



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 1, 2023 – 06:58 PM EDT

PDB ID : 3QHM
Title : Crystal analysis of the complex structure, E342A-cellobiohydrolase, of endocellulase from *Pyrococcus horikoshii*
Authors : Kim, H.-W.; Ishikawa, K.
Deposited on : 2011-01-26
Resolution : 2.01 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

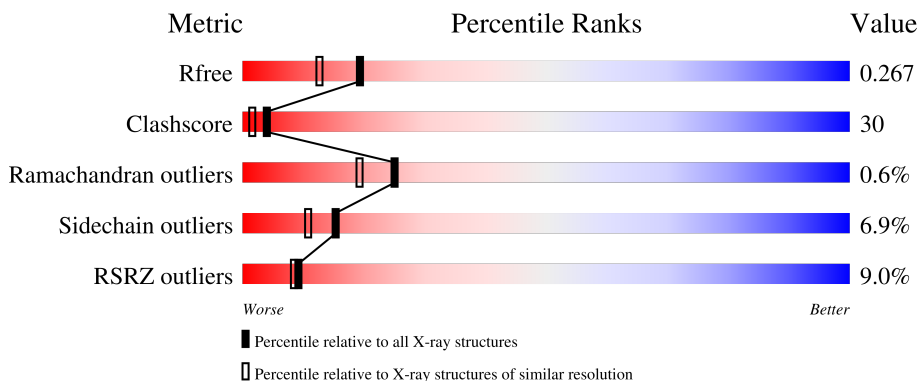
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	458	
1	B	458	
1	C	458	
2	D	4	
2	E	4	

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Mol	Chain	Length	Quality of chain
2	G	4	 75% 25%
3	F	2	 50% 50%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9565 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 458aa long hypothetical endo-1,4-beta-glucanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	377	3067	2000	502	555	10	0	0	0
1	B	377	3067	2000	502	555	10	0	0	0
1	C	377	3067	2000	502	555	10	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	342	ALA	GLU	engineered mutation	UNP O58925
B	342	ALA	GLU	engineered mutation	UNP O58925
C	342	ALA	GLU	engineered mutation	UNP O58925

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	D	4	45	24	21	0	0	0
2	E	4	45	24	21	0	0	0
2	G	4	45	24	21	0	0	0

- Molecule 3 is an oligosaccharide called beta-D-glucopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
3	F	2	23	12	11	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	67	Total	O	0	0
			67	67		
4	B	97	Total	O	0	0
			97	97		
4	C	42	Total	O	0	0
			42	42		

BGC1
BGC2
BGC3
BGC4

- Molecule 3: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain F: 

BGC1
BGC2

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	162.53Å 58.60Å 139.06Å 90.00° 109.69° 90.00°	Depositor
Resolution (Å)	35.25 – 2.01 35.25 – 2.01	Depositor EDS
% Data completeness (in resolution range)	87.7 (35.25-2.01) 87.5 (35.25-2.01)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 2.01Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.228 , 0.261 0.236 , 0.267	Depositor DCC
R_{free} test set	3729 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	28.7	Xtrriage
Anisotropy	0.723	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 63.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9565	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.39	0/3178	0.81	9/4339 (0.2%)
1	B	0.38	0/3178	0.69	1/4339 (0.0%)
1	C	0.32	0/3178	0.65	0/4339
All	All	0.37	0/9534	0.72	10/13017 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	347	TYR	N-CA-C	-8.56	87.89	111.00
1	A	393	ASP	N-CA-C	-8.50	88.06	111.00
1	A	130	LEU	CA-CB-CG	7.18	131.81	115.30
1	A	348	GLY	N-CA-C	6.17	128.53	113.10
1	A	130	LEU	N-CA-C	6.05	127.34	111.00
1	B	68	TRP	CA-CB-CG	5.86	124.83	113.70
1	A	347	TYR	CA-CB-CG	5.67	124.16	113.40
1	A	115	GLN	C-N-CD	5.49	139.94	128.40
1	A	389	ILE	N-CA-C	-5.37	96.51	111.00
1	A	197	ASP	N-CA-C	-5.08	97.28	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	347	TYR	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3067	0	2909	330	1
1	B	3067	0	2909	67	0
1	C	3067	0	2909	157	0
2	D	45	0	38	3	0
2	E	45	0	39	0	0
2	G	45	0	39	2	0
3	F	23	0	21	2	0
4	A	67	0	0	18	0
4	B	97	0	0	10	2
4	C	42	0	0	5	2
All	All	9565	0	8864	551	3

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 30.

