



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 16, 2020 – 04:09 PM JST

PDB ID : 6IK7
Title : Crystal structure of tomato beta-galactosidase (TBG) 4 in complex with beta-1,3-galactobiose
Authors : Matsuyama, K.; Nakae, S.; Igarashi, K.; Tada, T.; Ishimaru, M.
Deposited on : 2018-10-15
Resolution : 3.10 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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 : 2.15.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.15.1

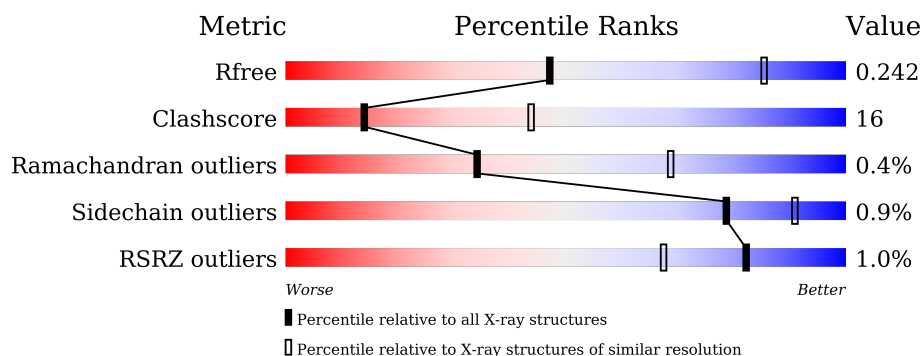
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	718	<div> <div></div> <div>73%</div> <div>25%</div> <div>..</div> </div>
1	B	718	<div> <div></div> <div>72%</div> <div>25%</div> <div>..</div> </div>
2	C	2	<div> <div>50%</div> <div>50%</div> </div>
2	E	2	<div> <div>100%</div> </div>
3	D	2	<div> <div>50%</div> <div>50%</div> </div>
3	F	2	<div> <div>50%</div> <div>50%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11293 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 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	705	Total	C	N	O	S	0	0	0
			5527	3533	944	1023	27			
1	A	705	Total	C	N	O	S	0	0	0
			5527	3533	944	1023	27			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	18	GLU	-	expression tag	UNP O81100
B	19	ALA	-	expression tag	UNP O81100
B	20	GLU	-	expression tag	UNP O81100
B	21	ALA	-	expression tag	UNP O81100
B	22	GLU	-	expression tag	UNP O81100
B	23	PHE	-	expression tag	UNP O81100
B	181	ALA	GLU	engineered mutation	UNP O81100
B	725	SER	-	expression tag	UNP O81100
B	726	ALA	-	expression tag	UNP O81100
B	727	ALA	-	expression tag	UNP O81100
B	728	ALA	-	expression tag	UNP O81100
B	729	ALA	-	expression tag	UNP O81100
B	730	SER	-	expression tag	UNP O81100
B	731	PHE	-	expression tag	UNP O81100
B	732	LEU	-	expression tag	UNP O81100
B	733	GLU	-	expression tag	UNP O81100
B	734	GLN	-	expression tag	UNP O81100
B	735	LYS	-	expression tag	UNP O81100
A	18	GLU	-	expression tag	UNP O81100
A	19	ALA	-	expression tag	UNP O81100
A	20	GLU	-	expression tag	UNP O81100
A	21	ALA	-	expression tag	UNP O81100
A	22	GLU	-	expression tag	UNP O81100
A	23	PHE	-	expression tag	UNP O81100
A	181	ALA	GLU	engineered mutation	UNP O81100

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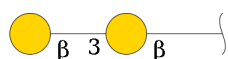
Chain	Residue	Modelled	Actual	Comment	Reference
A	725	SER	-	expression tag	UNP O81100
A	726	ALA	-	expression tag	UNP O81100
A	727	ALA	-	expression tag	UNP O81100
A	728	ALA	-	expression tag	UNP O81100
A	729	ALA	-	expression tag	UNP O81100
A	730	SER	-	expression tag	UNP O81100
A	731	PHE	-	expression tag	UNP O81100
A	732	LEU	-	expression tag	UNP O81100
A	733	GLU	-	expression tag	UNP O81100
A	734	GLN	-	expression tag	UNP O81100
A	735	LYS	-	expression tag	UNP O81100

