å 93.67Å 100.81Å 107.67° 108.30° 84.93°	Depositor
Resolution (Å)	46.62 – 3.40 46.62 – 3.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (46.62-3.40) 67.2 (46.62-3.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 3.40Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.237 , 0.275 (Not available) , 0.288	Depositor DCC
R_{free} test set	1963 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	73.9	Xtriage
Anisotropy	0.378	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 31.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	14876	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.50% 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

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	0.58	0/3784	0.92	2/5142 (0.0%)
1	B	0.59	0/3805	0.89	2/5157 (0.0%)
1	C	0.60	0/3730	1.01	8/5063 (0.2%)
1	D	0.60	0/3751	1.00	8/5090 (0.2%)
All	All	0.59	0/15070	0.96	20/20452 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	125	ASN	CA-C-N	10.47	127.20	119.66
1	C	125	ASN	C-N-CA	10.47	127.20	119.66
1	D	444	VAL	N-CA-C	-8.75	100.32	110.21
1	D	418	HIS	N-CA-C	-8.22	102.81	113.17
1	D	125	ASN	CA-C-N	7.94	125.38	119.66
1	D	125	ASN	C-N-CA	7.94	125.38	119.66
1	D	374	GLU	N-CA-C	-6.79	105.29	113.97
1	C	444	VAL	N-CA-C	-6.53	101.11	109.80
1	C	132	GLU	CA-C-N	6.09	124.07	120.24
1	C	132	GLU	C-N-CA	6.09	124.07	120.24
1	B	133	ILE	CB-CA-C	-6.07	107.91	114.35
1	C	370	PHE	N-CA-C	5.68	117.95	109.25
1	D	126	PRO	CA-C-N	5.62	125.55	119.76
1	D	126	PRO	C-N-CA	5.62	125.55	119.76
1	D	9	VAL	N-CA-C	5.47	115.83	108.17
1	B	446	GLY	N-CA-C	5.41	123.50	113.76
1	A	334	ILE	CA-C-N	5.31	125.12	119.28
1	A	334	ILE	C-N-CA	5.31	125.12	119.28
1	C	416	ASP	N-CA-C	-5.28	103.51	110.43
1	C	241	VAL	N-CA-C	5.14	116.30	109.37