All (551) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:PRO:HA	1:A:388:GLY:H	1.10	1.12
1:A:389:ILE:HD13	1:A:401:LYS:HE3	1.31	1.10
1:A:76:HIS:CD2	1:A:114:THR:HG21	1.90	1.05
1:A:402:TYR:HA	4:A:515:HOH:O	1.58	1.02
1:A:131:ASP:O	1:A:133:LEU:N	1.93	1.01
1:A:75:ASN:ND2	1:A:75:ASN:H	1.58	1.00
1:A:131:ASP:OD2	1:A:134:GLN:CG	2.11	0.99
1:A:112:PRO:O	1:A:131:ASP:HB3	1.62	0.99
1:A:131:ASP:OD2	1:A:134:GLN:HG2	1.62	0.99
1:A:87:GLU:HA	1:A:90:LEU:HD13	1.45	0.98
1:A:389:ILE:CD1	1:A:401:LYS:HE3	1.93	0.97
1:A:130:LEU:HB2	1:A:134:GLN:HG3	1.46	0.97
1:A:75:ASN:H	1:A:75:ASN:HD22	1.02	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:LEU:O	1:A:137:GLU:HG3	1.66	0.95
1:C:211:ALA:HA	1:C:216:THR:HG21	1.51	0.92
1:A:119:ILE:HG13	1:A:121:TYR:CE1	2.05	0.92
1:A:347:TYR:OH	1:A:360:ASN:CG	2.08	0.92
1:A:131:ASP:OD2	1:A:134:GLN:CD	2.10	0.91
1:A:390:LEU:HA	1:A:397:ILE:HA	1.53	0.90
1:B:380:ASN:H	1:B:380:ASN:HD22	1.18	0.89
1:A:131:ASP:C	1:A:133:LEU:H	1.76	0.88
1:A:88:ASP:O	1:A:92:GLN:HB2	1.74	0.88
1:A:381:PRO:HA	1:A:388:GLY:N	1.87	0.88
1:A:85:ASN:HD21	1:A:87:GLU:HG2	1.39	0.87
1:A:91:LEU:HA	4:A:513:HOH:O	1.72	0.87
1:A:77:VAL:HG23	1:A:116:PRO:HB3	1.56	0.86
1:A:68:TRP:CE2	1:A:78:VAL:HG11	2.11	0.86
1:A:75:ASN:HD22	1:A:75:ASN:N	1.75	0.85
1:A:146:LEU:HD11	4:A:473:HOH:O	1.76	0.85
1:C:201:GLU:OE2	1:C:297:HIS:HD2	1.58	0.85
1:C:318:ASN:OD1	1:C:319:LEU:HD12	1.75	0.85
1:C:269:TYR:C	1:C:270:ASN:HD22	1.80	0.83
1:A:224:ASN:HD21	1:B:75:ASN:HD21	1.27	0.83
1:A:78:VAL:HG12	4:A:502:HOH:O	1.80	0.82
1:C:198:LEU:HD21	1:C:249:ILE:CG2	2.10	0.82
1:B:146:LEU:HG	4:B:532:HOH:O	1.77	0.82
1:A:389:ILE:CG2	1:A:389:ILE:O	2.27	0.82
1:A:81:LEU:O	1:A:81:LEU:HD23	1.79	0.81
1:C:198:LEU:HD21	1:C:249:ILE:HG23	1.60	0.81
1:A:68:TRP:CZ2	1:A:78:VAL:HG11	2.15	0.81
1:A:364:ASP:O	1:A:368:GLU:HG3	1.81	0.81
1:A:131:ASP:O	1:A:134:GLN:N	2.14	0.80
1:B:384:GLY:HA3	4:B:524:HOH:O	1.81	0.80
1:C:85:ASN:HD22	1:C:88:ASP:H	1.29	0.80
1:C:214:ASP:OD1	1:C:216:THR:HB	1.82	0.80
1:C:254:THR:O	1:C:276:ASN:HB3	1.81	0.79
1:A:266:LYS:HE2	1:A:267:TRP:NE1	1.98	0.79
1:C:205:VAL:HG22	1:C:216:THR:O	1.83	0.78
1:C:270:ASN:HD22	1:C:270:ASN:N	1.80	0.78
1:A:85:ASN:HA	1:A:124:ASN:HD21	1.50	0.77
1:A:315:PHE:CE2	1:A:354:ARG:HG2	2.20	0.77
1:A:393:ASP:OD2	1:A:396:THR:HG22	1.85	0.77
1:C:64:PHE:HB3	1:C:409:MET:HE3	1.66	0.76
1:A:290:ASN:H	1:A:290:ASN:HD22	1.31	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:GLY:C	1:A:130:LEU:HD22	2.06	0.76
1:C:35:THR:HG22	1:C:38:GLY:H	1.49	0.76
1:C:35:THR:HG22	1:C:37:THR:H	1.51	0.75
1:A:381:PRO:HD2	1:A:394:TRP:CH2	2.20	0.75
1:A:347:TYR:CD1	1:A:356:VAL:HA	2.22	0.74
1:A:389:ILE:O	1:A:389:ILE:HG22	1.87	0.74
1:C:124:ASN:HB3	1:C:127:LEU:HD12	1.69	0.74
1:A:71:PHE:HB2	1:A:104:PRO:O	1.87	0.74
1:A:81:LEU:HD22	1:A:119:ILE:HB	1.70	0.74
1:A:190:TYR:N	1:A:190:TYR:HD2	1.85	0.74
1:A:347:TYR:OH	1:A:360:ASN:OD1	2.04	0.74
1:A:214:ASP:HB3	4:B:524:HOH:O	1.86	0.74
1:A:37:THR:HG22	1:A:39:ILE:H	1.53	0.74
1:A:75:ASN:HA	1:A:117:ILE:HB	1.70	0.74
1:B:253:GLY:O	1:B:274:GLY:HA2	1.87	0.74
1:A:381:PRO:HD2	1:A:394:TRP:CZ2	2.24	0.73
1:A:73:THR:HB	1:A:74:PRO:HD2	1.69	0.73
1:A:71:PHE:HZ	1:A:135:ILE:HD11	1.52	0.73
1:A:89:MET:O	1:A:93:ILE:HG12	1.89	0.72
1:A:190:TYR:N	1:A:190:TYR:CD2	2.56	0.72
1:A:406:LYS:HA	1:A:409:MET:CE	2.19	0.72
1:C:306:GLN:OE1	1:C:307:PRO:HD2	1.88	0.72
1:A:144:GLY:HA3	4:A:501:HOH:O	1.88	0.72
1:A:112:PRO:O	1:A:131:ASP:CB	2.37	0.72
1:C:35:THR:HG21	1:C:39:ILE:O	1.90	0.72
1:A:75:ASN:ND2	1:A:75:ASN:N	2.34	0.71
1:A:81:LEU:HB2	4:A:502:HOH:O	1.89	0.71
1:C:214:ASP:C	1:C:216:THR:H	1.93	0.71
1:A:76:HIS:O	1:A:132:SER:HB3	1.89	0.71
1:A:76:HIS:CD2	1:A:114:THR:CG2	2.71	0.71
1:C:214:ASP:O	1:C:216:THR:N	2.19	0.71
1:B:380:ASN:HD22	1:B:380:ASN:N	1.89	0.71
1:B:241:LEU:HD13	1:B:249:ILE:HD11	1.74	0.70
1:C:35:THR:HG22	1:C:37:THR:N	2.06	0.70
1:C:35:THR:CG2	1:C:38:GLY:H	2.05	0.70
1:A:349:HIS:C	1:A:351:GLY:N	2.42	0.70
1:A:407:ARG:O	1:A:410:ASP:HB3	1.92	0.69
1:A:347:TYR:CD2	1:A:356:VAL:HG13	2.28	0.69
1:C:201:GLU:OE2	1:C:297:HIS:CD2	2.44	0.69
1:A:86:TRP:O	1:A:89:MET:HB2	1.92	0.68
1:A:345:GLY:O	1:A:346:LYS:HG3	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:LYS:HA	1:A:409:MET:HE3	1.76	0.68
1:A:390:LEU:HD23	1:A:391:GLN:H	1.57	0.68
1:A:349:HIS:O	1:A:351:GLY:N	2.26	0.68
1:A:131:ASP:C	1:A:133:LEU:N	2.38	0.67
1:A:86:TRP:O	1:A:90:LEU:HD12	1.94	0.67
1:A:87:GLU:HA	1:A:90:LEU:CD1	2.23	0.67
1:C:391:GLN:HG2	1:C:398:TRP:CE2	2.30	0.67
1:C:330:VAL:HG12	1:C:336:TYR:HB2	1.75	0.67
1:A:68:TRP:HE3	1:A:380:ASN:HD21	1.41	0.66
1:A:199:LYS:HE2	1:A:200:ASN:O	1.94	0.66
1:A:405:LEU:HB3	4:A:515:HOH:O	1.95	0.66
1:A:68:TRP:CG	1:A:78:VAL:HG21	2.30	0.66
1:A:390:LEU:O	1:A:391:GLN:HG3	1.96	0.66
1:C:316:PRO:O	1:C:319:LEU:HD13	1.95	0.66
1:A:84:ARG:NH1	1:A:394:TRP:O	2.28	0.66
1:A:349:HIS:C	1:A:351:GLY:H	1.97	0.66
1:A:91:LEU:HD23	4:A:473:HOH:O	1.94	0.66
1:C:289:ARG:NH1	1:C:336:TYR:CE1	2.64	0.66
1:A:346:LYS:HA	1:A:401:LYS:HG3	1.78	0.66
1:B:201:GLU:OE2	1:B:297:HIS:HD2	1.79	0.66
1:B:403:ASN:HA	1:B:406:LYS:CD	2.26	0.65
1:B:85:ASN:HD22	1:B:88:ASP:H	1.44	0.65
1:A:397:ILE:HG22	4:A:478:HOH:O	1.97	0.65
1:A:93:ILE:HD11	1:A:379:TRP:CD2	2.32	0.65
1:A:71:PHE:CZ	1:A:135:ILE:HD11	2.31	0.64
1:B:246:HIS:CD2	1:B:246:HIS:H	2.15	0.64
1:B:364:ASP:O	1:B:368:GLU:HG3	1.97	0.64
1:A:81:LEU:HD22	1:A:119:ILE:CG2	2.27	0.64
1:A:402:TYR:OH	1:A:406:LYS:NZ	2.31	0.64
1:A:89:MET:HE1	1:A:394:TRP:HE3	1.63	0.64
1:A:84:ARG:HD3	1:A:89:MET:CE	2.28	0.64
1:C:354:ARG:HA	1:C:357:ILE:HD12	1.80	0.64
1:C:380:ASN:HD22	1:C:380:ASN:H	1.46	0.64
1:A:85:ASN:ND2	1:A:87:GLU:HG2	2.13	0.63
1:A:366:MET:HB2	1:A:408:LEU:HD11	1.81	0.63
1:B:403:ASN:HA	1:B:406:LYS:HD3	1.80	0.63
1:C:208:PRO:HA	1:C:211:ALA:H	1.64	0.63
1:A:364:ASP:OD1	1:A:407:ARG:NH2	2.27	0.62
1:A:119:ILE:HG13	1:A:121:TYR:HE1	1.62	0.62
1:C:198:LEU:HD22	1:C:198:LEU:H	1.65	0.62
1:A:85:ASN:HA	1:A:124:ASN:ND2	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:LYS:HE3	1:A:386:THR:C	2.20	0.62