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



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

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



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

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

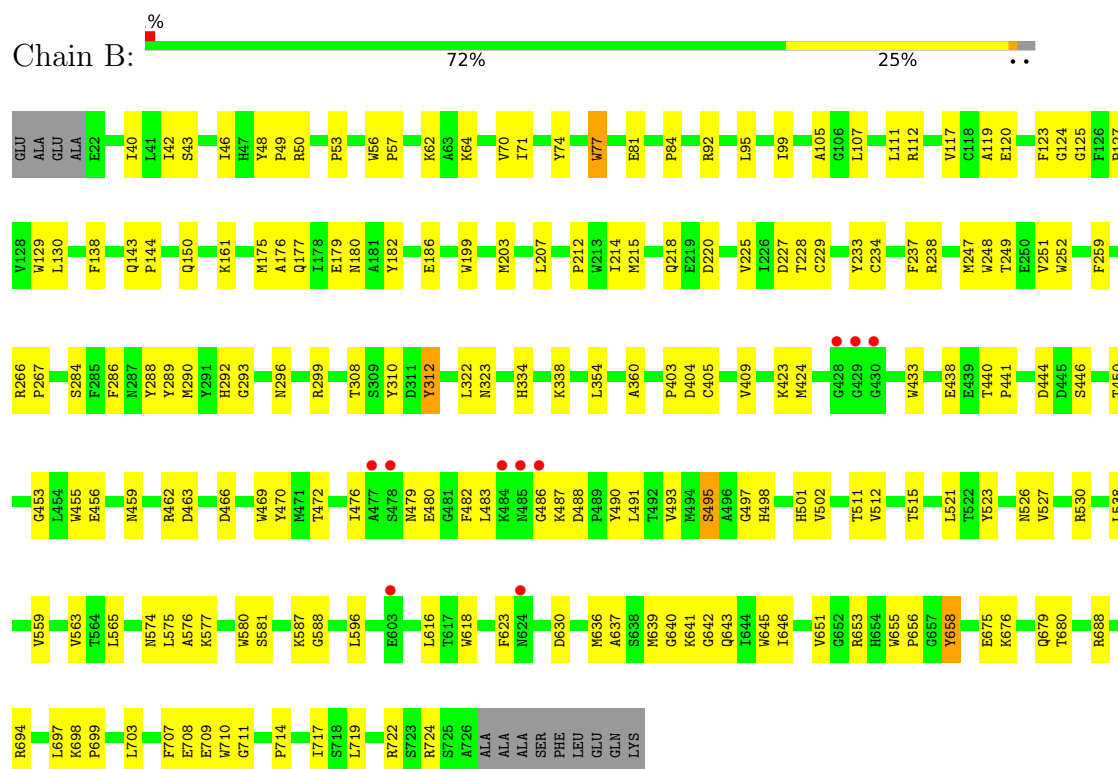
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	59	Total	O	0	0
			59	59		
5	A	50	Total	O	0	0
			50	50		

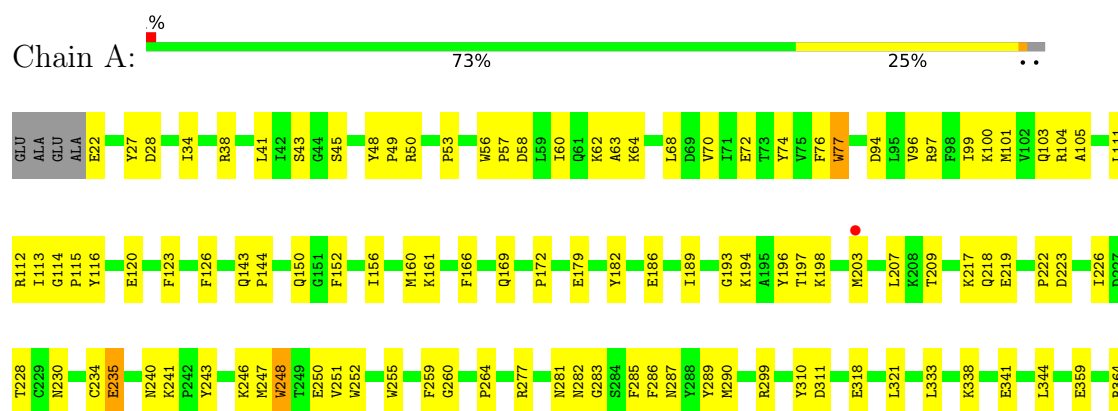
3 Residue-property plots

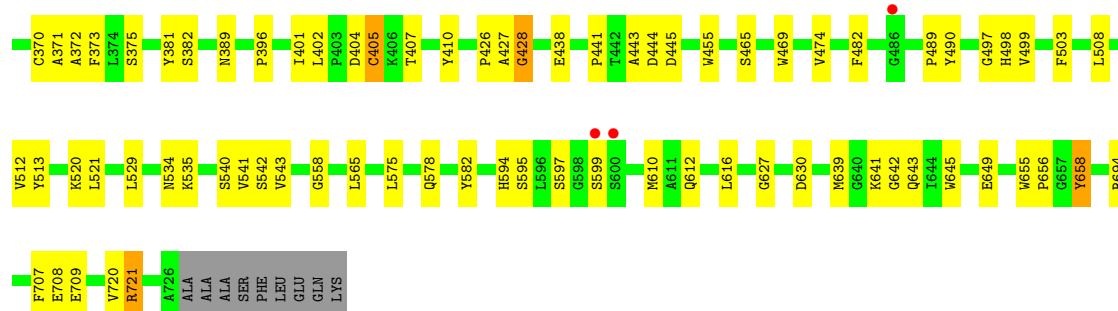
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Beta-galactosidase