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	3700	0	3562	118	0
1	B	3722	0	3663	90	0
1	C	3648	0	3540	88	0
1	D	3670	0	3560	111	0
2	E	34	0	29	8	0
2	F	34	0	30	7	0
2	G	34	0	29	5	0
2	H	34	0	30	6	0
All	All	14876	0	14443	422	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 (422) 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)
2:G:3:GLC:O5	2:G:3:GLC:C5	1.63	1.46
2:E:3:GLC:C5	2:E:3:GLC:O5	1.64	1.45
2:H:3:GLC:C5	2:H:3:GLC:O5	1.64	1.43
2:F:3:GLC:O5	2:F:3:GLC:C5	1.64	1.41
2:F:3:GLC:O5	2:F:3:GLC:C1	1.72	1.38
2:E:3:GLC:O5	2:E:3:GLC:C1	1.73	1.37
2:G:3:GLC:O5	2:G:3:GLC:C1	1.71	1.36
2:H:3:GLC:O5	2:H:3:GLC:C1	1.71	1.36
1:A:155:PRO:HD3	1:A:345:ARG:HG3	1.36	1.05
1:C:413:ARG:HG2	1:C:414:HIS:CE1	1.93	1.03
1:A:129:THR:HG22	1:A:250:THR:OG1	1.61	1.00
1:A:420:MET:HA	1:A:420:MET:HE3	1.43	0.99
1:A:272:PRO:HG2	1:A:273:ASN:ND2	1.79	0.97
1:B:144:GLY:O	1:B:145:LYS:NZ	2.03	0.91
1:B:377:LEU:HD21	1:B:396:LEU:HD21	1.52	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:403:ILE:HD12	1:C:428:ILE:HD13	1.52	0.89
2:E:3:GLC:C5	2:E:3:GLC:C1	2.51	0.88
1:A:420:MET:HA	1:A:420:MET:CE	2.06	0.85
2:G:3:GLC:C5	2:G:3:GLC:C1	2.54	0.85
2:F:3:GLC:C5	2:F:3:GLC:C1	2.56	0.84
1:A:129:THR:HG22	1:A:250:THR:HG1	1.43	0.84
1:A:155:PRO:CD	1:A:345:ARG:HG3	2.07	0.83
1:B:424:ILE:HG22	1:B:428:ILE:HG13	1.60	0.83
1:A:421:LEU:HD12	1:A:421:LEU:H	1.42	0.83
1:D:80:ILE:HG22	1:D:82:PRO:HD3	1.62	0.80
1:C:16:LYS:NZ	1:C:263:LEU:HD22	1.96	0.80
1:B:420:MET:HE3	1:B:421:LEU:H	1.47	0.80
1:A:376:LEU:HD11	1:A:400:LYS:HD2	1.63	0.79
2:H:3:GLC:C5	2:H:3:GLC:C1	2.60	0.78
1:B:170:PHE:CD2	1:B:334:ILE:HD11	2.18	0.78
1:C:11:TRP:HB3	1:C:44:LEU:HD12	1.64	0.78
1:A:420:MET:HE3	1:A:420:MET:CA	2.15	0.77
1:A:155:PRO:HD3	1:A:345:ARG:CG	2.16	0.75
1:D:405:LYS:HB3	1:D:459:MET:HE1	1.66	0.75
1:D:462:ILE:O	1:D:466:LEU:HD12	1.86	0.75
1:D:91:TYR:HE2	1:D:309:GLU:HG3	1.52	0.75
1:B:407:LYS:O	1:B:411:ASN:ND2	2.20	0.74
1:B:296:LYS:HZ2	1:B:296:LYS:HB3	1.51	0.73
1:C:5:GLU:O	1:C:273:ASN:ND2	2.20	0.73
1:D:146:SER:HB2	1:D:223:THR:HG23	1.69	0.73
1:B:296:LYS:HB3	1:B:296:LYS:NZ	2.03	0.73
1:C:342:TYR:HD1	1:C:345:ARG:HH21	1.37	0.72
1:D:430:THR:O	1:D:434:ARG:HG2	1.88	0.72
1:C:47:LYS:O	1:C:51:VAL:HG22	1.88	0.72
1:C:16:LYS:HZ1	1:C:263:LEU:HD22	1.53	0.72
1:C:399:ILE:HD11	1:C:451:ILE:HG12	1.73	0.71
1:B:336:GLN:HE21	1:B:373:LYS:HG3	1.55	0.71
1:A:391:TYR:HB2	1:A:396:LEU:HD12	1.72	0.71
1:D:16:LYS:NZ	2:H:1:GLC:O2	2.24	0.70
1:D:368:HIS:O	1:D:368:HIS:ND1	2.23	0.70
1:D:402:VAL:HG11	1:D:458:LEU:HG	1.72	0.70
1:C:179:ILE:HD12	1:C:179:ILE:H	1.57	0.70
1:D:34:ILE:HD13	1:D:276:LEU:HD21	1.71	0.69
1:C:149:MET:HE3	1:C:206:ASN:HD21	1.58	0.69
1:B:144:GLY:C	1:B:145:LYS:HZ3	2.00	0.69
1:B:382:GLU:O	1:B:385:THR:HG22	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:ILE:HG22	1:B:82:PRO:HD3	1.74	0.68
1:D:388:LEU:HB3	1:D:396:LEU:HD11	1.75	0.68
1:D:414:HIS:HD2	1:D:470:HIS:CD2	2.11	0.68
1:B:240:LYS:HA	1:C:240:LYS:HG3	1.74	0.68
1:A:155:PRO:HG3	1:A:345:ARG:HA	1.77	0.67
1:B:378:GLN:HG2	1:B:379:GLU:HG2	1.75	0.66
1:C:431:SER:OG	1:C:462:ILE:HG22	1.95	0.66
1:A:168:TYR:CE1	1:A:183:GLY:HA3	2.31	0.65
1:B:391:TYR:H	1:B:391:TYR:HD2	1.44	0.65
1:B:303:VAL:HG22	1:B:312:LEU:HD12	1.79	0.65
1:A:67:ARG:NH2	2:E:3:GLC:O3	2.30	0.65
1:B:72:ALA:HB2	1:B:105:ILE:HG21	1.79	0.65
1:D:444:VAL:O	1:D:445:LYS:HB2	1.98	0.64
1:B:377:LEU:HD21	1:B:396:LEU:CD2	2.26	0.63
1:D:475:ARG:HG2	1:D:475:ARG:HH11	1.63	0.63
1:A:51:VAL:HB	1:A:56:ASP:OD2	1.98	0.63
1:B:97:ALA:HB2	1:B:330:ILE:HD11	1.81	0.63
1:C:206:ASN:N	1:C:206:ASN:OD1	2.31	0.62
1:C:395:ASP:OD1	1:C:451:ILE:CG2	2.48	0.62
1:A:410:PHE:HE2	1:A:419:TYR:CD1	2.18	0.62
1:B:129:THR:HG22	1:B:132:GLU:HG3	1.80	0.62
1:D:182:VAL:O	1:D:366:GLN:NE2	2.32	0.62
1:B:355:ARG:HG3	1:B:355:ARG:HH11	1.64	0.61
1:D:332:PRO:O	1:D:337:MET:HG3	2.01	0.61
1:D:403:ILE:HD12	1:D:428:ILE:HD13	1.81	0.61
1:D:413:ARG:HB2	1:D:414:HIS:ND1	2.15	0.61
1:A:410:PHE:HE2	1:A:419:TYR:CE1	2.18	0.61
1:D:69:GLY:HA3	1:D:333:ASN:O	2.01	0.61
1:D:384:LEU:HD23	1:D:425:GLU:HB2	1.82	0.61
1:A:376:LEU:HD13	1:A:376:LEU:O	2.01	0.60
1:B:424:ILE:CG2	1:B:428:ILE:HG13	2.30	0.60
1:D:98:VAL:O	1:D:105:ILE:HG13	2.01	0.60
1:A:450:THR:HG23	1:A:453:SER:H	1.66	0.60
1:C:137:ASP:OD2	1:C:204:HIS:HA	2.02	0.59
1:C:458:LEU:O	1:C:462:ILE:HG23	2.02	0.59
1:A:49:PRO:HG3	1:A:71:TYR:CE1	2.38	0.59
1:C:69:GLY:HA3	1:C:333:ASN:O	2.01	0.59
2:G:3:GLC:O5	2:G:3:GLC:C6	2.47	0.59
1:A:334:ILE:HG12	1:A:337:MET:HG2	1.84	0.59
1:C:343:ALA:HB2	1:C:369:ASN:ND2	2.18	0.59
1:C:80:ILE:HG22	1:C:82:PRO:HD3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:462:ILE:HG22	1:D:466:LEU:HD11	1.85	0.59
1:A:345:ARG:NH2	2:E:3:GLC:O6	2.36	0.59
1:A:135:ALA:O	1:A:139:GLU:HG3	2.04	0.58
1:A:169:ALA:O	1:A:170:PHE:CD2	2.57	0.57
1:A:410:PHE:CE2	1:A:419:TYR:CD1	2.92	0.57
1:B:131:GLU:N	1:B:131:GLU:OE1	2.38	0.57
1:A:246:THR:OG1	1:A:247:VAL:N	2.36	0.57
1:B:380:PHE:HD1	1:B:384:LEU:HD12	1.70	0.57
1:A:391:TYR:HB2	1:A:396:LEU:CD1	2.34	0.57
1:D:4:GLU:O	1:D:273:ASN:ND2	2.30	0.57
1:D:414:HIS:CD2	1:D:470:HIS:CD2	2.91	0.57
1:C:182:VAL:HG12	1:C:184:VAL:HG23	1.87	0.56
1:D:414:HIS:NE2	1:D:467:GLU:OE2	2.35	0.56
1:C:34:ILE:HD13	1:C:276:LEU:HD21	1.88	0.56
1:A:350:ASN:HB3	1:A:356:GLN:HG2	1.88	0.56
1:A:410:PHE:CE2	1:A:419:TYR:CE1	2.94	0.56
1:D:442:ARG:HG2	1:D:449:GLU:HG2	1.87	0.56
1:A:130:TRP:HA	1:A:133:ILE:HD13	1.88	0.56
1:A:192:GLY:HA2	1:A:251:PHE:CE2	2.41	0.56
1:D:381:PRO:HG2	1:D:384:LEU:HB2	1.86	0.56
1:B:336:GLN:NE2	1:B:373:LYS:HG3	2.21	0.55
1:D:407:LYS:HG3	1:D:424:ILE:HD12	1.88	0.55
1:C:444:VAL:O	1:C:445:LYS:HD3	2.06	0.55
1:C:48:PHE:HA	1:C:51:VAL:CG2	2.36	0.55
1:C:457:TYR:O	1:C:461:THR:OG1	2.23	0.55
1:D:123:LEU:HD22	1:D:224:ALA:HB1	1.88	0.55
1:C:16:LYS:HZ2	1:C:263:LEU:HD22	1.71	0.55
1:A:272:PRO:HG2	1:A:273:ASN:HD21	1.70	0.55
1:B:130:TRP:HB3	1:B:195:PHE:HE2	1.71	0.55
1:B:170:PHE:CE2	1:B:334:ILE:HD11	2.42	0.55
1:C:90:LEU:HD23	1:C:108:PRO:HG2	1.87	0.55
1:D:428:ILE:HG22	1:D:432:LEU:HD12	1.89	0.55
1:A:381:PRO:O	1:A:385:THR:HG23	2.07	0.54
1:B:16:LYS:NZ	2:F:1:GLC:O2	2.39	0.54
1:D:134:PRO:HA	1:D:204:HIS:CD2	2.42	0.54
1:A:209:THR:HA	1:A:213:ILE:HD12	1.89	0.54
1:B:391:TYR:N	1:B:391:TYR:CD2	2.73	0.54
1:C:435:LEU:HD21	1:C:451:ILE:HD13	1.88	0.54
1:D:372:ASP:N	1:D:372:ASP:OD1	2.40	0.54
1:B:49:PRO:HG3	1:B:71:TYR:CE1	2.42	0.54
1:D:155:PRO:HG3	1:D:345:ARG:HA	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:185:ASP:O	1:D:190:LYS:NZ	2.41	0.54
1:D:399:ILE:HD11	1:D:451:ILE:HD11	1.90	0.54
1:B:65:HIS:CE1	1:B:331:MET:HB2	2.43	0.54
1:B:172:TYR:HE1	1:B:175:GLY:HA2	1.73	0.54
1:A:46:GLU:O	1:A:49:PRO:HD2	2.08	0.54
1:D:395:ASP:OD1	1:D:451:ILE:CG2	2.56	0.54
1:A:373:LYS:O	1:A:377:LEU:HG	2.07	0.53
1:B:443:GLY:HA2	1:B:447:GLN:O	2.07	0.53
1:B:218:PHE:HA	1:B:223:THR:HG22	1.89	0.53
1:C:229:GLY:HA3	1:C:231:TRP:CH2	2.44	0.53
1:B:65:HIS:CD2	1:B:97:ALA:HB1	2.43	0.53
1:D:457:TYR:O	1:D:461:THR:OG1	2.25	0.53
1:B:66:ASP:N	1:B:66:ASP:OD1	2.40	0.53
1:C:439:ILE:HD11	1:C:454:MET:HG3	1.91	0.53
1:D:260:VAL:HB	1:D:330:ILE:HD13	1.91	0.53
1:D:129:THR:OG1	1:D:132:GLU:HG3	2.08	0.52
1:D:178:ASP:OD2	1:D:181:ASP:HB2	2.09	0.52
1:D:230:PRO:HA	1:D:233:TRP:CE2	2.44	0.52
1:C:29:GLU:HG3	1:C:35:LYS:HA	1.90	0.52
1:B:49:PRO:HG3	1:B:71:TYR:HE1	1.74	0.52
1:D:118:TYR:CE1	1:D:126:PRO:HD3	2.45	0.52
1:D:378:GLN:O	1:D:379:GLU:HG2	2.09	0.52
1:B:378:GLN:NE2	1:B:379:GLU:H	2.07	0.52
1:B:414:HIS:NE2	1:B:467:GLU:OE2	2.39	0.52
1:C:128:LYS:HD2	1:C:128:LYS:N	2.24	0.52
1:D:384:LEU:O	1:D:388:LEU:HD12	2.10	0.52
1:B:287:THR:HG23	1:B:290:GLY:H	1.73	0.52
1:D:47:LYS:O	1:D:51:VAL:HG22	2.09	0.52
1:D:199:LEU:HD13	1:D:205:MET:HE2	1.92	0.52
1:A:29:GLU:OE2	1:A:35:LYS:HD3	2.10	0.51
1:A:115:SER:OG	1:A:324:ASN:ND2	2.43	0.51
1:C:116:LEU:HB2	1:C:248:LEU:HD23	1.92	0.51
1:C:234:SER:HB3	1:C:299:PRO:HD3	1.93	0.51
1:C:406:ALA:HB3	1:C:462:ILE:HD11	1.93	0.51
1:B:129:THR:CG2	1:B:132:GLU:HG3	2.39	0.51
2:E:2:GLC:O3	2:E:3:GLC:O2	2.28	0.51
1:D:134:PRO:HA	1:D:204:HIS:HD2	1.75	0.51
1:D:471:ALA:O	1:D:475:ARG:HB2	2.11	0.51
1:C:312:LEU:HB3	1:C:318:ILE:HD13	1.93	0.51
2:H:3:GLC:O5	2:H:3:GLC:C6	2.51	0.51
1:D:407:LYS:O	1:D:411:ASN:HB2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:424:ILE:O	1:D:427:GLU:HG2	2.11	0.51
1:A:168:TYR:CD1	1:A:168:TYR:C	2.89	0.50
1:A:155:PRO:CG	1:A:345:ARG:HG3	2.41	0.50
1:C:179:ILE:H	1:C:179:ILE:CD1	2.18	0.50
1:A:399:ILE:HG23	1:A:458:LEU:HD21	1.93	0.50
1:A:424:ILE:HG23	1:A:466:LEU:HD21	1.93	0.50
1:D:32:THR:HG22	1:D:34:ILE:HD12	1.93	0.50
1:A:155:PRO:HG3	1:A:345:ARG:CA	2.41	0.50
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.46	0.50
1:A:129:THR:HB	1:A:131:GLU:HG2	1.94	0.50
1:D:86:PHE:O	1:D:89:LYS:HB2	2.12	0.50
1:D:424:ILE:HD13	1:D:466:LEU:HD21	1.94	0.50
1:A:368:HIS:ND1	1:A:369:ASN:N	2.60	0.50
1:A:192:GLY:HA2	1:A:251:PHE:HE2	1.76	0.50
1:B:69:GLY:HA3	1:B:333:ASN:O	2.11	0.50
1:B:399:ILE:O	1:B:403:ILE:HG12	2.12	0.50
1:C:413:ARG:HG2	1:C:414:HIS:ND1	2.25	0.49
1:C:407:LYS:O	1:C:411:ASN:HB2	2.12	0.49
1:A:182:VAL:HG12	1:A:184:VAL:HG23	1.94	0.49
1:A:63:TRP:HB3	1:A:68:PHE:HE1	1.77	0.49
1:A:148:LEU:HD22	1:A:150:PHE:HD2	1.78	0.49
1:A:195:PHE:O	1:A:198:ASP:HB2	2.12	0.49
1:D:391:TYR:HB2	1:D:396:LEU:HD12	1.94	0.49
1:A:471:ALA:O	1:A:475:ARG:HG3	2.12	0.49
1:C:218:PHE:HA	1:C:223:THR:HG22	1.95	0.49
1:B:439:ILE:HG23	1:B:449:GLU:HB3	1.94	0.49
1:C:123:LEU:HD11	1:C:224:ALA:HB1	1.95	0.49
1:D:118:TYR:CE2	1:D:126:PRO:HG3	2.48	0.49
1:A:405:LYS:HG2	1:A:459:MET:HE1	1.94	0.49
1:C:48:PHE:HA	1:C:51:VAL:HG23	1.95	0.49
1:C:49:PRO:HG3	1:C:71:TYR:CE1	2.48	0.49
1:A:376:LEU:CD1	1:A:400:LYS:HD2	2.38	0.48