1:A:77:VAL:HG12	1:A:78:VAL:O	1.99	0.62
1:A:246:HIS:H	1:A:246:HIS:CD2	2.16	0.61
1:A:347:TYR:CE1	1:A:356:VAL:O	2.53	0.61
1:C:306:GLN:HB2	1:C:309:PHE:CE2	2.35	0.61
1:C:246:HIS:H	1:C:246:HIS:CD2	2.17	0.61
1:A:62:HIS:CD2	1:A:372:CYS:SG	2.94	0.61
1:C:49:TYR:CE1	1:C:60:PRO:HG3	2.35	0.61
1:A:81:LEU:HD22	1:A:119:ILE:CB	2.30	0.61
1:B:142:LYS:HE2	1:B:146:LEU:HD11	1.83	0.61
1:C:316:PRO:HG3	1:C:357:ILE:CG2	2.31	0.61
1:A:160:THR:HB	1:A:161:HIS:HD2	1.66	0.60
1:C:39:ILE:HD13	1:C:191:TRP:HB2	1.83	0.60
1:C:76:HIS:NE2	1:C:159:CYS:SG	2.74	0.60
1:A:72:GLU:O	1:A:72:GLU:HG2	2.02	0.60
1:A:89:MET:HE3	1:A:394:TRP:CZ3	2.37	0.60
1:A:392:ASP:O	1:A:394:TRP:N	2.35	0.60
1:C:276:ASN:OD1	1:C:278:MET:HG3	2.01	0.60
1:A:344:GLY:HA3	1:A:377:TRP:O	2.02	0.60
1:A:399:GLU:CD	1:A:399:GLU:H	2.03	0.60
1:B:297:HIS:HE1	4:B:484:HOH:O	1.84	0.60
1:C:137:GLU:OE2	1:C:186:ARG:HD3	2.01	0.60
1:A:80:GLY:HA3	1:A:394:TRP:CE3	2.37	0.59
1:A:55:SER:OG	1:A:57:GLU:HG3	2.02	0.59
1:B:79:HIS:HD2	1:B:382:ASP:OD2	1.85	0.59
1:A:93:ILE:HD13	1:A:379:TRP:CE3	2.38	0.59
1:A:133:LEU:O	1:A:136:MET:HB3	2.03	0.59
1:C:219:THR:HG23	1:C:228:ASP:OD1	2.02	0.59
1:C:306:GLN:HB2	1:C:309:PHE:CD2	2.37	0.59
1:A:383:SER:O	1:A:385:ASP:N	2.36	0.59
1:A:89:MET:HE1	1:A:394:TRP:CE3	2.37	0.59
1:C:271:ALA:CB	1:C:276:ASN:HB2	2.33	0.59
1:A:156:ARG:HB3	1:A:158:GLY:O	2.03	0.58
1:C:163:GLU:HB3	4:C:462:HOH:O	2.01	0.58
1:A:84:ARG:HD3	1:A:89:MET:SD	2.43	0.58
1:A:86:TRP:H	1:A:124:ASN:HD21	1.51	0.58
1:B:96:LEU:O	1:B:409:MET:HE1	2.03	0.58
1:C:319:LEU:HD12	1:C:319:LEU:H	1.67	0.58
1:A:77:VAL:HG23	1:A:116:PRO:CB	2.30	0.58
1:A:84:ARG:HD3	1:A:89:MET:HE1	1.85	0.58
1:C:34:GLN:HG3	1:C:53:VAL:HG21	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:LEU:O	1:A:93:ILE:HB	2.02	0.58
1:A:121:TYR:O	1:A:125:PRO:HG3	2.03	0.58
1:A:379:TRP:CE3	1:A:389:ILE:HG22	2.39	0.58
1:C:315:PHE:CE1	1:C:316:PRO:HB3	2.39	0.58
1:A:352:ASP:OD2	1:A:353:PRO:N	2.36	0.58
1:A:391:GLN:O	1:A:393:ASP:O	2.21	0.58
1:C:180:TRP:O	1:C:183:VAL:HG22	2.03	0.58
1:A:96:LEU:HD22	1:A:402:TYR:CZ	2.39	0.57
1:A:135:ILE:HG13	1:A:136:MET:N	2.19	0.57
1:A:68:TRP:CE3	1:A:68:TRP:HA	2.39	0.57
1:B:142:LYS:HE2	1:B:146:LEU:CD1	2.34	0.57
1:C:138:LYS:HD2	4:C:471:HOH:O	2.03	0.57
1:C:407:ARG:HG2	1:C:407:ARG:HH11	1.69	0.57
1:C:120:ASP:OD1	1:C:122:SER:HB3	2.04	0.57
1:C:347:TYR:HB2	1:C:401:LYS:HD3	1.86	0.57
1:A:67:ASN:OD1	1:A:102:ARG:HB3	2.05	0.57
1:A:352:ASP:OD2	1:A:354:ARG:N	2.38	0.57
1:A:380:ASN:C	1:A:388:GLY:HA3	2.24	0.57
1:A:400:ASP:OD1	1:A:400:ASP:N	2.38	0.57
1:A:391:GLN:OE1	1:A:396:THR:HG23	2.05	0.57
1:A:392:ASP:O	1:A:393:ASP:C	2.41	0.57
1:B:85:ASN:HD21	1:B:87:GLU:HB3	1.68	0.57
1:A:96:LEU:O	1:A:406:LYS:HE2	2.05	0.57
1:C:219:THR:HG22	1:C:227:THR:C	2.25	0.57
1:A:66:VAL:CG2	1:A:101:ILE:HG22	2.34	0.57
1:C:324:TYR:HA	1:C:328:GLY:HA3	1.87	0.57
1:C:108:GLU:HG3	4:C:493:HOH:O	2.03	0.56
1:A:89:MET:CE	1:A:394:TRP:CE3	2.88	0.56
1:C:380:ASN:HD22	1:C:380:ASN:N	2.03	0.56
1:A:68:TRP:HE3	1:A:68:TRP:HA	1.71	0.56
1:A:376:TYR:HB3	4:A:517:HOH:O	2.06	0.56
1:C:35:THR:CG2	1:C:37:THR:H	2.18	0.56
1:C:180:TRP:CH2	1:C:196:ALA:HB1	2.41	0.56
1:A:72:GLU:HG3	1:A:155:HIS:CD2	2.40	0.56
1:C:297:HIS:HE1	4:C:474:HOH:O	1.89	0.56
1:A:347:TYR:OH	1:A:360:ASN:ND2	2.38	0.56
1:B:403:ASN:O	1:B:406:LYS:HG2	2.06	0.56
1:A:65:GLY:HA2	1:A:100:ALA:O	2.06	0.56
1:A:267:TRP:CZ3	1:A:307:PRO:HG2	2.41	0.56
1:A:66:VAL:HG21	1:A:93:ILE:HD12	1.86	0.56
1:A:359:GLN:NE2	1:A:376:TYR:OH	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:TRP:CE2	1:C:230:ASN:HB3	2.41	0.56
1:C:209:PRO:HA	1:C:212:TYR:CD2	2.41	0.55
1:A:352:ASP:OD2	1:A:352:ASP:C	2.43	0.55
1:A:80:GLY:HA3	1:A:394:TRP:CD2	2.42	0.55
1:A:66:VAL:HG23	1:A:101:ILE:HG22	1.89	0.55
1:C:270:ASN:N	1:C:270:ASN:ND2	2.53	0.55
1:A:390:LEU:CD2	1:A:391:GLN:H	2.20	0.55
1:A:94:LYS:HZ2	1:A:94:LYS:HB2	1.72	0.55
1:A:93:ILE:CD1	1:A:379:TRP:CD2	2.90	0.54
1:A:404:ASN:O	1:A:407:ARG:HD3	2.06	0.54
1:C:201:GLU:O	1:C:203:HIS:HD2	1.90	0.54
1:B:406:LYS:HB2	1:B:406:LYS:NZ	2.23	0.54
1:A:62:HIS:HD2	1:A:372:CYS:SG	2.30	0.54
1:A:102:ARG:O	1:A:104:PRO:HD3	2.08	0.54
1:A:297:HIS:HE1	4:A:489:HOH:O	1.89	0.54
1:A:406:LYS:HA	1:A:409:MET:HE2	1.89	0.54
1:A:94:LYS:NZ	1:A:146:LEU:HB3	2.23	0.53
1:A:180:TRP:CH2	1:A:196:ALA:HB1	2.43	0.53
1:C:219:THR:OG1	1:C:222:MET:HG2	2.08	0.53
1:C:400:ASP:OD1	1:C:401:LYS:N	2.41	0.53
1:C:404:ASN:O	1:C:407:ARG:HB2	2.08	0.53
1:A:76:HIS:HB3	1:A:132:SER:OG	2.08	0.53
1:A:117:ILE:CG2	1:A:118:GLY:N	2.72	0.53
1:C:409:MET:O	1:C:410:ASP:HB2	2.07	0.53
1:C:364:ASP:O	1:C:368:GLU:HG3	2.08	0.53
1:B:246:HIS:H	1:B:246:HIS:HD2	1.56	0.53
1:A:380:ASN:O	1:A:388:GLY:N	2.42	0.53
1:A:64:PHE:CZ	1:A:409:MET:HA	2.43	0.53
1:A:131:ASP:O	1:A:134:GLN:HG2	2.08	0.53
1:A:346:LYS:HE3	1:A:386:THR:CA	2.39	0.53
1:A:289:ARG:HB2	1:C:145:ASP:HA	1.91	0.52
1:A:161:HIS:HA	4:A:508:HOH:O	2.08	0.52
1:C:35:THR:OG1	1:C:246:HIS:HB3	2.09	0.52
1:A:94:LYS:HD3	4:A:499:HOH:O	2.08	0.52
1:A:135:ILE:C	1:A:135:ILE:HD12	2.30	0.52
1:C:219:THR:CG2	1:C:227:THR:OG1	2.57	0.52
1:C:220:TRP:HB3	1:C:257:THR:HB	1.90	0.52
1:C:246:HIS:H	1:C:246:HIS:HD2	1.57	0.52
1:A:347:TYR:CA	1:A:356:VAL:HG22	2.39	0.52
1:A:367:ILE:HG13	1:A:408:LEU:HG	1.91	0.52
1:C:37:THR:HG22	1:C:39:ILE:HD12	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:LEU:HD22	1:A:130:LEU:N	2.24	0.52
1:A:86:TRP:HD1	1:A:124:ASN:ND2	2.08	0.51
1:A:94:LYS:HE2	4:A:488:HOH:O	2.10	0.51
1:C:214:ASP:C	1:C:216:THR:N	2.59	0.51
1:A:381:PRO:O	1:A:387:GLY:HA2	2.10	0.51
1:B:265:TYR:CE2	1:B:267:TRP:HB2	2.45	0.51
1:A:90:LEU:HD21	1:A:142:LYS:HB3	1.91	0.51
1:A:68:TRP:CZ2	1:A:78:VAL:CG1	2.89	0.51
1:A:87:GLU:O	1:A:91:LEU:N	2.37	0.51
1:A:379:TRP:CE3	1:A:389:ILE:CG2	2.94	0.51
1:B:357:ILE:HD12	1:B:357:ILE:H	1.75	0.51
1:A:81:LEU:CD2	1:A:119:ILE:HB	2.39	0.51
1:C:198:LEU:HD22	1:C:250:PHE:O	2.09	0.51
1:A:347:TYR:CB	1:A:356:VAL:HG22	2.41	0.51
1:A:389:ILE:O	1:A:389:ILE:HG23	2.09	0.51
1:C:156:ARG:HG2	1:C:162:ILE:HG22	1.91	0.51
1:A:93:ILE:HD11	1:A:379:TRP:CG	2.45	0.51
1:A:73:THR:HB	1:A:75:ASN:ND2	2.25	0.51