• Molecule 1: Beta-galactosidase





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

Chain C: 50% 50%

MAG1
MAG2

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

Chain E: 100%

MAG1
MAG2

- Molecule 3: beta-D-galactopyranose-(1-3)-beta-D-galactopyranose

Chain D: 50% 50%

GAL1
GAL2

- Molecule 3: beta-D-galactopyranose-(1-3)-beta-D-galactopyranose

Chain F: 50% 50%

GAL1
GAL2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.61Å 97.95Å 161.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.43 – 3.10 43.40 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.43-3.10) 99.9 (43.40-3.10)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.88 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.8.0232	Depositor
R, R_{free}	0.199 , 0.244 0.201 , 0.242	Depositor DCC
R_{free} test set	1375 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.089	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.028 for k,h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	11293	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% 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: GAL, NAG

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	0.62	0/5690	0.75	2/7734 (0.0%)
1	B	0.61	0/5690	0.75	0/7734
All	All	0.61	0/11380	0.75	2/15468 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	TRP	CA-CB-CG	5.89	124.90	113.70
1	A	77	TRP	CA-CB-CG	-5.84	102.60	113.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5527	0	5349	170	0
1	B	5527	0	5349	174	0
2	C	28	0	25	0	0
2	E	28	0	25	1	0
3	D	23	0	21	1	0
3	F	23	0	21	2	0
4	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	14	0	13	4	0
5	A	50	0	0	4	0
5	B	59	0	0	5	0
All	All	11293	0	10816	345	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 16.