1:D:474:MET:O	1:D:477:LYS:HD3	2.13	0.48
1:D:462:ILE:HG22	1:D:466:LEU:CD1	2.43	0.48
1:A:155:PRO:HD2	2:E:2:GLC:O6	2.13	0.48
1:C:460:GLN:O	1:C:464:THR:OG1	2.31	0.48
1:D:350:ASN:HB3	1:D:356:GLN:HB2	1.95	0.48
1:A:49:PRO:HG3	1:A:71:TYR:HE1	1.77	0.48
1:C:395:ASP:OD1	1:C:451:ILE:HG21	2.11	0.48
1:B:230:PRO:HA	1:B:233:TRP:CE2	2.48	0.48
1:A:149:MET:HB2	1:A:223:THR:HG21	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:LEU:HD12	1:A:207:ALA:HA	1.95	0.48
1:A:376:LEU:HD11	1:A:400:LYS:CD	2.40	0.48
1:B:355:ARG:HG3	1:B:355:ARG:NH1	2.28	0.48
1:A:129:THR:OG1	1:A:132:GLU:HG2	2.13	0.48
1:A:168:TYR:CE1	1:A:183:GLY:N	2.82	0.48
1:C:233:TRP:HB2	1:C:299:PRO:HG2	1.95	0.48
1:C:255:PRO:HG2	1:C:257:LYS:NZ	2.28	0.48
1:C:395:ASP:OD1	1:C:451:ILE:HG22	2.14	0.48
1:A:169:ALA:C	1:A:170:PHE:CD2	2.92	0.48
1:B:382:GLU:HA	1:B:385:THR:HG22	1.96	0.48
1:D:384:LEU:HD11	1:D:428:ILE:HG21	1.95	0.48
1:A:168:TYR:CE1	1:A:183:GLY:CA	2.96	0.48
1:C:428:ILE:O	1:C:432:LEU:HD12	2.14	0.48
2:F:3:GLC:O5	2:F:3:GLC:C6	2.50	0.48
1:B:248:LEU:HB3	1:B:256:SER:HB2	1.96	0.47
1:C:377:LEU:C	1:C:377:LEU:HD23	2.39	0.47
1:D:190:LYS:HD2	1:D:359:ASP:OD2	2.13	0.47
1:C:439:ILE:CD1	1:C:454:MET:HG3	2.44	0.47
1:D:435:LEU:HD22	1:D:458:LEU:HD22	1.96	0.47
1:D:442:ARG:HG2	1:D:449:GLU:CG	2.44	0.47
1:A:90:LEU:CD1	1:A:108:PRO:HG2	2.45	0.47
1:A:410:PHE:CE2	1:A:419:TYR:HD1	2.32	0.47
1:A:421:LEU:HD12	1:A:421:LEU:N	2.22	0.47
1:B:41:PRO:HG2	1:B:44:LEU:HB3	1.97	0.47
1:B:91:TYR:HE2	1:B:309:GLU:HG2	1.78	0.47
1:D:190:LYS:O	1:D:194:THR:OG1	2.32	0.47
1:A:424:ILE:HG22	1:A:428:ILE:HG13	1.96	0.47
1:A:159:TRP:CH2	1:A:340:PHE:HZ	2.33	0.46
1:A:83:ASP:OD1	1:A:83:ASP:N	2.48	0.46
1:A:151:ASN:HB3	1:A:211:TYR:N	2.29	0.46
1:B:312:LEU:HB3	1:B:318:ILE:HD13	1.97	0.46
1:A:19:ASN:HB2	1:A:297:ASP:OD2	2.15	0.46
1:A:123:LEU:HD13	1:A:123:LEU:HA	1.68	0.46
1:B:391:TYR:HB2	1:B:396:LEU:HD12	1.98	0.46
1:D:475:ARG:HG2	1:D:475:ARG:NH1	2.29	0.46
1:B:42:ASP:N	1:B:42:ASP:OD2	2.48	0.46
1:C:114:LEU:HD21	1:C:157:PHE:CD2	2.51	0.46
1:D:414:HIS:ND1	1:D:414:HIS:N	2.63	0.46
1:D:170:PHE:CD2	1:D:334:ILE:HD11	2.51	0.46
1:A:3:ILE:HG13	1:A:57:GLY:O	2.16	0.46
1:A:456:ALA:O	1:A:460:GLN:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:PHE:HD1	1:C:284:TYR:CD2	2.34	0.46
1:C:287:THR:HG23	1:C:290:GLY:H	1.79	0.46
1:D:91:TYR:CE2	1:D:309:GLU:HG3	2.42	0.46
1:A:472:LEU:O	1:A:476:ARG:HB2	2.16	0.46
1:B:420:MET:HE3	1:B:421:LEU:N	2.24	0.46
1:B:130:TRP:HB3	1:B:195:PHE:CE2	2.50	0.45
1:B:155:PRO:HD3	1:B:345:ARG:HG3	1.97	0.45
1:C:229:GLY:HA3	1:C:231:TRP:CZ3	2.51	0.45
1:C:396:LEU:HD23	1:C:396:LEU:HA	1.80	0.45
1:D:28:PHE:HA	1:D:284:TYR:CE1	2.51	0.45
1:D:458:LEU:O	1:D:462:ILE:HG13	2.16	0.45
1:B:114:LEU:HD12	1:B:114:LEU:HA	1.76	0.45
1:B:315:ASP:HA	1:B:316:PRO:HD2	1.83	0.45
1:B:67:ARG:NH2	2:F:3:GLC:O3	2.36	0.45
1:B:296:LYS:NZ	1:B:296:LYS:CB	2.73	0.45
1:A:90:LEU:HB2	1:A:95:TRP:NE1	2.32	0.45
1:A:200:ILE:HD13	1:A:207:ALA:HB2	1.98	0.45
1:A:457:TYR:O	1:A:461:THR:OG1	2.32	0.45
1:B:116:LEU:HB2	1:B:248:LEU:HD12	1.97	0.45
1:C:107:TYR:CD1	1:C:281:LEU:HD13	2.51	0.45
1:A:172:TYR:CG	1:A:172:TYR:O	2.69	0.45
1:D:154:GLU:OE1	2:H:2:GLC:O6	2.21	0.45
1:A:368:HIS:ND1	1:A:368:HIS:C	2.74	0.45
1:B:252:LYS:HB3	1:B:252:LYS:HE2	1.78	0.45
1:C:194:THR:HA	1:C:358:VAL:HG21	1.99	0.45
1:C:382:GLU:HA	1:C:385:THR:OG1	2.16	0.45
1:D:32:THR:CG2	1:D:34:ILE:HD12	2.46	0.45
1:C:382:GLU:O	1:C:385:THR:OG1	2.29	0.45
1:D:194:THR:HA	1:D:358:VAL:HG21	1.98	0.45
1:A:145:LYS:HD2	1:A:221:GLY:O	2.17	0.44
1:C:411:ASN:O	1:C:414:HIS:O	2.34	0.44
1:A:162:ILE:HG23	1:A:189:ALA:HB1	1.99	0.44
1:B:193:LEU:O	1:B:197:VAL:HG23	2.18	0.44
1:D:63:TRP:HB3	1:D:68:PHE:HE1	1.82	0.44
1:B:261:GLY:HA3	1:B:331:MET:HE2	2.00	0.44
1:A:384:LEU:O	1:A:388:LEU:HD23	2.17	0.44
1:B:91:TYR:CE1	1:B:306:LYS:HG2	2.51	0.44
1:C:417:MET:O	1:C:418:HIS:C	2.60	0.44
1:A:158:THR:HG21	1:A:196:LEU:HD22	1.99	0.44
1:B:61:ILE:HG13	1:B:62:PHE:N	2.31	0.44
1:C:407:LYS:HE3	1:C:420:MET:SD	2.58	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:GLY:HA3	1:A:333:ASN:O	2.17	0.44
1:B:273:ASN:HB3	1:B:276:LEU:HB2	2.00	0.44
1:C:43:LYS:HB3	1:C:46:GLU:HG3	1.99	0.44
1:C:67:ARG:NH2	2:G:3:GLC:O3	2.34	0.44
1:A:84:LYS:NZ	1:A:88:ASP:OD2	2.51	0.44
1:A:131:GLU:HG2	1:A:131:GLU:H	1.42	0.44
1:A:168:TYR:CD1	1:A:168:TYR:O	2.70	0.44
1:A:417:MET:HE2	1:A:417:MET:HB3	1.87	0.44
1:D:395:ASP:OD1	1:D:451:ILE:HG22	2.17	0.44
1:D:429:LEU:O	1:D:433:LYS:HG3	2.18	0.44
1:A:169:ALA:O	1:A:170:PHE:CG	2.71	0.44
1:C:449:GLU:OE2	1:C:449:GLU:HA	2.18	0.44
1:B:172:TYR:CE1	1:B:175:GLY:HA2	2.53	0.43
1:D:129:THR:HG22	1:D:250:THR:OG1	2.18	0.43
1:D:384:LEU:HD11	1:D:428:ILE:CG2	2.48	0.43
1:D:384:LEU:HD12	1:D:388:LEU:HD11	1.99	0.43
1:D:450:THR:O	1:D:453:SER:N	2.50	0.43
1:B:133:ILE:HD12	1:B:225:MET:HE1	2.00	0.43
1:D:150:PHE:HZ	1:D:196:LEU:HD11	1.83	0.43
1:A:383:GLN:HB3	1:A:425:GLU:OE2	2.18	0.43
1:B:12:ILE:HA	1:B:62:PHE:HB2	1.99	0.43
1:A:154:GLU:HA	1:A:155:PRO:HD3	1.85	0.43
1:A:205:MET:HB3	1:A:205:MET:HE2	1.83	0.43
1:D:383:GLN:NE2	1:D:425:GLU:OE1	2.51	0.43
1:A:210:ASP:H	1:A:213:ILE:HD12	1.82	0.43
1:B:312:LEU:HB3	1:B:318:ILE:CD1	2.48	0.43
1:B:414:HIS:HD2	1:B:470:HIS:CD2	2.35	0.43
1:B:424:ILE:HG22	1:B:424:ILE:O	2.16	0.43
1:D:151:ASN:HB3	1:D:157:PHE:CD1	2.53	0.43
1:B:122:LEU:HD23	1:B:224:ALA:HB2	1.99	0.43
1:D:41:PRO:HB2	1:D:47:LYS:NZ	2.34	0.43
1:D:427:GLU:OE2	1:D:427:GLU:N	2.52	0.43
1:B:112:GLU:HG2	1:B:261:GLY:O	2.18	0.43
1:B:114:LEU:HG	1:B:227:ILE:HG22	2.00	0.43
1:C:29:GLU:CG	1:C:35:LYS:HA	2.48	0.43
1:D:41:PRO:HG2	1:D:44:LEU:HB3	2.00	0.43
1:A:10:ILE:HB	1:A:38:VAL:HG22	2.00	0.43
1:A:372:ASP:HB3	1:A:375:MET:HB3	2.01	0.43
1:D:229:GLY:HA3	1:D:231:TRP:CH2	2.53	0.43
1:C:49:PRO:HG3	1:C:71:TYR:HE1	1.82	0.43
1:A:43:LYS:HD3	1:A:46:GLU:OE2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:LEU:H	1:A:421:LEU:CD1	2.20	0.43
1:A:185:ASP:CB	1:A:366:GLN:OE1	2.67	0.42
1:D:157:PHE:O	1:D:160:PRO:HD2	2.19	0.42
1:A:84:LYS:HD2	1:A:84:LYS:HA	1.81	0.42
1:B:202:ASN:O	1:B:203:LYS:HB2	2.19	0.42
1:D:107:TYR:OH	1:D:278:LYS:HD3	2.19	0.42
1:B:128:LYS:HE2	1:B:128:LYS:HB3	1.88	0.42
1:C:255:PRO:HG2	1:C:257:LYS:HZ2	1.83	0.42
1:A:427:GLU:HG3	1:A:466:LEU:HD23	2.00	0.42
1:B:134:PRO:HB3	1:B:204:HIS:NE2	2.35	0.42
1:B:392:ASP:OD1	1:B:392:ASP:N	2.52	0.42
1:C:438:ALA:O	1:C:442:ARG:HB2	2.19	0.42
1:D:62:PHE:HE1	1:D:109:ILE:HG13	1.83	0.42
1:D:83:ASP:OD1	1:D:83:ASP:N	2.52	0.42
1:A:315:ASP:HB3	1:A:318:ILE:HG12	2.01	0.42
1:B:259:PHE:HD1	1:B:259:PHE:HA	1.71	0.42
1:A:410:PHE:CE2	1:A:419:TYR:HE1	2.38	0.42
1:C:148:LEU:HB2	1:C:225:MET:HE3	2.02	0.42
1:D:155:PRO:HG3	1:D:345:ARG:CA	2.49	0.42
1:C:141:LYS:HB2	1:C:141:LYS:HE2	1.79	0.42
1:A:65:HIS:CE1	1:A:331:MET:HB2	2.55	0.42
1:A:291:LEU:HA	1:A:291:LEU:HD23	1.81	0.42
1:B:155:PRO:HD2	2:F:2:GLC:O6	2.20	0.42
1:D:395:ASP:OD1	1:D:451:ILE:HG21	2.19	0.42
1:D:424:ILE:HD13	1:D:424:ILE:HG21	1.79	0.42
1:D:454:MET:HE3	1:D:454:MET:HB3	1.74	0.42
1:C:250:THR:HG22	1:C:255:PRO:HA	2.02	0.42
1:D:298:LYS:HA	1:D:299:PRO:HD3	1.89	0.42
1:D:431:SER:HA	1:D:434:ARG:HG3	2.01	0.42
1:C:12:ILE:HA	1:C:62:PHE:HB2	2.02	0.41
1:D:43:LYS:HB2	1:D:43:LYS:HE3	1.89	0.41
1:A:66:ASP:OD2	2:E:2:GLC:O3	2.38	0.41
1:B:117:ILE:HG22	1:B:218:PHE:CZ	2.54	0.41
1:C:377:LEU:HD21	1:C:385:THR:CG2	2.50	0.41
1:A:159:TRP:NE1	1:A:163:ALA:HB2	2.35	0.41
1:A:196:LEU:O	1:A:196:LEU:HG	2.20	0.41
1:D:87:GLN:HG3	1:D:95:TRP:CE2	2.55	0.41
1:B:463:LEU:HD23	1:B:463:LEU:HA	1.90	0.41
1:A:123:LEU:HD13	1:A:124:PRO:HD2	2.02	0.41
1:C:66:ASP:HA	1:C:333:ASN:HA	2.02	0.41
1:C:91:TYR:HA	1:C:92:PRO:HD3	1.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:11:TRP:HB3	1:D:44:LEU:HD13	2.03	0.41
1:D:154:GLU:HA	1:D:155:PRO:HD3	1.96	0.41
1:B:380:PHE:HD1	1:B:384:LEU:CD1	2.34	0.41
1:C:407:LYS:HE3	1:C:420:MET:HA	2.02	0.41
1:A:308:TYR:CE2	1:A:312:LEU:HD11	2.56	0.41
1:C:76:LEU:N	1:C:76:LEU:HD13	2.36	0.41
1:C:393:TYR:CE2	1:C:394:ARG:NH1	2.89	0.41
1:C:424:ILE:O	1:C:428:ILE:HG13	2.21	0.41
1:D:373:LYS:HG3	1:D:393:TYR:HD1	1.85	0.41
1:D:107:TYR:HA	1:D:108:PRO:HD3	1.87	0.41
1:D:170:PHE:HE1	1:D:182:VAL:HG22	1.85	0.41
1:A:404:LEU:HD12	1:A:404:LEU:HA	1.88	0.41
1:D:216:ALA:HB1	1:D:220:LYS:NZ	2.36	0.41
1:D:424:ILE:HG22	1:D:428:ILE:HG13	2.03	0.41
1:A:168:TYR:CD1	1:A:183:GLY:HA3	2.55	0.41
1:B:194:THR:HA	1:B:358:VAL:HG21	2.02	0.41
1:B:378:GLN:CD	1:B:379:GLU:H	2.28	0.41
1:D:48:PHE:HB3	1:D:49:PRO:HD3	2.03	0.41
1:A:184:VAL:HG12	1:A:362:LEU:HD22	2.02	0.40
1:C:179:ILE:CG2	1:C:336:GLN:HG2	2.50	0.40
1:C:382:GLU:H	1:C:382:GLU:HG2	1.45	0.40
1:D:90:LEU:HB2	1:D:95:TRP:NE1	2.36	0.40
1:A:11:TRP:HB3	1:A:44:LEU:HD13	2.03	0.40
1:B:168:TYR:CE2	1:B:183:GLY:HA3	2.56	0.40
1:C:52:ALA:HA	1:C:56:ASP:O	2.21	0.40
1:D:68:PHE:HB3	1:D:105:ILE:HD13	2.03	0.40
1:A:61:ILE:HG13	1:A:62:PHE:N	2.35	0.40
1:B:271:SER:HA	1:B:272:PRO:HD3	1.83	0.40
1:D:377:LEU:O	1:D:380:PHE:HB2	2.22	0.40
1:A:167:GLY:HA3	1:A:189:ALA:HB2	2.04	0.40
1:B:233:TRP:HB2	1:B:299:PRO:HG2	2.04	0.40
1:C:47:LYS:HE3	1:C:47:LYS:HB2	1.82	0.40
1:C:109:ILE:HA	1:C:109:ILE:HD13	1.84	0.40
1:C:193:LEU:O	1:C:197:VAL:HG23	2.21	0.40
1:D:170:PHE:CE1	1:D:182:VAL:HG22	2.57	0.40
1:A:232:ALA:O	1:A:236:ILE:HG13	2.21	0.40
1:C:235:ASN:OD1	1:C:298:LYS:HE3	2.21	0.40
1:D:376:LEU:HD23	1:D:396:LEU:HD23	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	A	476/487 (98%)	463 (97%)	13 (3%)	0	100	100
1	B	477/487 (98%)	472 (99%)	5 (1%)	0	100	100
1	C	474/487 (97%)	469 (99%)	5 (1%)	0	100	100
1	D	468/487 (96%)	464 (99%)	4 (1%)	0	100	100
All	All	1895/1948 (97%)	1868 (99%)	27 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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](#)