1:A:380:ASN:O	1:A:388:GLY:CA	2.59	0.51
1:C:35:THR:HG21	1:C:39:ILE:H	1.75	0.51
1:C:131:ASP:OD1	1:C:134:GLN:HG3	2.11	0.51
1:A:130:LEU:HB3	1:A:134:GLN:HE21	1.76	0.51
1:A:77:VAL:CG1	1:A:78:VAL:N	2.74	0.50
1:B:320:PRO:HD3	1:B:361:LYS:HZ3	1.76	0.50
1:C:110:VAL:HA	1:C:133:LEU:HD13	1.92	0.50
1:C:85:ASN:HD21	1:C:87:GLU:HB2	1.75	0.50
1:A:94:LYS:HG3	1:A:95:SER:N	2.25	0.50
1:C:208:PRO:HB2	1:C:212:TYR:CE2	2.46	0.50
1:C:219:THR:HG22	1:C:227:THR:O	2.12	0.50
1:A:201:GLU:O	1:A:203:HIS:HD2	1.95	0.50
1:C:201:GLU:O	1:C:203:HIS:CD2	2.65	0.50
1:A:152:LEU:HD11	1:A:187:PHE:HB3	1.93	0.50
1:A:405:LEU:CB	4:A:515:HOH:O	2.56	0.50
1:C:316:PRO:HG3	1:C:357:ILE:HG21	1.94	0.50
1:C:377:TRP:CZ2	2:G:1:BGC:H3	2.47	0.50
1:A:84:ARG:CZ	1:A:394:TRP:O	2.59	0.49
1:A:156:ARG:CB	1:A:158:GLY:O	2.59	0.49
1:B:200:ASN:ND2	1:B:201:GLU:HG3	2.27	0.49
1:C:220:TRP:NE1	1:C:230:ASN:HB3	2.26	0.49
1:B:181:ILE:O	1:B:185:LYS:HG3	2.12	0.49
1:A:389:ILE:O	1:A:390:LEU:HB2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:THR:OG1	1:A:75:ASN:ND2	2.45	0.49
1:A:347:TYR:CE2	1:A:356:VAL:HG13	2.48	0.49
1:A:87:GLU:O	1:A:91:LEU:HG	2.13	0.49
1:A:253:GLY:O	1:A:274:GLY:HA2	2.13	0.49
1:A:397:ILE:HD13	1:A:397:ILE:O	2.13	0.49
1:B:177:ILE:O	1:B:181:ILE:HG13	2.12	0.49
1:C:85:ASN:ND2	1:C:88:ASP:H	2.05	0.49
1:C:219:THR:HG23	1:C:228:ASP:HA	1.94	0.49
1:B:68:TRP:HZ2	4:B:549:HOH:O	1.95	0.49
1:A:68:TRP:HB3	1:A:103:LEU:CD2	2.42	0.49
1:A:212:TYR:CD2	1:A:259:PRO:HD3	2.46	0.49
1:A:389:ILE:O	1:A:390:LEU:CB	2.57	0.49
1:A:393:ASP:O	1:A:395:THR:N	2.42	0.49
1:B:68:TRP:CZ3	1:B:79:HIS:O	2.66	0.49
1:A:266:LYS:HE2	1:A:267:TRP:HE1	1.77	0.49
1:A:135:ILE:HD12	1:A:139:ILE:HG13	1.95	0.49
1:B:203:HIS:HE1	4:B:512:HOH:O	1.96	0.49
1:B:259:PRO:O	1:B:263:SER:OG	2.26	0.49
1:B:380:ASN:HB3	1:B:382:ASP:OD1	2.13	0.49
1:A:90:LEU:HD22	1:A:142:LYS:CG	2.43	0.48
1:B:201:GLU:O	1:B:203:HIS:HD2	1.95	0.48
1:C:71:PHE:CE1	1:C:77:VAL:HA	2.48	0.48
1:A:94:LYS:HB2	1:A:94:LYS:NZ	2.28	0.48
1:A:380:ASN:C	1:A:388:GLY:CA	2.82	0.48
1:B:200:ASN:CG	1:B:201:GLU:HG3	2.34	0.48
1:C:270:ASN:OD1	1:C:307:PRO:HG3	2.14	0.48
1:A:85:ASN:CA	1:A:124:ASN:HD21	2.24	0.48
1:A:130:LEU:N	1:A:130:LEU:CD2	2.77	0.48
1:A:155:HIS:HA	1:A:200:ASN:CB	2.44	0.48
1:A:234:GLU:OE1	1:A:286:ASN:HB3	2.14	0.48
1:A:77:VAL:HG12	1:A:78:VAL:N	2.28	0.48
1:A:331:LYS:HB2	1:A:338:VAL:CG2	2.44	0.48
1:A:73:THR:HB	1:A:75:ASN:HD21	1.77	0.47
1:C:187:PHE:HA	1:C:190:TYR:HD2	1.79	0.47
1:A:391:GLN:HB2	1:A:396:THR:HG23	1.95	0.47
1:C:332:LEU:HD21	1:C:369:ASN:HB3	1.97	0.47
1:A:376:TYR:CD2	1:A:389:ILE:HG12	2.50	0.47
1:B:86:TRP:CH2	1:B:142:LYS:HG3	2.50	0.47
1:C:166:TRP:NE1	1:C:173:GLU:OE1	2.43	0.47
1:A:69:PHE:CZ	1:A:382:ASP:OD1	2.67	0.47
1:A:73:THR:CB	1:A:75:ASN:ND2	2.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:TYR:HA	1:A:356:VAL:HA	1.96	0.47
1:A:385:ASP:OD1	2:D:3:BGC:H5	2.14	0.47
1:A:68:TRP:CE3	1:A:380:ASN:ND2	2.81	0.47
1:A:69:PHE:O	1:A:78:VAL:HA	2.13	0.47
1:A:132:SER:O	1:A:136:MET:HB2	2.15	0.47
1:A:155:HIS:HA	1:A:200:ASN:HB3	1.97	0.47
1:B:107:THR:O	1:B:111:LYS:HG2	2.15	0.47
1:A:76:HIS:HB2	1:A:116:PRO:HA	1.96	0.47
1:A:366:MET:CB	1:A:408:LEU:HD11	2.43	0.47
1:B:75:ASN:ND2	4:B:516:HOH:O	2.47	0.47
1:C:381:PRO:HD2	1:C:394:TRP:CH2	2.49	0.47
1:A:49:TYR:OH	1:C:54:THR:HG21	2.15	0.47
1:A:69:PHE:CD2	1:A:377:TRP:CZ2	3.03	0.47
1:A:161:HIS:N	1:A:161:HIS:CD2	2.82	0.47
1:A:346:LYS:HE3	1:A:386:THR:HA	1.95	0.47
1:B:222:MET:SD	1:B:259:PRO:HD3	2.55	0.47
1:C:406:LYS:O	1:C:410:ASP:N	2.47	0.47
1:A:267:TRP:CE3	1:A:308:TYR:HD2	2.32	0.47
1:C:340:ILE:HD12	1:C:366:MET:HE3	1.96	0.47
1:A:347:TYR:CG	1:A:356:VAL:HA	2.50	0.46
1:A:402:TYR:OH	1:A:406:LYS:CE	2.64	0.46
1:B:73:THR:OG1	1:B:79:HIS:HE1	1.98	0.46
1:A:128:ARG:C	1:A:130:LEU:HD22	2.36	0.46
1:A:42:GLU:OE2	1:A:44:ARG:NE	2.43	0.46
1:A:302:ASP:HB3	1:A:385:ASP:O	2.15	0.46
1:A:349:HIS:O	1:A:350:GLY:C	2.54	0.46
1:B:79:HIS:CD2	1:B:382:ASP:OD2	2.67	0.46
1:B:316:PRO:HG3	1:B:357:ILE:CG2	2.45	0.46
1:C:295:SER:HA	1:C:339:VAL:O	2.15	0.46
1:C:380:ASN:O	1:C:388:GLY:HA3	2.15	0.46
1:C:69:PHE:HB2	1:C:72:GLU:OE2	2.16	0.46
1:C:315:PHE:HD1	1:C:319:LEU:HD11	1.79	0.46
1:A:69:PHE:HZ	1:A:382:ASP:OD1	1.99	0.46
1:B:409:MET:O	1:B:410:ASP:OD1	2.33	0.46
1:A:91:LEU:N	4:A:473:HOH:O	2.47	0.46
1:A:135:ILE:O	1:A:139:ILE:HG12	2.15	0.46
1:A:352:ASP:HA	1:A:353:PRO:HD3	1.28	0.46
1:B:380:ASN:N	1:B:380:ASN:ND2	2.60	0.46
1:A:72:GLU:HG3	1:A:155:HIS:HD2	1.81	0.46
1:A:80:GLY:HA3	1:A:394:TRP:CZ3	2.51	0.46
1:A:316:PRO:O	1:A:319:LEU:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:315:PHE:HA	1:C:316:PRO:HA	1.70	0.46
1:A:37:THR:HG22	1:A:39:ILE:N	2.27	0.46
1:A:71:PHE:CD2	1:A:103:LEU:HD22	2.51	0.46
1:C:246:HIS:CD2	1:C:246:HIS:N	2.84	0.46
1:A:131:ASP:OD2	1:A:134:GLN:NE2	2.49	0.46
1:A:135:ILE:HD12	1:A:139:ILE:CG1	2.46	0.45
1:C:68:TRP:HD1	1:C:380:ASN:HD21	1.63	0.45
1:C:163:GLU:CD	1:C:163:GLU:H	2.19	0.45
1:A:69:PHE:CD1	1:A:69:PHE:C	2.89	0.45
1:A:331:LYS:HB2	1:A:338:VAL:HG23	1.99	0.45
1:B:287:LEU:HD12	1:B:292:LEU:HD13	1.98	0.45
1:C:281:LYS:NZ	1:C:281:LYS:CB	2.79	0.45
1:C:209:PRO:HA	1:C:212:TYR:HD2	1.82	0.45
1:A:93:ILE:CD1	1:A:379:TRP:CE3	2.99	0.45
1:A:123:LYS:O	1:A:125:PRO:HD3	2.16	0.45
1:B:75:ASN:ND2	1:B:79:HIS:CE1	2.84	0.45
1:B:299:TYR:HE2	4:B:529:HOH:O	1.99	0.45
1:A:291:LYS:HA	1:A:291:LYS:HE2	1.97	0.45
1:C:74:PRO:HD3	1:C:159:CYS:O	2.17	0.45
1:A:103:LEU:HA	1:A:103:LEU:HD23	1.81	0.45
1:C:35:THR:HG21	1:C:39:ILE:N	2.32	0.45
1:C:304:TYR:CD2	3:F:1:BGC:H5	2.51	0.45
1:B:85:ASN:ND2	1:B:87:GLU:HB3	2.31	0.45
1:C:219:THR:HG21	1:C:227:THR:OG1	2.16	0.45
1:C:331:LYS:HB2	1:C:338:VAL:HG23	1.99	0.45
1:A:68:TRP:HH2	1:A:89:MET:HE3	1.82	0.45
1:A:100:ALA:HA	1:A:148:ILE:HG23	1.98	0.45
1:B:57:GLU:HG3	1:B:58:GLU:N	2.32	0.45
1:A:94:LYS:HE3	1:A:98:PHE:O	2.17	0.45
1:A:143:ALA:O	1:A:148:ILE:HB	2.17	0.45
1:A:220:TRP:CE2	1:A:230:ASN:HB3	2.52	0.44
1:C:319:LEU:HD12	1:C:319:LEU:N	2.30	0.44
1:A:77:VAL:HG22	1:A:135:ILE:HG12	1.99	0.44
1:C:162:ILE:O	1:C:162:ILE:HG13	2.16	0.44
1:C:292:LEU:HG	1:C:293:VAL:N	2.31	0.44