All (345) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:TRP:CZ3	1:B:81:GLU:CD	1.95	1.40
1:B:641:LYS:HD2	1:B:655:TRP:CZ3	1.57	1.40
1:A:101:MET:CE	1:A:104:ARG:HH21	1.35	1.38
1:A:370:CYS:C	1:A:405:CYS:SG	2.10	1.29
1:B:641:LYS:HG3	1:B:655:TRP:CE3	1.73	1.22
1:B:641:LYS:CG	1:B:655:TRP:CE3	2.27	1.18
1:B:77:TRP:CH2	1:B:81:GLU:CD	2.20	1.14
1:B:641:LYS:CG	1:B:655:TRP:HE3	1.57	1.13
1:B:248:TRP:CD1	1:B:251:VAL:HG22	1.84	1.12
1:B:77:TRP:CH2	1:B:81:GLU:OE2	2.03	1.11
1:B:77:TRP:CZ3	1:B:81:GLU:OE2	2.01	1.10
1:A:101:MET:HE1	1:A:104:ARG:HH21	1.16	1.06
1:A:235:GLU:OE1	1:A:277:ARG:HD3	1.56	1.06
1:B:641:LYS:CD	1:B:655:TRP:CZ3	2.38	1.05
1:B:77:TRP:CH2	1:B:81:GLU:OE1	2.09	1.04
1:A:101:MET:CE	1:A:104:ARG:NH2	2.20	1.04
1:A:248:TRP:NE1	1:A:251:VAL:HG22	1.73	1.04
1:A:101:MET:HE3	1:A:104:ARG:HH21	1.17	1.04
1:A:259:PHE:HB3	1:A:655:TRP:CZ3	1.93	1.03
1:A:370:CYS:O	1:A:405:CYS:SG	2.16	1.02
1:A:48:TYR:OH	1:A:76:PHE:N	1.94	0.99
1:B:641:LYS:HG3	1:B:655:TRP:HE3	1.04	0.98
1:B:248:TRP:NE1	1:B:251:VAL:HG22	1.76	0.98
1:B:77:TRP:CE3	1:B:81:GLU:OE2	2.19	0.94
1:A:655:TRP:CZ3	1:A:658:TYR:HB3	2.00	0.93
1:B:641:LYS:HD2	1:B:655:TRP:HZ3	1.13	0.92
1:B:655:TRP:HE1	1:B:658:TYR:HD2	1.16	0.92
1:A:655:TRP:CZ3	1:A:658:TYR:CB	2.45	0.90
1:B:77:TRP:CZ2	1:B:81:GLU:OE2	2.23	0.90
1:A:235:GLU:OE1	1:A:277:ARG:CD	2.20	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:CYS:HB2	1:A:248:TRP:CE3	2.08	0.88
1:B:459:ASN:HD22	4:B:803:NAG:H83	1.40	0.86
1:A:299:ARG:HE	1:A:643:GLN:HE22	1.25	0.84
1:B:641:LYS:CD	1:B:655:TRP:CE3	2.59	0.83
1:A:101:MET:HE3	1:A:104:ARG:NH2	1.84	0.83
1:B:77:TRP:CZ3	1:B:81:GLU:OE1	2.28	0.82
1:A:234:CYS:HB2	1:A:248:TRP:CZ3	2.14	0.82
1:A:259:PHE:CB	1:A:655:TRP:CZ3	2.61	0.82
1:B:641:LYS:CB	1:B:655:TRP:HE3	1.93	0.81
1:A:259:PHE:HB3	1:A:655:TRP:CE3	2.16	0.80
1:A:101:MET:HE1	1:A:104:ARG:NH2	1.89	0.80
1:A:259:PHE:CD1	1:A:655:TRP:CE3	2.71	0.79
1:A:161:LYS:HA	1:A:169:GLN:HE22	1.48	0.79
1:B:77:TRP:CE2	1:B:81:GLU:OE2	2.35	0.79
1:B:698:LYS:HB2	1:B:699:PRO:HD2	1.64	0.79
1:B:641:LYS:HB2	1:B:655:TRP:CE3	2.17	0.78
1:B:636:MET:HE3	1:B:639:MET:HG3	1.64	0.78
1:B:77:TRP:CD2	1:B:81:GLU:OE2	2.36	0.78
1:B:641:LYS:CB	1:B:655:TRP:CE3	2.66	0.77
1:B:655:TRP:NE1	1:B:658:TYR:HB2	2.01	0.76
1:A:48:TYR:CZ	1:A:76:PHE:HD2	2.05	0.73
1:B:81:GLU:OE2	1:B:127:PRO:HG2	1.89	0.73
1:B:565:LEU:HB2	1:B:575:LEU:HD11	1.72	0.72
1:A:438:GLU:HB3	1:A:616:LEU:HB3	1.72	0.71
1:B:488:ASP:O	1:B:526:ASN:ND2	2.23	0.71
1:A:299:ARG:HE	1:A:643:GLN:NE2	1.87	0.71
1:A:259:PHE:CA	1:A:655:TRP:CZ3	2.74	0.70
1:A:48:TYR:HH	1:A:76:PHE:H	1.34	0.70
1:A:234:CYS:SG	1:A:248:TRP:HZ3	2.15	0.70