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	1.26	1 (8%)	17,17,17	1.60	4 (23%)
2	GLC	E	2	2	11,11,12	1.98	4 (36%)	15,15,17	1.61	4 (26%)
2	GLC	E	3	2	11,11,12	7.23	7 (63%)	15,15,17	4.66	8 (53%)
2	GLC	F	1	2	12,12,12	1.19	0	17,17,17	1.30	3 (17%)
2	GLC	F	2	2	11,11,12	1.66	3 (27%)	15,15,17	1.21	1 (6%)
2	GLC	F	3	2	11,11,12	7.02	8 (72%)	15,15,17	3.42	5 (33%)
2	GLC	G	1	2	12,12,12	1.18	3 (25%)	17,17,17	1.79	6 (35%)
2	GLC	G	2	2	11,11,12	1.67	3 (27%)	15,15,17	1.63	4 (26%)
2	GLC	G	3	2	11,11,12	6.87	7 (63%)	15,15,17	3.72	4 (26%)
2	GLC	H	1	2	12,12,12	1.23	1 (8%)	17,17,17	1.40	2 (11%)
2	GLC	H	2	2	11,11,12	1.73	3 (27%)	15,15,17	1.12	2 (13%)
2	GLC	H	3	2	11,11,12	6.95	8 (72%)	15,15,17	2.86	5 (33%)