1:B:357:ILE:HD12	1:B:357:ILE:N	2.32	0.44
1:A:64:PHE:CG	1:A:409:MET:HG2	2.53	0.44
1:A:86:TRP:N	1:A:124:ASN:HD21	2.14	0.44
1:A:134:GLN:HG2	1:A:134:GLN:H	1.55	0.44
1:A:306:GLN:O	1:A:309:PHE:HB2	2.16	0.44
1:A:383:SER:HB3	1:A:387:GLY:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:406:LYS:HB2	1:B:406:LYS:HZ2	1.81	0.44
1:A:36:PRO:HG2	1:A:37:THR:H	1.83	0.44
1:B:380:ASN:H	1:B:380:ASN:ND2	1.97	0.44
1:A:65:GLY:CA	1:A:100:ALA:O	2.66	0.44
1:C:330:VAL:HG12	1:C:336:TYR:CB	2.45	0.44
1:A:144:GLY:CA	4:A:501:HOH:O	2.56	0.44
1:A:152:LEU:HD11	1:A:187:PHE:CB	2.48	0.44
1:C:219:THR:HG22	1:C:227:THR:OG1	2.17	0.44
1:A:153:ASP:OD1	1:A:153:ASP:C	2.55	0.44
1:A:267:TRP:CZ2	1:A:313:LYS:HG2	2.52	0.44
1:C:316:PRO:HA	1:C:319:LEU:CD1	2.48	0.44
1:A:295:SER:HA	1:A:339:VAL:O	2.18	0.44
1:A:86:TRP:H	1:A:124:ASN:ND2	2.14	0.43
1:A:134:GLN:O	1:A:135:ILE:C	2.54	0.43
1:A:397:ILE:HG12	1:A:398:TRP:N	2.32	0.43
1:C:316:PRO:CA	1:C:319:LEU:HD13	2.48	0.43
1:A:272:TRP:CZ3	1:A:304:TYR:HB3	2.52	0.43
1:A:381:PRO:HD2	1:A:394:TRP:CZ3	2.53	0.43
1:B:238:LYS:NZ	4:B:514:HOH:O	2.48	0.43
1:C:188:GLY:O	1:C:246:HIS:HE1	2.00	0.43
1:A:121:TYR:N	1:A:121:TYR:CD1	2.86	0.43
1:A:123:LYS:C	1:A:125:PRO:HD3	2.38	0.43
1:C:35:THR:CG2	1:C:39:ILE:O	2.60	0.43
1:C:262:ASP:HA	1:C:278:MET:HE3	2.00	0.43
1:A:68:TRP:CD2	1:A:78:VAL:HG21	2.54	0.43
1:B:406:LYS:NZ	1:B:406:LYS:CB	2.82	0.43
1:A:36:PRO:HG2	1:A:37:THR:N	2.34	0.43
1:A:383:SER:C	1:A:385:ASP:N	2.70	0.43
1:A:389:ILE:HD13	1:A:389:ILE:HA	1.57	0.43
1:C:37:THR:HG21	1:C:191:TRP:CB	2.49	0.43
1:A:346:LYS:HG2	4:A:492:HOH:O	2.18	0.43
1:A:354:ARG:HG3	1:A:357:ILE:HD12	2.01	0.43
1:C:286:ASN:ND2	4:C:464:HOH:O	2.42	0.43
1:A:164:PRO:HG3	1:A:205:VAL:CG1	2.49	0.43
1:A:289:ARG:NH2	1:A:336:TYR:CZ	2.87	0.43
1:C:35:THR:HA	1:C:36:PRO:HD3	1.83	0.43
1:A:119:ILE:HD11	1:A:121:TYR:OH	2.19	0.43
1:A:379:TRP:CG	1:A:379:TRP:O	2.72	0.43
1:A:141:LYS:HD3	1:A:141:LYS:HA	1.83	0.43
1:A:398:TRP:HB3	1:A:401:LYS:CB	2.49	0.43
1:B:220:TRP:CE2	1:B:230:ASN:HB3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:289:ARG:O	1:C:289:ARG:CG	2.67	0.43
1:A:80:GLY:CA	1:A:394:TRP:CE2	3.02	0.42
1:C:65:GLY:HA2	1:C:100:ALA:O	2.18	0.42
1:C:302:ASP:HB3	1:C:385:ASP:O	2.18	0.42
1:B:188:GLY:O	1:B:246:HIS:HE1	2.02	0.42
1:C:265:TYR:CD1	1:C:266:LYS:N	2.87	0.42
1:A:90:LEU:CD2	1:A:142:LYS:HB3	2.49	0.42
1:A:381:PRO:CG	1:A:394:TRP:CE2	3.02	0.42
1:B:142:LYS:O	1:B:146:LEU:HD13	2.19	0.42
1:C:49:TYR:HE1	1:C:60:PRO:HG3	1.82	0.42
1:A:38:GLY:O	1:A:53:VAL:HG22	2.20	0.42
1:A:386:THR:C	1:A:387:GLY:O	2.57	0.42
1:A:68:TRP:HE3	1:A:380:ASN:ND2	2.14	0.42
1:A:93:ILE:HG22	1:A:94:LYS:N	2.34	0.42
1:A:379:TRP:CZ3	1:A:389:ILE:CG2	3.03	0.42
1:A:381:PRO:CA	1:A:388:GLY:N	2.72	0.42
1:C:34:GLN:OE1	1:C:34:GLN:HA	2.19	0.42
1:C:201:GLU:CD	2:G:1:BGC:O1	2.58	0.42
1:A:68:TRP:HB3	1:A:103:LEU:HD23	2.01	0.42
1:A:110:VAL:O	1:A:133:LEU:HD22	2.19	0.42
1:A:316:PRO:HG3	1:A:357:ILE:CG2	2.50	0.42
1:B:86:TRP:CZ3	1:B:142:LYS:HG3	2.55	0.42
1:B:187:PHE:HA	1:B:190:TYR:HD2	1.84	0.42
1:B:87:GLU:OE2	1:B:142:LYS:NZ	2.47	0.42
1:C:292:LEU:HD21	1:C:294:TYR:CE2	2.54	0.42
1:B:35:THR:HA	1:B:36:PRO:HD3	1.76	0.42
1:B:161:HIS:HE1	4:B:508:HOH:O	2.02	0.42
1:C:42:GLU:OE2	1:C:51:ILE:CD1	2.68	0.42
1:C:172:SER:O	1:C:175:ASP:HB2	2.20	0.42
1:C:319:LEU:H	1:C:319:LEU:CD1	2.33	0.42
1:C:319:LEU:HD21	1:C:358:TRP:HD1	1.85	0.42
1:A:72:GLU:OE2	2:D:4:BGC:O6	2.25	0.42
1:A:73:THR:HB	1:A:74:PRO:CD	2.45	0.42
1:A:94:LYS:HD3	1:A:94:LYS:C	2.40	0.42
1:A:346:LYS:CE	1:A:386:THR:C	2.88	0.42
1:C:198:LEU:CD2	1:C:250:PHE:O	2.68	0.42
1:C:242:LYS:HE3	1:C:242:LYS:HB2	1.86	0.42
1:C:283:TYR:N	1:C:284:PRO:CD	2.83	0.42
1:C:345:GLY:HA3	1:C:355:ASP:OD2	2.20	0.42
1:A:35:THR:HB	1:A:246:HIS:HB3	2.01	0.41
1:A:131:ASP:O	1:A:133:LEU:CA	2.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:MET:SD	1:A:259:PRO:HD2	2.60	0.41
1:A:346:LYS:HA	1:A:401:LYS:CG	2.49	0.41
1:B:246:HIS:CD2	1:B:246:HIS:N	2.86	0.41
1:C:271:ALA:HB1	1:C:276:ASN:HB2	2.01	0.41
1:C:316:PRO:HG3	1:C:357:ILE:HG22	2.02	0.41
1:A:74:PRO:O	1:A:76:HIS:ND1	2.39	0.41
1:C:241:LEU:HD23	1:C:241:LEU:HA	1.83	0.41
1:A:81:LEU:HB3	1:A:119:ILE:HG22	2.02	0.41
1:A:199:LYS:HB3	1:A:229:TRP:CH2	2.55	0.41
1:A:376:TYR:OH	1:A:401:LYS:HD3	2.21	0.41
1:C:279:ALA:O	1:C:283:TYR:HB2	2.21	0.41
1:A:68:TRP:CD1	1:A:78:VAL:HG21	2.55	0.41
1:A:93:ILE:O	1:A:96:LEU:HB2	2.20	0.41
1:A:106:CYS:HA	1:A:156:ARG:O	2.20	0.41
1:A:135:ILE:HD12	1:A:135:ILE:O	2.21	0.41
1:A:267:TRP:CE2	1:A:308:TYR:HB3	2.55	0.41
1:B:42:GLU:OE2	1:B:58:GLU:OE2	2.38	0.41
1:A:379:TRP:CZ3	1:A:389:ILE:HG22	2.55	0.41
1:B:77:VAL:O	1:B:78:VAL:C	2.59	0.41
1:C:257:THR:OG1	1:C:258:ASN:N	2.52	0.41
1:C:323:TRP:CH2	1:C:362:LEU:HA	2.55	0.41
1:C:368:GLU:C	1:C:370:LYS:H	2.24	0.41
1:C:304:TYR:CE2	3:F:1:BGC:H5	2.56	0.41
1:A:72:GLU:OE1	1:A:72:GLU:N	2.44	0.41
1:A:100:ALA:HA	1:A:149:PHE:O	2.21	0.41
1:A:344:GLY:HA2	1:A:376:TYR:CE1	2.56	0.41
1:A:383:SER:OG	1:A:385:ASP:OD2	2.37	0.41
1:A:398:TRP:HB3	1:A:401:LYS:HB3	2.01	0.41
1:B:206:THR:O	1:B:211:ALA:HB2	2.21	0.41
1:C:180:TRP:CE3	1:C:183:VAL:HG21	2.56	0.41
1:C:205:VAL:CG2	1:C:216:THR:O	2.63	0.41
1:A:130:LEU:CB	1:A:134:GLN:HE21	2.34	0.41
1:C:184:ALA:HA	1:C:247:TRP:CH2	2.55	0.41
1:A:316:PRO:HG3	1:A:357:ILE:HG21	2.04	0.40
1:A:72:GLU:OE2	2:D:4:BGC:C6	2.70	0.40
1:A:133:LEU:HD12	1:A:133:LEU:HA	1.70	0.40
1:A:343:PHE:O	1:A:377:TRP:N	2.40	0.40
1:A:347:TYR:CG	1:A:356:VAL:HG13	2.56	0.40
1:C:198:LEU:HD21	1:C:249:ILE:HG22	1.98	0.40
1:A:315:PHE:HA	1:A:316:PRO:HA	1.85	0.40
1:C:46:ASP:OD1	1:C:46:ASP:C	2.60	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:173:GLU:O	1:C:177:ILE:HG13	2.22	0.40
1:C:197:ASP:HA	1:C:250:PHE:HB2	2.03	0.40
1:B:299:TYR:CD1	1:B:303:VAL:HG21	2.56	0.40
1:C:331:LYS:HB2	1:C:338:VAL:CG2	2.51	0.40
1:C:131:ASP:OD1	1:C:134:GLN:CG	2.70	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:554:HOH:O	4:C:500:HOH:O[4_545]	2.03	0.17
4:B:553:HOH:O	4:C:553:HOH:O[4_545]	2.10	0.10
1:A:95:SER:OG	1:A:354:ARG:NH1[4_556]	2.16	0.04