1:A:247:MET:HG2	1:A:285:PHE:HB3	1.74	0.69
1:A:259:PHE:CD1	1:A:655:TRP:HE3	2.10	0.69
1:A:248:TRP:HD1	1:A:286:PHE:HB2	1.58	0.69
1:B:179:GLU:HG3	1:B:180:ASN:H	1.56	0.69
1:B:248:TRP:HE1	1:B:251:VAL:HG22	1.55	0.68
1:B:77:TRP:CE3	1:B:81:GLU:CD	2.61	0.68
1:A:259:PHE:CG	1:A:655:TRP:CE3	2.82	0.68
1:A:259:PHE:C	1:A:655:TRP:CZ3	2.67	0.68
1:A:64:LYS:NZ	1:A:105:ALA:O	2.27	0.68
1:A:259:PHE:CB	1:A:655:TRP:CE3	2.76	0.68
1:A:234:CYS:SG	1:A:248:TRP:CZ3	2.87	0.67
1:B:512:VAL:HG13	1:B:521:LEU:HD13	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:595:SER:O	1:A:599:SER:HB2	1.95	0.67
1:A:248:TRP:CD1	1:A:251:VAL:HG22	2.30	0.67
1:B:215:MET:HB3	1:B:218:GLN:HG3	1.77	0.66
1:A:48:TYR:CZ	1:A:76:PHE:CD2	2.83	0.66
1:B:433:TRP:NE1	1:B:714:PRO:O	2.29	0.66
1:B:234:CYS:SG	1:B:237:PHE:HB2	2.36	0.66
1:B:50:ARG:NH1	1:B:296:ASN:OD1	2.29	0.66
1:A:260:GLY:N	1:A:655:TRP:HZ3	1.93	0.66
1:A:193:GLY:O	1:A:197:THR:HG23	1.96	0.66
1:A:260:GLY:N	1:A:655:TRP:CZ3	2.65	0.65
1:B:81:GLU:HG3	1:B:81:GLU:O	1.96	0.65
1:A:499:VAL:HB	1:A:541:VAL:HB	1.77	0.65
1:A:259:PHE:HB3	1:A:655:TRP:CH2	2.32	0.64
1:B:470:TYR:HB2	1:B:538:LEU:HB2	1.79	0.64
1:A:234:CYS:CB	1:A:248:TRP:CZ3	2.80	0.64
1:B:175:MET:HE2	1:B:214:ILE:HG23	1.80	0.64
1:A:250:GLU:OE1	3:F:2:GAL:H3	1.98	0.63
1:A:235:GLU:OE1	1:A:277:ARG:NE	2.31	0.63
1:B:77:TRP:CZ3	1:B:81:GLU:CG	2.82	0.63
1:A:299:ARG:HH21	1:A:465:SER:HA	1.64	0.62
1:B:175:MET:CE	1:B:214:ILE:HG23	2.29	0.61
1:A:248:TRP:CE2	1:A:251:VAL:HG22	2.36	0.61
1:B:655:TRP:N	1:B:656:PRO:HD3	2.15	0.61
1:A:474:VAL:HG13	1:A:534:ASN:HB2	1.83	0.60
1:B:248:TRP:CD1	1:B:251:VAL:CG2	2.74	0.59
1:A:512:VAL:HG23	1:A:521:LEU:HD12	1.85	0.59
1:B:62:LYS:HG2	1:B:322:LEU:HD21	1.83	0.59
1:B:512:VAL:CG1	1:B:521:LEU:HD13	2.32	0.59
1:A:543:VAL:O	1:A:543:VAL:HG12	2.02	0.58
1:B:641:LYS:CG	1:B:655:TRP:CZ3	2.77	0.58
1:B:48:TYR:CG	1:B:49:PRO:HD3	2.38	0.58
1:A:248:TRP:NE1	1:A:251:VAL:CG2	2.60	0.57
1:A:48:TYR:CG	1:A:49:PRO:HD3	2.39	0.57
1:B:490:TYR:O	1:B:565:LEU:HD12	2.04	0.57
1:A:48:TYR:CE1	1:A:76:PHE:CD2	2.92	0.57
1:A:655:TRP:CE3	1:A:658:TYR:HB2	2.40	0.57
1:B:182:TYR:HB3	1:B:218:GLN:OE1	2.05	0.57
1:A:240:ASN:O	1:A:241:LYS:HD2	2.05	0.57
1:A:627:GLY:O	1:A:721:ARG:NH2	2.38	0.57
1:B:234:CYS:N	1:B:248:TRP:CH2	2.72	0.57
1:A:426:PRO:HA	1:A:720:VAL:HG12	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:TYR:HB2	1:A:126:PHE:CZ	2.41	0.56
1:A:655:TRP:CE3	1:A:658:TYR:CB	2.89	0.55
1:A:166:PHE:CE1	1:A:209:THR:HB	2.41	0.55
1:B:74:TYR:CE2	1:B:119:ALA:HB2	2.42	0.55
1:A:655:TRP:CZ3	1:A:658:TYR:HB2	2.37	0.55
1:A:655:TRP:CD2	1:A:658:TYR:HB2	2.42	0.55