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	-	1/2/22/22	0/1/1/1
2	GLC	E	2	2	-	0/2/19/22	0/1/1/1
2	GLC	E	3	2	-	2/2/19/22	0/1/1/1
2	GLC	F	1	2	-	2/2/22/22	0/1/1/1
2	GLC	F	2	2	-	0/2/19/22	0/1/1/1
2	GLC	F	3	2	-	2/2/19/22	0/1/1/1
2	GLC	G	1	2	-	0/2/22/22	0/1/1/1
2	GLC	G	2	2	-	1/2/19/22	0/1/1/1
2	GLC	G	3	2	-	2/2/19/22	0/1/1/1
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	GLC	H	3	2	-	2/2/19/22	0/1/1/1

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	3	GLC	O5-C1	17.68	1.73	1.43
2	F	3	GLC	O5-C1	16.84	1.72	1.43
2	G	3	GLC	O5-C1	16.59	1.71	1.43
2	H	3	GLC	O5-C1	16.58	1.71	1.43
2	H	3	GLC	O5-C5	10.94	1.64	1.43
2	F	3	GLC	O5-C5	10.82	1.64	1.43
2	E	3	GLC	O5-C5	10.56	1.64	1.43
2	G	3	GLC	O5-C5	10.29	1.63	1.43
2	H	3	GLC	O3-C3	7.89	1.62	1.43
2	E	3	GLC	O3-C3	7.79	1.62	1.43
2	F	3	GLC	O3-C3	7.78	1.62	1.43
2	G	3	GLC	O3-C3	7.60	1.61	1.43
2	E	3	GLC	C2-C3	-6.84	1.42	1.52
2	F	3	GLC	C2-C3	-6.53	1.42	1.52
2	G	3	GLC	C2-C3	-6.39	1.42	1.52
2	H	3	GLC	C2-C3	-5.87	1.43	1.52
2	E	3	GLC	C4-C3	-4.96	1.39	1.52
2	G	3	GLC	C4-C3	-4.29	1.41	1.52
2	H	3	GLC	C4-C3	-4.00	1.41	1.52
2	F	3	GLC	C4-C3	-3.99	1.42	1.52
2	E	2	GLC	O5-C1	3.62	1.49	1.43
2	G	2	GLC	O5-C1	3.60	1.49	1.43
2	H	2	GLC	O5-C1	3.60	1.49	1.43
2	F	2	GLC	O5-C1	3.51	1.49	1.43
2	E	3	GLC	O2-C2	3.16	1.50	1.43
2	H	3	GLC	O2-C2	3.08	1.49	1.43
2	E	2	GLC	C2-C3	-3.06	1.47	1.52
2	G	3	GLC	O2-C2	3.00	1.49	1.43
2	E	2	GLC	O5-C5	2.81	1.48	1.43
2	F	3	GLC	O2-C2	2.79	1.49	1.43
2	E	1	GLC	O5-C1	2.73	1.49	1.42
2	G	3	GLC	C6-C5	-2.63	1.43	1.51
2	H	3	GLC	O4-C4	2.60	1.49	1.43
2	F	3	GLC	C6-C5	-2.56	1.43	1.51
2	F	2	GLC	O5-C5	2.56	1.48	1.43
2	E	2	GLC	O4-C4	2.49	1.49	1.43
2	H	2	GLC	O5-C5	2.45	1.48	1.43
2	E	3	GLC	C6-C5	-2.44	1.43	1.51
2	H	3	GLC	C6-C5	-2.43	1.43	1.51
2	F	3	GLC	O4-C4	2.39	1.48	1.43
2	F	2	GLC	C2-C3	-2.31	1.48	1.52
2	G	2	GLC	O5-C5	2.29	1.47	1.43
2	H	1	GLC	O5-C1	2.28	1.48	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	2	GLC	O4-C4	2.13	1.48	1.43
2	G	1	GLC	O2-C2	2.08	1.48	1.43
2	G	2	GLC	C2-C3	-2.08	1.49	1.52
2	G	1	GLC	O3-C3	2.04	1.48	1.43
2	G	1	GLC	O5-C1	2.04	1.47	1.42