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	375/458 (82%)	338 (90%)	31 (8%)	6 (2%)	9 4
1	B	375/458 (82%)	361 (96%)	14 (4%)	0	100 100
1	C	375/458 (82%)	354 (94%)	20 (5%)	1 (0%)	41 37
All	All	1125/1374 (82%)	1053 (94%)	65 (6%)	7 (1%)	25 19

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	SER
1	A	384	GLY
1	A	390	LEU
1	A	112	PRO
1	C	215	GLY

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Mol	Chain	Res	Type
1	A	327	PHE
1	A	350	GLY

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	323/395 (82%)	285 (88%)	38 (12%)	5	3
1	B	323/395 (82%)	311 (96%)	12 (4%)	34	32
1	C	323/395 (82%)	306 (95%)	17 (5%)	22	18
All	All	969/1185 (82%)	902 (93%)	67 (7%)	15	11

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ASP
1	A	68	TRP
1	A	69	PHE
1	A	75	ASN
1	A	78	VAL
1	A	79	HIS
1	A	87	GLU
1	A	88	ASP
1	A	92	GLN
1	A	94	LYS
1	A	112	PRO
1	A	126	ASP
1	A	130	LEU
1	A	134	GLN
1	A	135	ILE
1	A	141	LYS
1	A	153	ASP
1	A	154	TYR
1	A	180	TRP
1	A	190	TYR

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Mol	Chain	Res	Type
1	A	281	LYS
1	A	290	ASN
1	A	297	HIS
1	A	313	LYS
1	A	319	LEU
1	A	326	HIS
1	A	343	PHE
1	A	346	LYS
1	A	347	TYR
1	A	354	ARG
1	A	389	ILE
1	A	390	LEU
1	A	392	ASP
1	A	393	ASP
1	A	396	THR
1	A	397	ILE
1	A	399	GLU
1	A	408	LEU
1	B	57	GLU
1	B	68	TRP
1	B	115	GLN
1	B	142	LYS
1	B	153	ASP
1	B	154	TYR
1	B	241	LEU
1	B	246	HIS
1	B	297	HIS
1	B	380	ASN
1	B	399	GLU
1	B	406	LYS
1	C	58	GLU
1	C	138	LYS
1	C	154	TYR
1	C	160	THR
1	C	214	ASP
1	C	246	HIS
1	C	260	LYS
1	C	270	ASN
1	C	282	ASP
1	C	287	LEU
1	C	289	ARG
1	C	297	HIS

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Mol	Chain	Res	Type
1	C	317	ASP
1	C	334	LEU
1	C	380	ASN
1	C	407	ARG
1	C	410	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	HIS
1	A	75	ASN
1	A	85	ASN
1	A	124	ASN
1	A	134	GLN
1	A	161	HIS
1	A	203	HIS
1	A	246	HIS
1	A	290	ASN
1	A	297	HIS
1	A	325	HIS
1	A	359	GLN
1	B	62	HIS
1	B	75	ASN
1	B	79	HIS
1	B	85	ASN
1	B	115	GLN
1	B	161	HIS
1	B	178	ASN
1	B	203	HIS
1	B	246	HIS
1	B	270	ASN
1	B	297	HIS
1	B	325	HIS
1	B	326	HIS
1	B	349	HIS
1	B	359	GLN
1	B	380	ASN
1	B	391	GLN
1	C	62	HIS
1	C	75	ASN
1	C	79	HIS
1	C	85	ASN

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Mol	Chain	Res	Type
1	C	203	HIS
1	C	246	HIS
1	C	255	GLN
1	C	270	ASN
1	C	297	HIS
1	C	325	HIS
1	C	349	HIS
1	C	359	GLN
1	C	380	ASN

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](#)

14 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	BGC	D	1	2	12,12,12	1.23	2 (16%)	17,17,17	1.58	4 (23%)
2	BGC	D	2	2	11,11,12	0.79	0	15,15,17	1.15	2 (13%)
2	BGC	D	3	2	11,11,12	0.90	0	15,15,17	1.49	2 (13%)
2	BGC	D	4	2	11,11,12	0.68	0	15,15,17	1.27	3 (20%)
2	BGC	E	1	2	12,12,12	1.25	0	17,17,17	1.75	3 (17%)
2	BGC	E	2	2	11,11,12	0.89	0	15,15,17	1.14	2 (13%)
2	BGC	E	3	2	11,11,12	1.03	1 (9%)	15,15,17	1.20	1 (6%)
2	BGC	E	4	2	11,11,12	0.73	0	15,15,17	0.96	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BGC	F	1	3	12,12,12	0.51	0	17,17,17	0.93	2 (11%)
3	BGC	F	2	3	11,11,12	0.32	0	15,15,17	0.99	1 (6%)
2	BGC	G	1	2	12,12,12	1.39	0	17,17,17	2.15	5 (29%)
2	BGC	G	2	2	11,11,12	0.88	0	15,15,17	1.08	1 (6%)
2	BGC	G	3	2	11,11,12	0.89	1 (9%)	15,15,17	1.24	2 (13%)
2	BGC	G	4	2	11,11,12	0.81	0	15,15,17	1.06	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	D	1	2	-	0/2/22/22	0/1/1/1
2	BGC	D	2	2	-	0/2/19/22	0/1/1/1
2	BGC	D	3	2	-	0/2/19/22	0/1/1/1
2	BGC	D	4	2	-	0/2/19/22	0/1/1/1
2	BGC	E	1	2	-	0/2/22/22	0/1/1/1
2	BGC	E	2	2	-	0/2/19/22	0/1/1/1
2	BGC	E	3	2	-	0/2/19/22	0/1/1/1
2	BGC	E	4	2	-	0/2/19/22	0/1/1/1
3	BGC	F	1	3	-	0/2/22/22	0/1/1/1
3	BGC	F	2	3	-	0/2/19/22	0/1/1/1
2	BGC	G	1	2	-	0/2/22/22	0/1/1/1
2	BGC	G	2	2	-	0/2/19/22	0/1/1/1
2	BGC	G	3	2	-	0/2/19/22	0/1/1/1
2	BGC	G	4	2	-	0/2/19/22	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	3	BGC	C4-C5	2.15	1.57	1.53
2	G	3	BGC	O5-C1	2.15	1.47	1.43
2	D	1	BGC	O5-C5	2.09	1.49	1.44
2	D	1	BGC	O4-C4	2.03	1.47	1.43

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	1	BGC	C1-C2-C3	-4.79	100.37	110.31
2	G	1	BGC	C1-O5-C5	-4.38	105.40	113.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3	BGC	C1-C2-C3	-4.03	104.72	109.67
2	E	1	BGC	C1-O5-C5	-3.85	106.39	113.66
2	E	1	BGC	C1-C2-C3	-3.80	102.44	110.31
2	D	1	BGC	C1-O5-C5	-3.73	106.62	113.66
2	D	1	BGC	C1-C2-C3	-2.89	104.32	110.31
2	D	4	BGC	C1-C2-C3	-2.87	106.14	109.67
2	G	3	BGC	C1-C2-C3	-2.82	106.20	109.67
2	G	1	BGC	O4-C4-C5	-2.54	103.00	109.30
3	F	1	BGC	C1-O5-C5	-2.52	108.91	113.66
2	D	2	BGC	C1-C2-C3	-2.48	106.62	109.67
2	D	1	BGC	C6-C5-C4	-2.44	107.28	113.00
2	D	2	BGC	C6-C5-C4	-2.35	107.50	113.00
2	G	2	BGC	C1-C2-C3	-2.33	106.81	109.67
2	G	4	BGC	C6-C5-C4	-2.21	107.82	113.00
2	G	3	BGC	C6-C5-C4	-2.20	107.86	113.00
2	D	3	BGC	C1-O5-C5	-2.12	109.31	112.19
3	F	1	BGC	C3-C4-C5	-2.12	106.46	110.24
2	E	2	BGC	C6-C5-C4	-2.10	108.09	113.00
2	E	2	BGC	C1-C2-C3	-2.09	107.10	109.67
2	G	1	BGC	O4-C4-C3	2.07	115.14	110.35
2	E	4	BGC	C2-C3-C4	-2.07	107.32	110.89
2	D	1	BGC	O2-C2-C3	-2.06	105.59	110.35
3	F	2	BGC	O5-C1-C2	-2.05	107.61	110.77
2	E	1	BGC	C6-C5-C4	-2.05	108.21	113.00
2	E	3	BGC	C6-C5-C4	-2.04	108.23	113.00
2	G	1	BGC	O2-C2-C3	-2.03	105.66	110.35
2	D	4	BGC	C2-C3-C4	-2.03	107.39	110.89
2	D	4	BGC	C1-O5-C5	-2.02	109.46	112.19

There are no chirality outliers.