1:A:444:ASP:OD1	1:A:445:ASP:N	2.40	0.55
1:A:45:SER:HA	1:A:72:GLU:O	2.06	0.55
1:A:482:PHE:CD1	1:A:489:PRO:HD3	2.42	0.55
1:B:179:GLU:HG2	1:B:182:TYR:HB2	1.89	0.55
1:A:53:PRO:HA	1:A:56:TRP:CD2	2.43	0.54
1:B:84:PRO:HG3	1:B:129:TRP:CE3	2.43	0.54
1:A:594:HIS:HB2	1:A:645:TRP:CD2	2.43	0.54
1:A:381:TYR:C	1:A:396:PRO:HG3	2.28	0.54
1:A:372:ALA:HB3	1:A:402:LEU:HB2	1.88	0.54
1:A:641:LYS:O	1:A:709:GLU:N	2.31	0.53
1:A:299:ARG:NH1	1:A:643:GLN:OE1	2.41	0.53
1:A:655:TRP:CE2	1:A:658:TYR:HB2	2.42	0.53
1:B:43:SER:HB2	1:B:70:VAL:HB	1.90	0.53
1:B:308:THR:OG1	1:B:653:ARG:O	2.24	0.53
1:A:228:THR:HG22	1:A:247:MET:HB2	1.89	0.53
1:A:248:TRP:CD1	1:A:251:VAL:CG2	2.92	0.53
1:A:427:ALA:O	1:A:428:GLY:C	2.46	0.53
1:B:120:GLU:OE1	1:B:310:TYR:OH	2.19	0.53
1:B:642:GLY:HA3	1:B:707:PHE:O	2.08	0.53
1:B:95:LEU:N	5:B:903:HOH:O	2.41	0.53
1:A:594:HIS:HB2	1:A:645:TRP:CG	2.44	0.53
1:B:636:MET:SD	1:B:717:ILE:HG12	2.49	0.53
1:A:112:ARG:NH2	1:A:250:GLU:OE2	2.38	0.53
1:A:57:PRO:HB3	1:A:101:MET:CE	2.39	0.52
1:B:99:ILE:HD11	1:B:111:LEU:HD13	1.91	0.52
1:B:228:THR:HG22	1:B:247:MET:HB2	1.91	0.52
1:B:175:MET:CE	1:B:212:PRO:HB2	2.40	0.52
1:B:497:GLY:HA3	1:B:498:HIS:HD2	1.74	0.52
1:B:233:TYR:C	1:B:248:TRP:CH2	2.83	0.52
1:B:456:GLU:H	4:B:803:NAG:H81	1.75	0.52
1:B:252:TRP:HA	1:B:289:TYR:O	2.10	0.52
1:B:143:GLN:HB3	1:B:144:PRO:HD3	1.92	0.52
1:B:179:GLU:HG3	1:B:180:ASN:N	2.25	0.52
1:B:354:LEU:HD11	1:B:360:ALA:HB2	1.90	0.52
1:A:94:ASP:HB3	1:A:97:ARG:HB3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:438:GLU:HB3	1:B:616:LEU:HB3	1.93	0.51
1:B:630:ASP:O	1:B:694:ARG:HD3	2.11	0.51
1:A:371:ALA:N	1:A:405:CYS:SG	2.79	0.51
1:B:515:THR:HG22	1:B:658:TYR:OH	2.11	0.51
1:B:64:LYS:HD2	1:B:105:ALA:O	2.10	0.51
1:A:575:LEU:O	1:A:578:GLN:HG2	2.10	0.51
1:A:77:TRP:HZ3	1:A:156:ILE:HG13	1.76	0.51
1:B:441:PRO:HD2	1:B:469:TRP:CZ2	2.46	0.51
1:B:655:TRP:CH2	1:B:710:TRP:CZ2	2.99	0.51
1:B:495:SER:HB2	1:B:559:VAL:HG22	1.93	0.50
1:A:161:LYS:CA	1:A:169:GLN:HE22	2.21	0.50
1:A:641:LYS:HD3	1:A:709:GLU:CD	2.32	0.50
1:B:77:TRP:HB2	5:B:901:HOH:O	2.10	0.50
1:A:161:LYS:NZ	1:A:207:LEU:O	2.44	0.50
1:A:311:ASP:OD1	1:A:318:GLU:N	2.42	0.50
1:A:338:LYS:NZ	5:A:901:HOH:O	2.36	0.50
1:A:48:TYR:HE1	1:A:74:TYR:O	1.93	0.50
1:A:99:ILE:HD11	1:A:111:LEU:HD13	1.93	0.50
1:B:53:PRO:HA	1:B:56:TRP:CD2	2.47	0.50
1:B:77:TRP:HB3	1:B:125:GLY:O	2.12	0.50
1:A:490:TYR:O	1:A:565:LEU:HD12	2.11	0.49
1:A:455:TRP:HB2	1:A:582:TYR:CE2	2.47	0.49
1:B:111:LEU:O	1:B:176:ALA:HA	2.12	0.49
1:B:323:ASN:ND2	5:B:904:HOH:O	2.44	0.49
1:B:215:MET:O	1:B:227:ASP:HA	2.12	0.49
1:A:642:GLY:HA3	1:A:707:PHE:O	2.13	0.49