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	3	GLC	C1-O5-C5	-11.58	96.66	112.19
2	E	3	GLC	C2-C3-C4	10.06	128.55	110.86
2	G	3	GLC	C1-O5-C5	-9.90	98.92	112.19
2	F	3	GLC	C1-O5-C5	-9.69	99.20	112.19
2	G	3	GLC	C2-C3-C4	8.51	125.83	110.86
2	H	3	GLC	C1-O5-C5	-7.68	101.90	112.19
2	F	3	GLC	C2-C3-C4	7.16	123.46	110.86
2	H	3	GLC	C2-C3-C4	6.52	122.32	110.86
2	E	3	GLC	O5-C5-C6	5.49	118.34	107.66
2	G	3	GLC	O3-C3-C2	-4.82	100.21	110.05
2	E	3	GLC	O3-C3-C2	-4.19	101.50	110.05
2	F	3	GLC	O3-C3-C2	-3.60	102.70	110.05
2	G	1	GLC	C4-C3-C2	3.51	116.98	110.83
2	E	2	GLC	O3-C3-C2	-3.27	103.38	110.05
2	E	3	GLC	C6-C5-C4	-3.25	105.03	113.02
2	E	1	GLC	O5-C5-C4	3.07	115.23	109.70
2	H	2	GLC	C1-O5-C5	2.98	116.19	112.19
2	G	2	GLC	C1-C2-C3	2.98	113.98	109.64
2	F	2	GLC	O4-C4-C3	-2.97	103.39	110.38
2	E	3	GLC	C1-C2-C3	2.94	113.92	109.64
2	E	2	GLC	C1-O5-C5	2.92	116.10	112.19
2	F	1	GLC	O4-C4-C3	-2.88	103.59	110.38
2	G	1	GLC	C1-C2-C3	2.86	116.20	110.36
2	E	3	GLC	O3-C3-C4	-2.84	103.67	110.38
2	H	1	GLC	O5-C5-C4	2.84	114.81	109.70
2	G	2	GLC	O5-C5-C4	2.74	117.50	110.83
2	G	1	GLC	C1-O5-C5	-2.69	108.46	113.65
2	G	2	GLC	O4-C4-C5	-2.68	102.73	109.32
2	H	3	GLC	O3-C3-C2	-2.65	104.65	110.05
2	G	3	GLC	C1-C2-C3	2.62	113.45	109.64
2	H	1	GLC	C4-C3-C2	2.61	115.42	110.83
2	F	1	GLC	C1-O5-C5	-2.58	108.66	113.65
2	G	1	GLC	O5-C5-C4	2.58	114.34	109.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	1	GLC	C3-C4-C5	2.56	114.86	110.23
2	E	2	GLC	C2-C3-C4	2.53	115.31	110.86
2	E	3	GLC	O2-C2-C1	2.47	114.87	109.22
2	F	3	GLC	O5-C5-C4	2.47	116.83	110.83
2	E	1	GLC	C1-C2-C3	2.42	115.29	110.36
2	F	1	GLC	O1-C1-O5	-2.30	103.59	110.41
2	H	3	GLC	O5-C5-C4	2.29	116.41	110.83
2	E	1	GLC	O3-C3-C2	-2.19	105.22	110.38
2	G	2	GLC	C1-O5-C5	2.13	115.05	112.19
2	E	1	GLC	C6-C5-C4	-2.11	107.85	113.02
2	H	2	GLC	O5-C5-C4	2.07	115.87	110.83
2	E	2	GLC	O2-C2-C3	-2.07	105.87	110.15
2	H	3	GLC	C6-C5-C4	-2.06	107.95	113.02
2	G	1	GLC	O4-C4-C5	-2.01	104.37	109.32
2	F	3	GLC	C6-C5-C4	-2.01	108.08	113.02

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	3	GLC	O5-C5-C6-O6
2	E	3	GLC	O5-C5-C6-O6
2	G	3	GLC	O5-C5-C6-O6
2	H	3	GLC	O5-C5-C6-O6
2	F	3	GLC	C4-C5-C6-O6
2	F	1	GLC	C4-C5-C6-O6
2	G	3	GLC	C4-C5-C6-O6
2	E	3	GLC	C4-C5-C6-O6
2	H	1	GLC	C4-C5-C6-O6
2	F	1	GLC	O5-C5-C6-O6
2	H	3	GLC	C4-C5-C6-O6
2	H	1	GLC	O5-C5-C6-O6
2	E	1	GLC	O5-C5-C6-O6
2	G	2	GLC	C4-C5-C6-O6

There are no ring outliers.

9 monomers are involved in 26 short contacts:

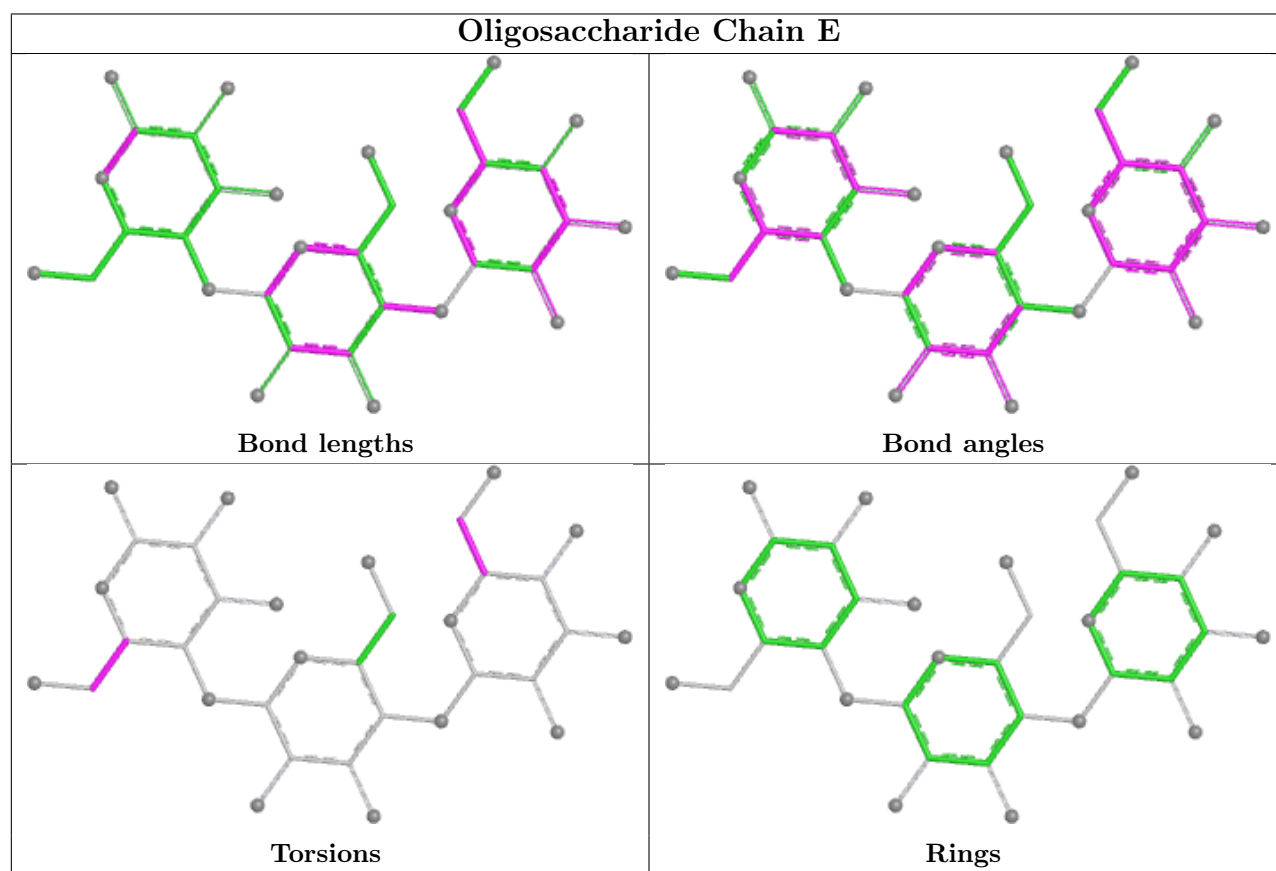
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	3	GLC	6	0
2	F	1	GLC	1	0

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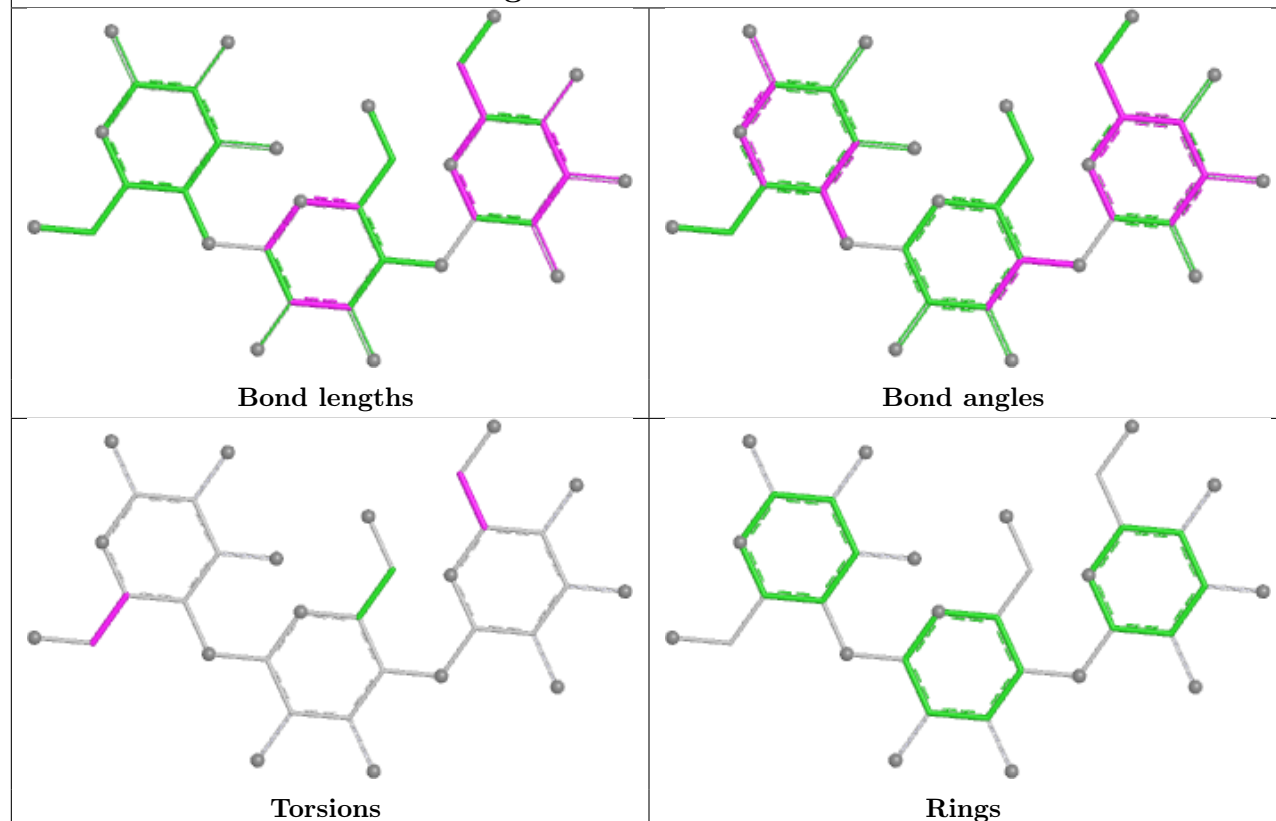
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	1	GLC	1	0
2	F	3	GLC	5	0
2	E	2	GLC	3	0
2	F	2	GLC	1	0
2	H	2	GLC	1	0
2	H	3	GLC	4	0
2	G	3	GLC	5	0