There are no torsion outliers.

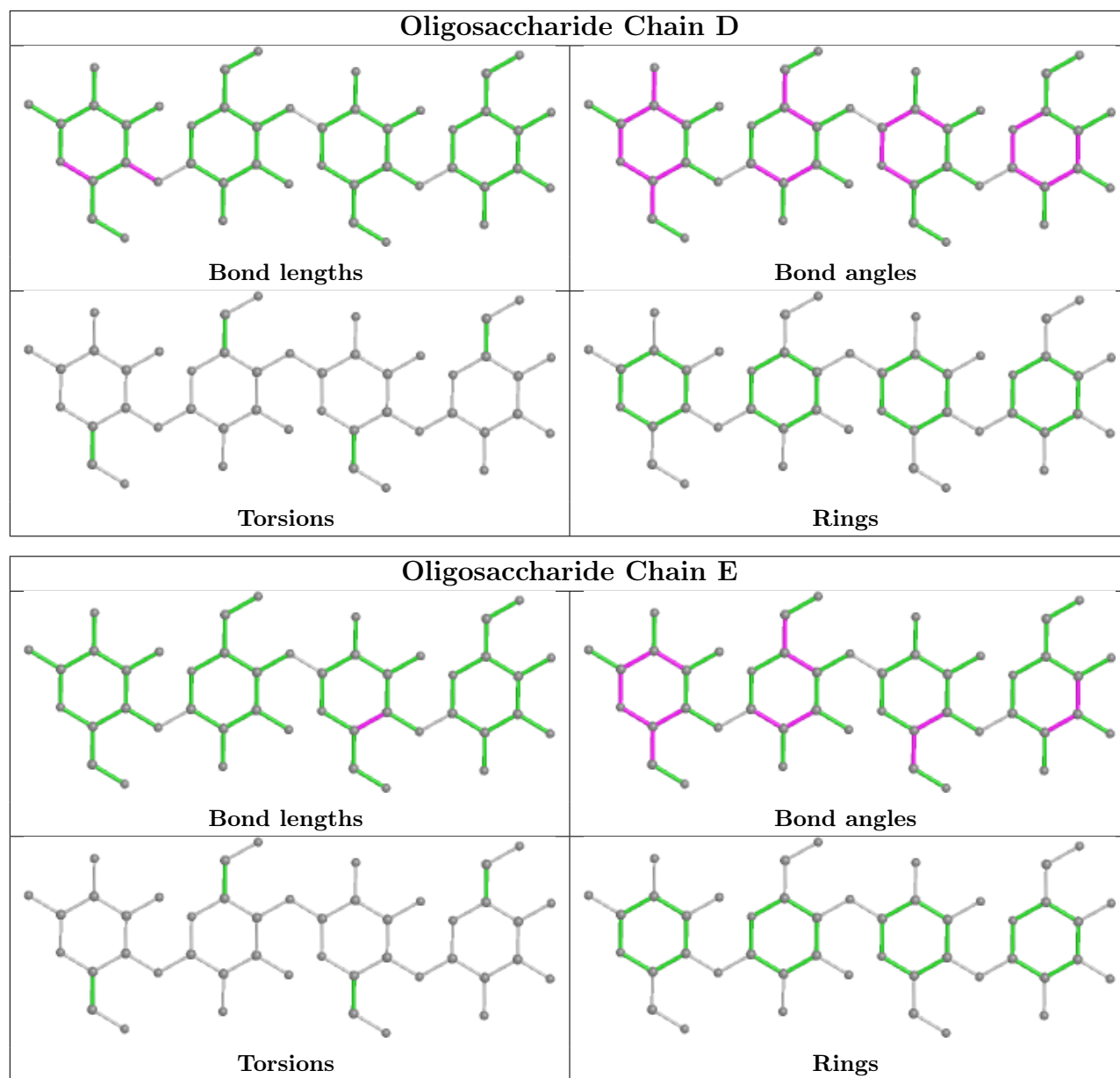
There are no ring outliers.

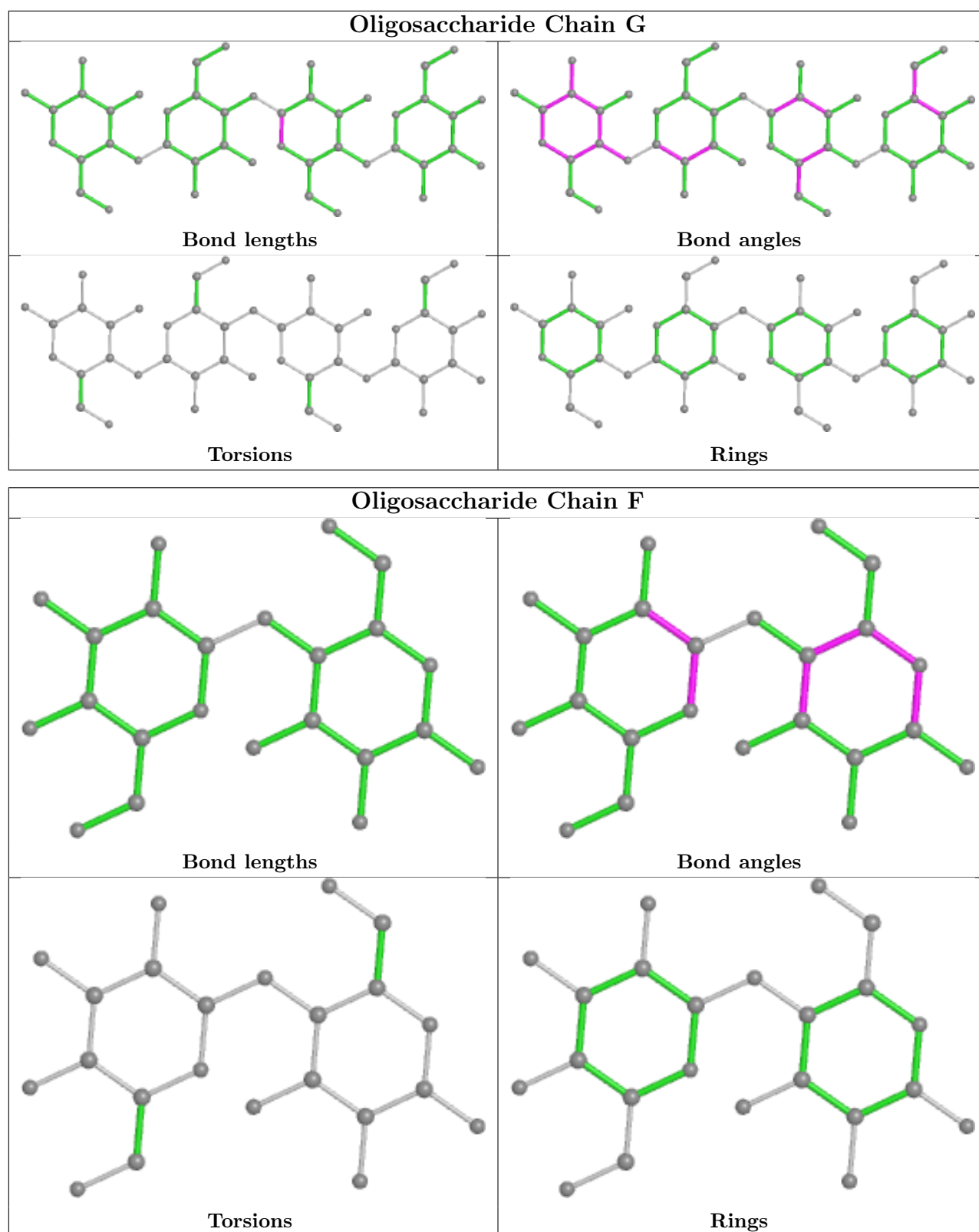
4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	1	BGC	2	0
2	D	3	BGC	1	0
3	F	1	BGC	2	0
2	D	4	BGC	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

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	377/458 (82%)	0.54	35 (9%) 8 8	26, 52, 84, 113	2 (0%)
1	B	377/458 (82%)	-0.11	10 (2%) 54 53	20, 33, 55, 85	3 (0%)
1	C	377/458 (82%)	0.62	57 (15%) 2 1	29, 56, 81, 101	2 (0%)
All	All	1131/1374 (82%)	0.35	102 (9%) 9 8	20, 46, 79, 113	7 (0%)