1:B:438:GLU:OE2	1:B:587:LYS:N	2.36	0.49
1:A:179:GLU:OE1	1:A:196:TYR:OH	2.29	0.49
1:A:639:MET:O	1:A:656:PRO:HG2	2.13	0.49
1:A:498:HIS:CD2	1:A:542:SER:HA	2.47	0.49
1:A:259:PHE:C	1:A:655:TRP:CH2	2.86	0.49
1:B:623:PHE:CE2	1:B:719:LEU:HD11	2.48	0.49
1:A:219:GLU:O	1:A:240:ASN:HB3	2.13	0.49
1:A:359:GLU:HB2	1:A:375:SER:OG	2.13	0.49
1:B:77:TRP:HH2	1:B:81:GLU:OE1	1.86	0.49
1:A:156:ILE:O	1:A:160:MET:HG2	2.12	0.48
1:A:74:TYR:HA	1:A:112:ARG:O	2.13	0.48
1:A:248:TRP:CD1	1:A:286:PHE:HB2	2.43	0.48
1:A:72:GLU:CG	1:A:112:ARG:HD2	2.44	0.48
1:B:161:LYS:NZ	1:B:207:LEU:O	2.43	0.48
1:A:50:ARG:NH1	1:A:310:TYR:CE2	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:MET:HB3	1:B:218:GLN:CG	2.44	0.48
1:A:103:GLN:HB2	1:A:172:PRO:HG3	1.94	0.48
1:A:289:TYR:HA	1:A:290:MET:HA	1.69	0.48
1:B:77:TRP:O	1:B:81:GLU:HG2	2.14	0.48
1:B:455:TRP:CD1	4:B:803:NAG:H82	2.49	0.48
1:B:643:GLN:HB2	1:B:645:TRP:CH2	2.49	0.47
1:B:709:GLU:O	1:B:709:GLU:HG2	2.14	0.47
1:A:194:LYS:O	1:A:197:THR:OG1	2.32	0.47
1:A:43:SER:HB2	1:A:70:VAL:HB	1.95	0.47
1:B:248:TRP:HD1	1:B:286:PHE:HB2	1.77	0.47
1:A:404:ASP:O	1:A:405:CYS:HB2	2.15	0.47
1:A:641:LYS:HG3	1:A:655:TRP:HD1	1.80	0.47
2:E:1:NAG:H61	2:E:2:NAG:C7	2.44	0.47
1:A:150:GLN:OE1	1:A:203:MET:HG2	2.15	0.47
1:B:175:MET:HE1	1:B:212:PRO:HB2	1.97	0.47
1:A:341:GLU:HA	1:A:344:LEU:HD12	1.97	0.47
1:B:655:TRP:HH2	1:B:710:TRP:CZ2	2.32	0.47
1:B:48:TYR:CE1	1:B:95:LEU:HD13	2.50	0.47
1:B:502:VAL:HG21	1:B:523:TYR:CD2	2.50	0.47
1:A:610:MET:O	1:A:612:GLN:HG2	2.15	0.47
1:B:215:MET:HG3	1:B:225:VAL:HG11	1.96	0.46
1:B:182:TYR:CE2	1:B:186:GLU:HB2	2.50	0.46
1:A:186:GLU:OE1	1:A:218:GLN:NE2	2.48	0.46
1:A:252:TRP:HA	1:A:289:TYR:O	2.15	0.46
1:B:490:TYR:O	1:B:565:LEU:HA	2.15	0.46
1:B:450:THR:HG22	1:B:581:SER:CB	2.46	0.46
1:B:289:TYR:HA	1:B:290:MET:HA	1.69	0.46
1:B:482:PHE:HB2	1:B:487:LYS:O	2.14	0.46
1:B:424:MET:HE1	1:B:722:ARG:HB2	1.97	0.46
1:A:281:ASN:O	1:A:283:GLY:N	2.48	0.46
1:B:299:ARG:NH1	1:B:463:ASP:O	2.33	0.46
1:B:42:ILE:HG23	1:B:286:PHE:CE1	2.51	0.46
1:B:42:ILE:HG23	1:B:286:PHE:CZ	2.50	0.46
1:A:113:ILE:O	1:A:115:PRO:HA	2.16	0.46
1:A:96:VAL:HG12	1:A:100:LYS:HE2	1.98	0.46
1:B:497:GLY:HA3	1:B:498:HIS:CD2	2.51	0.45
1:B:459:ASN:ND2	4:B:803:NAG:H83	2.21	0.45
1:B:476:ILE:HG21	1:B:482:PHE:HZ	1.80	0.45
1:A:56:TRP:O	1:A:60:ILE:HD12	2.15	0.45
1:A:120:GLU:OE1	3:F:2:GAL:H62	2.16	0.45
1:A:639:MET:HB3	1:A:708:GLU:OE1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:GLU:OE2	1:B:127:PRO:CG	2.63	0.45
1:B:215:MET:CB	1:B:218:GLN:HG3	2.45	0.45
1:B:423:LYS:HD3	1:B:423:LYS:HA	1.80	0.45
1:A:645:TRP:HA	1:A:649:GLU:O	2.16	0.45