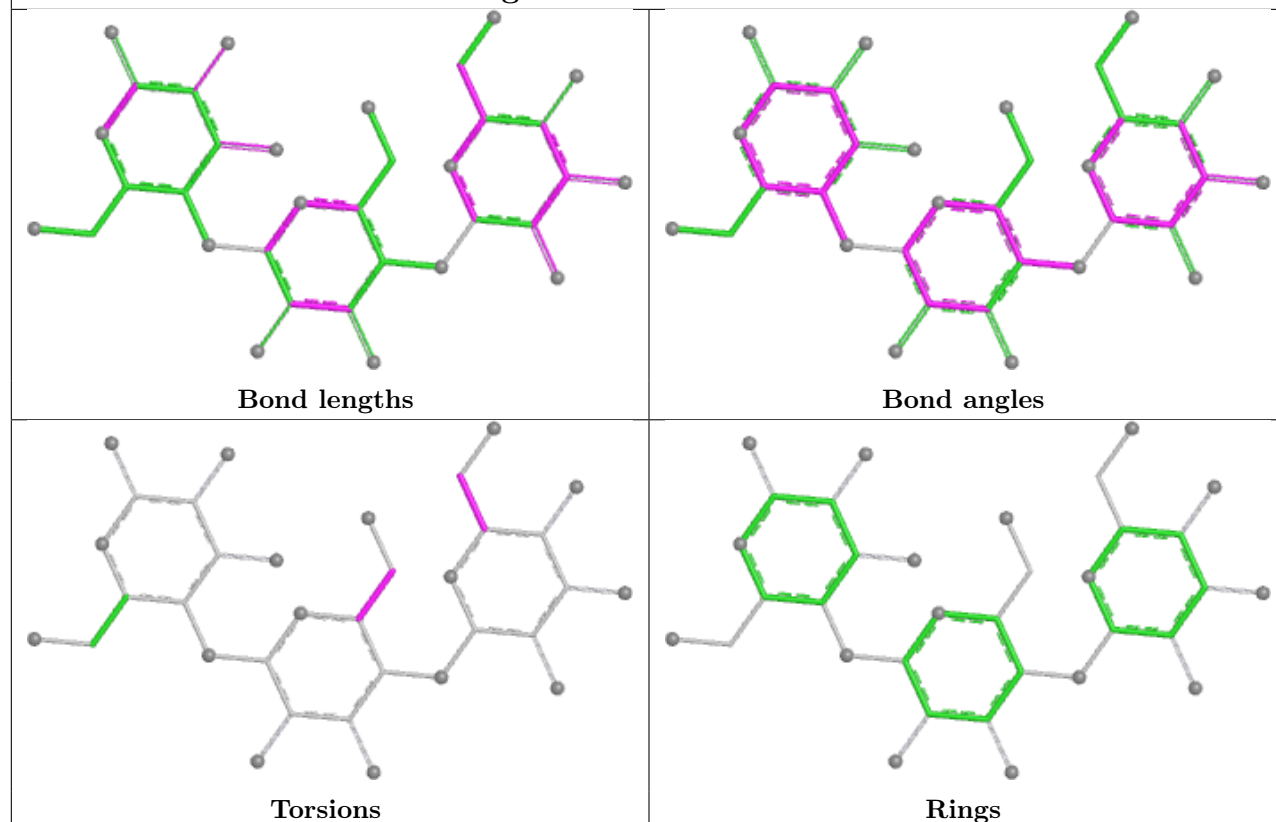
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

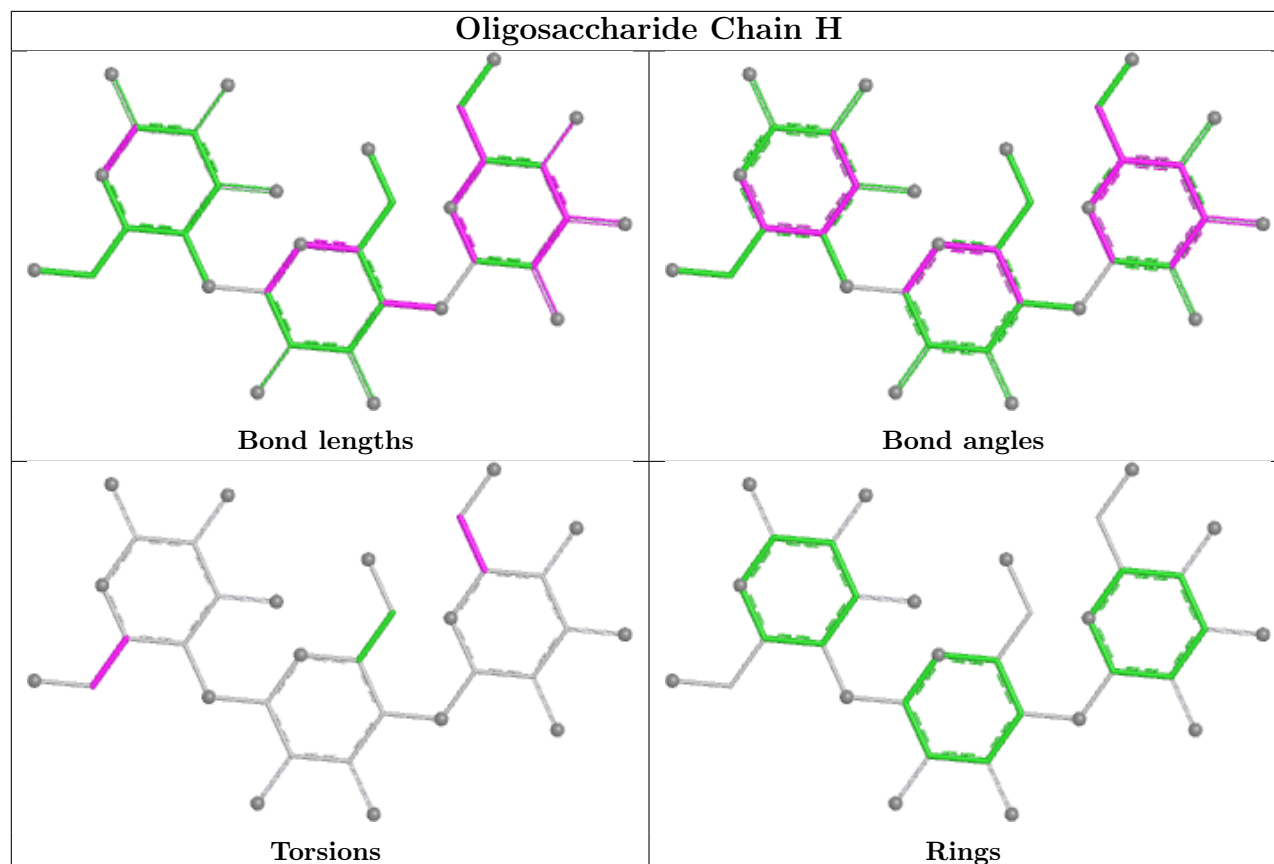


Oligosaccharide Chain F



Oligosaccharide Chain G





5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	478/487 (98%)	-0.18	5 (1%) 79 66	31, 61, 95, 121	0
1	B	479/487 (98%)	-0.11	11 (2%) 61 46	39, 66, 99, 135	0
1	C	476/487 (97%)	-0.23	5 (1%) 78 65	41, 68, 105, 140	0
1	D	472/487 (96%)	-0.08	7 (1%) 72 57	45, 73, 100, 138	0
All	All	1905/1948 (97%)	-0.15	28 (1%) 72 57	31, 67, 101, 140	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	321	THR	3.7
1	D	435	LEU	3.7
1	B	69	GLY	3.4
1	C	301	GLY	3.2
1	A	264	SER	3.1
1	D	109	ILE	3.1
1	B	347	ALA	3.0
1	D	164	ALA	2.8
1	B	392	ASP	2.7
1	B	324	ASN	2.7
1	B	349	ILE	2.7
1	D	340	PHE	2.6
1	C	165	ASP	2.5
1	A	291	LEU	2.5
1	C	375	MET	2.4
1	D	330	ILE	2.4
1	B	55	GLY	2.4
1	D	313	ALA	2.3
1	B	371	ASP	2.3
1	A	398	ILE	2.3
1	B	148	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	365	ALA	2.2
1	C	169	ALA	2.2
1	A	313	ALA	2.2
1	D	328	GLY	2.1
1	C	409	SER	2.1
1	B	322	MET	2.0
1	B	337	MET	2.0