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	127	LEU	6.8
1	C	269	TYR	5.6
1	A	122	SER	5.5
1	C	151	LEU	5.4
1	C	342	ALA	4.8
1	C	216	THR	4.6
1	C	53	VAL	4.6
1	C	340	ILE	4.6
1	A	349	HIS	4.5
1	C	217	GLY	4.4
1	C	205	VAL	4.3
1	C	375	PHE	4.1
1	A	128	ARG	4.1
1	C	265	TYR	3.9
1	C	102	ARG	3.8
1	C	264	SER	3.8
1	A	129	GLY	3.8
1	B	170	ASP	3.8
1	C	315	PHE	3.8
1	C	101	ILE	3.7
1	C	152	LEU	3.7
1	C	66	VAL	3.6
1	C	267	TRP	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	295	SER	3.5
1	C	251	VAL	3.4
1	C	250	PHE	3.4
1	C	263	SER	3.3
1	A	394	TRP	3.3
1	C	208	PRO	3.3
1	A	379	TRP	3.3
1	C	341	GLY	3.2
1	A	377	TRP	3.1
1	C	332	LEU	3.1
1	C	311	PRO	3.1
1	C	339	VAL	3.1
1	A	118	GLY	3.1
1	C	103	LEU	3.1
1	A	53	VAL	3.0
1	C	206	THR	3.0
1	C	54	THR	2.9
1	C	334	LEU	2.9
1	A	342	ALA	2.9
1	A	340	ILE	2.9
1	A	344	GLY	2.9
1	A	135	ILE	2.8
1	C	104	PRO	2.8
1	C	150	VAL	2.8
1	C	214	ASP	2.8
1	B	104	PRO	2.7
1	A	375	PHE	2.7
1	C	212	TYR	2.7
1	A	142	LYS	2.7
1	B	151	LEU	2.7
1	A	267	TRP	2.6
1	C	196	ALA	2.6
1	A	121	TYR	2.6
1	C	211	ALA	2.6
1	C	336	TYR	2.5
1	A	88	ASP	2.4
1	A	80	GLY	2.4
1	C	45	GLY	2.4
1	A	136	MET	2.4
1	A	35	THR	2.4
1	A	82	TRP	2.4
1	C	58	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	380	ASN	2.4
1	A	95	SER	2.4
1	B	171	PHE	2.3
1	C	343	PHE	2.3
1	C	243	VAL	2.3
1	C	262	ASP	2.3
1	C	194	ILE	2.3
1	C	63	LEU	2.3
1	B	68	TRP	2.3
1	C	260	LYS	2.3
1	C	170	ASP	2.2
1	A	38	GLY	2.2
1	C	207	SER	2.2
1	A	119	ILE	2.2
1	C	374	PHE	2.2
1	A	296	PRO	2.2
1	B	295	SER	2.2
1	A	130	LEU	2.2
1	C	195	GLY	2.2
1	B	169	GLU	2.1
1	A	138	LYS	2.1
1	A	346	LYS	2.1
1	B	375	PHE	2.1
1	C	283	TYR	2.1
1	B	250	PHE	2.1
1	C	153	ASP	2.1
1	C	65	GLY	2.1
1	A	89	MET	2.1
1	A	159	CYS	2.1
1	C	143	ALA	2.0
1	C	171	PHE	2.0
1	C	309	PHE	2.0
1	B	102	ARG	2.0
1	C	197	ASP	2.0
1	A	390	LEU	2.0
1	C	198	LEU	2.0
1	A	83	LYS	2.0

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

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

6.3 Carbohydrates

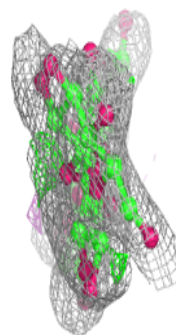
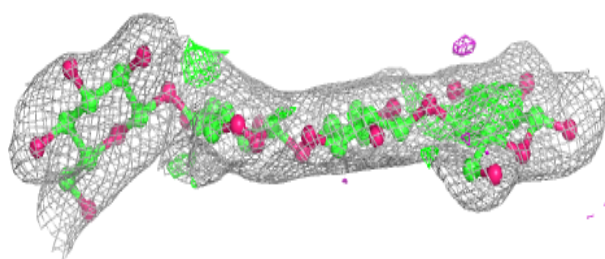
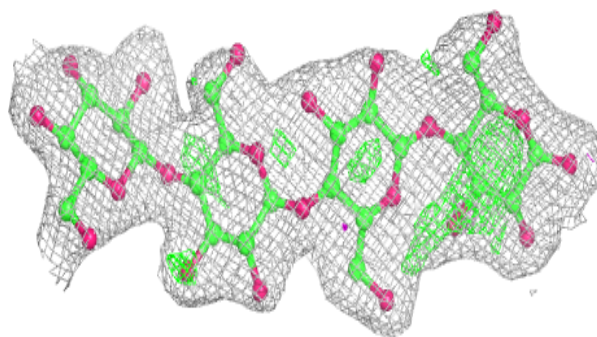
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	BGC	G	1	12/12	0.85	0.17	50,57,67,68	0
3	BGC	F	2	11/12	0.85	0.13	63,68,75,77	0
2	BGC	G	4	11/12	0.87	0.17	49,58,69,76	0
2	BGC	D	3	11/12	0.88	0.14	56,61,65,71	0
3	BGC	F	1	12/12	0.89	0.11	58,68,75,78	0
2	BGC	E	1	12/12	0.89	0.14	36,42,49,66	0
2	BGC	D	1	12/12	0.90	0.14	38,48,59,71	0
2	BGC	D	2	11/12	0.90	0.14	39,44,60,64	0
2	BGC	D	4	11/12	0.91	0.16	46,56,70,72	0
2	BGC	E	2	11/12	0.92	0.11	28,35,42,47	0
2	BGC	G	2	11/12	0.93	0.10	40,46,51,52	0
2	BGC	G	3	11/12	0.94	0.15	53,61,67,68	0
2	BGC	E	4	11/12	0.94	0.18	31,33,41,43	0
2	BGC	E	3	11/12	0.95	0.15	23,26,37,55	0

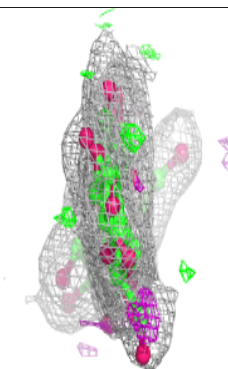
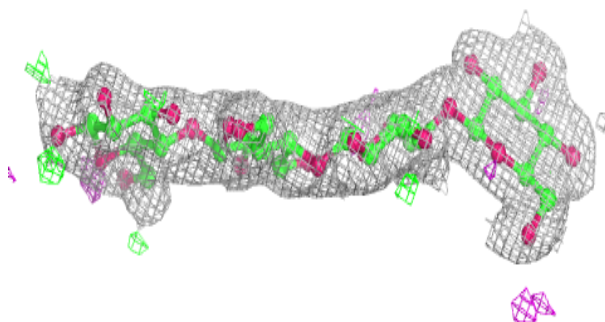
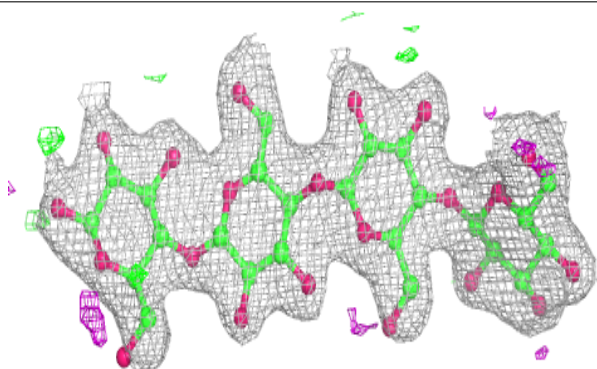
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

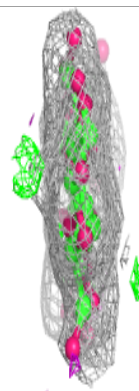
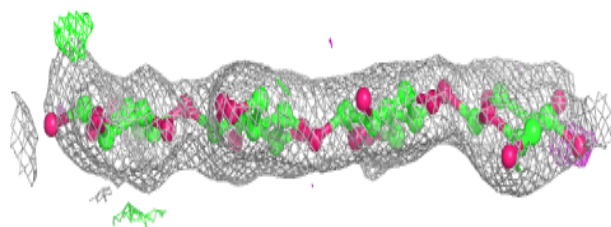
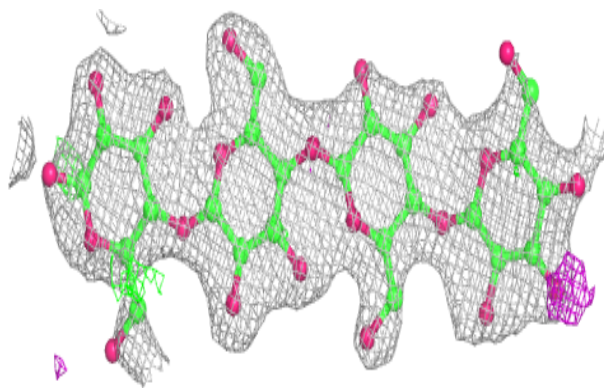
**Electron density around Chain E:**

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

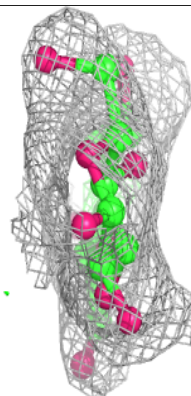
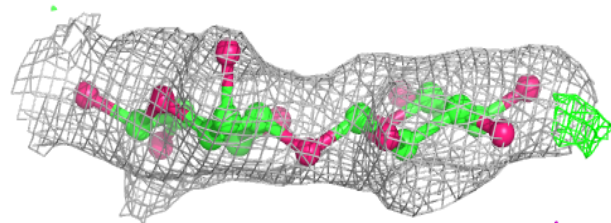
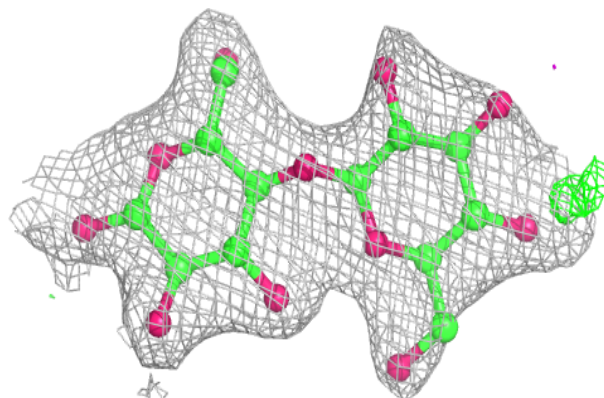


Electron density around Chain G:

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

**Electron density around Chain 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.