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

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

6.3 Carbohydrates [i](#)

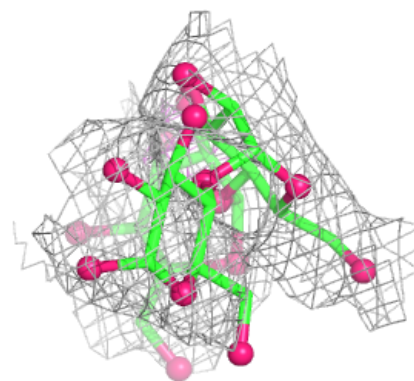
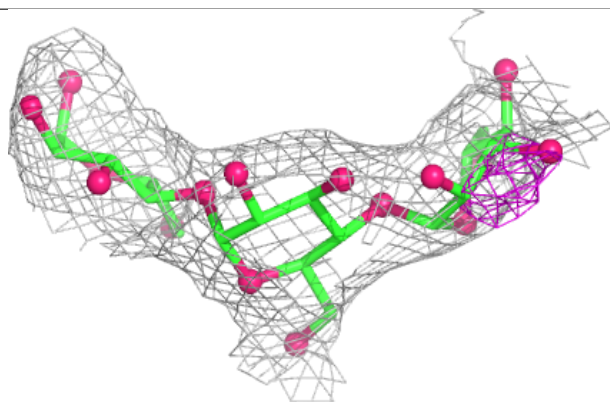
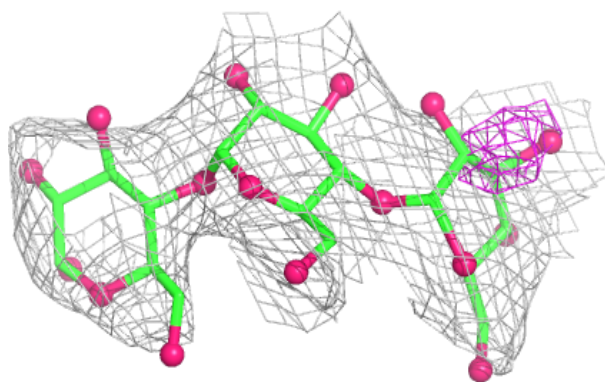
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	H	1	12/12	0.70	0.18	49,60,68,72	0
2	GLC	H	3	11/12	0.73	0.11	45,65,74,74	0
2	GLC	G	1	12/12	0.77	0.15	57,64,72,78	0
2	GLC	H	2	11/12	0.78	0.08	47,55,59,66	0
2	GLC	F	3	11/12	0.79	0.09	52,61,67,76	0
2	GLC	E	3	11/12	0.80	0.10	37,55,66,76	0
2	GLC	F	1	12/12	0.82	0.14	47,52,59,61	0
2	GLC	G	3	11/12	0.87	0.07	61,67,79,81	0
2	GLC	F	2	11/12	0.88	0.09	33,47,57,60	0
2	GLC	E	1	12/12	0.91	0.07	36,44,53,54	0
2	GLC	E	2	11/12	0.93	0.07	30,33,36,37	0
2	GLC	G	2	11/12	0.95	0.05	53,60,69,70	0

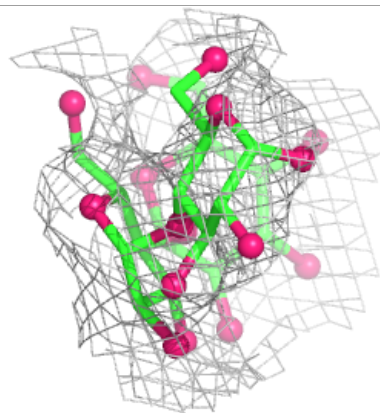
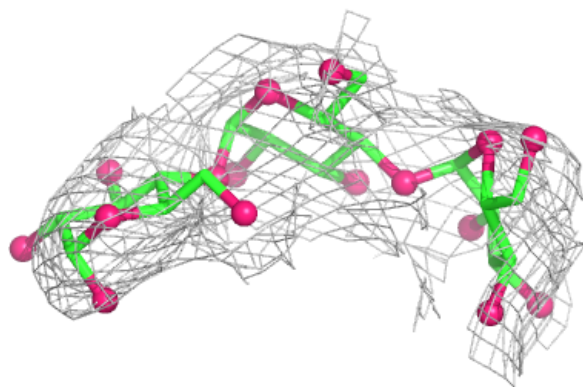
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

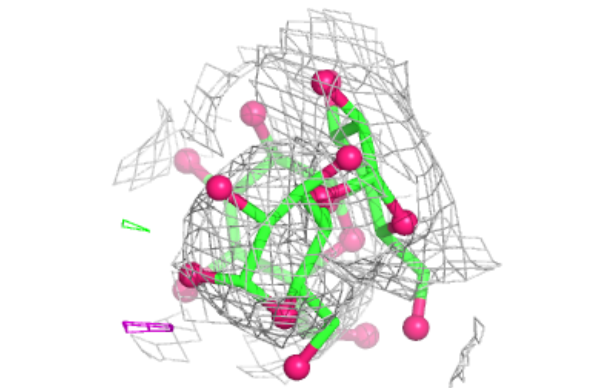
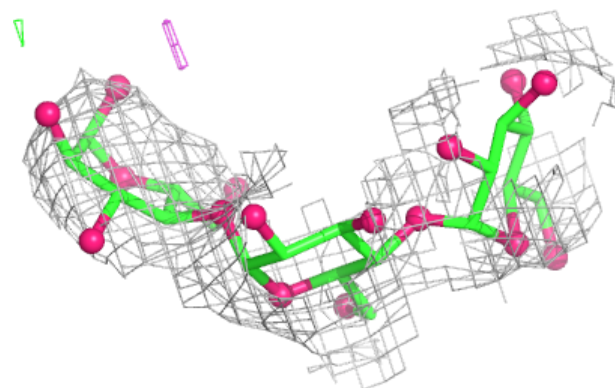
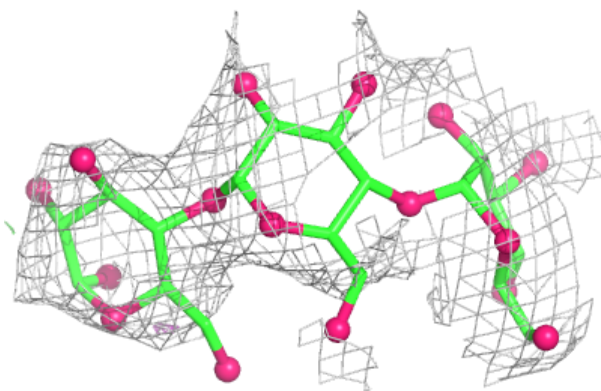
**Electron density around Chain F:**

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

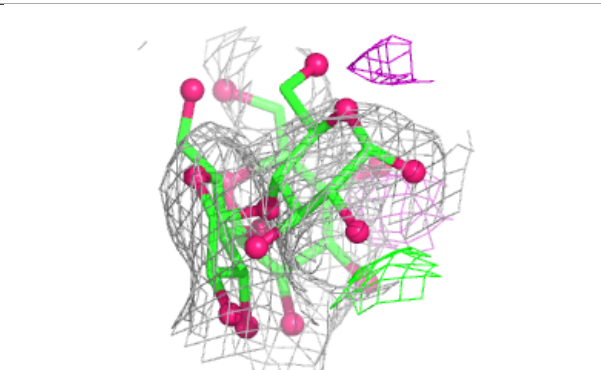
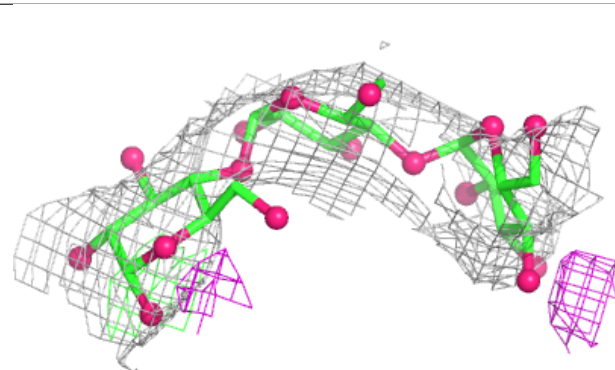
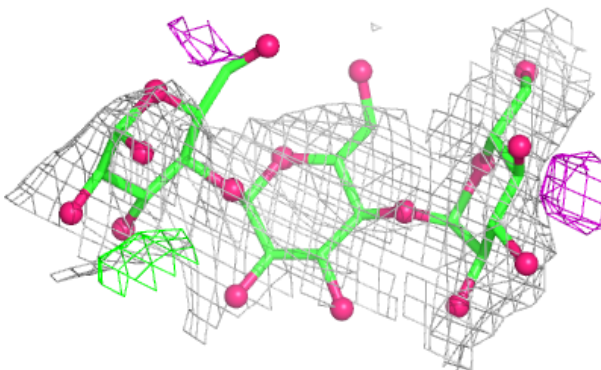


Electron density around Chain G:

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

**Electron density around Chain H:**

$2mF_o-DF_c$ (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.