1:B:130:LEU:HD23	1:B:138:PHE:CE1	2.52	0.45
1:B:259:PHE:CE1	1:B:655:TRP:HB3	2.51	0.45
1:B:646:ILE:HG21	1:B:697:LEU:CD2	2.47	0.45
1:B:707:PHE:CE2	1:B:709:GLU:HB2	2.52	0.45
1:B:530:ARG:NH1	5:B:908:HOH:O	2.49	0.45
1:B:641:LYS:CD	1:B:655:TRP:HZ3	2.02	0.45
1:A:41:LEU:HB2	1:A:285:PHE:HD1	1.82	0.45
1:A:255:TRP:HB3	1:A:264:PRO:HG2	1.99	0.45
1:B:50:ARG:NH2	1:B:119:ALA:O	2.48	0.45
1:B:229:CYS:SG	1:B:248:TRP:HE3	2.40	0.45
1:A:498:HIS:HD2	1:A:542:SER:HA	1.82	0.45
1:B:480:GLU:HB2	1:B:483:LEU:HD12	1.98	0.45
1:B:488:ASP:HB2	1:B:526:ASN:HB3	1.99	0.45
1:A:503:PHE:CE2	1:A:508:LEU:HB2	2.51	0.44
1:A:655:TRP:N	1:A:656:PRO:CD	2.80	0.44
1:B:640:GLY:O	1:B:656:PRO:HD2	2.18	0.44
1:B:71:ILE:HG12	1:B:107:LEU:HD23	1.99	0.44
1:A:540:SER:OG	1:A:558:GLY:HA2	2.18	0.44
1:B:404:ASP:O	1:B:405:CYS:HB2	2.18	0.44
1:B:42:ILE:HG21	1:B:334:HIS:CD2	2.53	0.44
1:A:27:TYR:CD2	1:A:226:ILE:HD11	2.53	0.44
1:A:321:LEU:HD22	5:A:943:HOH:O	2.17	0.44
1:A:63:ALA:HB1	1:A:68:LEU:HD22	2.00	0.44
1:B:641:LYS:CB	1:B:655:TRP:CZ3	3.00	0.44
1:A:34:ILE:O	1:A:34:ILE:HG22	2.18	0.44
1:A:364:ARG:HG3	1:A:370:CYS:HB2	2.00	0.44
1:A:72:GLU:OE1	1:A:287:ASN:ND2	2.51	0.44
1:B:233:TYR:C	1:B:248:TRP:HH2	2.21	0.44
1:A:114:GLY:O	1:A:179:GLU:N	2.33	0.43
1:B:574:ASN:OD1	1:B:577:LYS:HG3	2.17	0.43
1:B:199:TRP:O	1:B:203:MET:HG2	2.18	0.43
1:B:646:ILE:HG12	1:B:651:VAL:HG21	1.99	0.43
1:A:182:TYR:HB3	1:A:218:GLN:OE1	2.18	0.43
1:A:382:SER:N	1:A:396:PRO:HG3	2.32	0.43
1:B:637:ALA:HB2	1:B:688:ARG:HG3	1.99	0.43
1:A:373:PHE:CE1	1:A:401:ILE:HG23	2.54	0.43
1:A:513:TYR:HE2	1:A:658:TYR:CE1	2.35	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:466:ASP:CG	1:B:653:ARG:HH21	2.21	0.43
1:A:143:GLN:HB3	1:A:144:PRO:HD3	1.99	0.43
1:A:77:TRP:CD1	1:A:152:PHE:CE2	3.07	0.43
1:B:56:TRP:HB2	1:B:57:PRO:HD3	2.01	0.43
1:B:42:ILE:HD12	1:B:338:LYS:NZ	2.33	0.43
1:A:333:LEU:HB2	1:A:410:TYR:CZ	2.54	0.43
1:B:491:LEU:HB2	1:B:527:VAL:HG21	2.01	0.43
1:A:77:TRP:CZ2	1:A:152:PHE:HA	2.54	0.43
1:A:103:GLN:HB2	1:A:172:PRO:CG	2.49	0.42
1:A:246:LYS:NZ	1:A:281:ASN:HB3	2.34	0.42
1:A:48:TYR:CD1	1:A:49:PRO:HD3	2.53	0.42
1:B:403:PRO:HD3	1:B:409:VAL:HG22	2.02	0.42
1:B:675:GLU:HG2	1:B:676:LYS:HG2	2.01	0.42
1:B:472:THR:HB	1:B:580:TRP:CE3	2.54	0.42
1:B:708:GLU:HG3	1:B:711:GLY:H	1.84	0.42
1:B:424:MET:HE1	1:B:722:ARG:CB	2.50	0.42
1:A:520:LYS:HB3	5:A:936:HOH:O	2.19	0.42
1:B:214:ILE:HD12	1:B:249:THR:HG21	2.00	0.42
1:B:312:TYR:OH	3:D:2:GAL:O6	2.25	0.42
1:A:120:GLU:OE2	1:A:310:TYR:OH	2.34	0.42
1:B:655:TRP:N	1:B:656:PRO:CD	2.83	0.42
1:A:441:PRO:HD2	1:A:469:TRP:CH2	2.55	0.42
1:A:497:GLY:HA2	1:A:521:LEU:HD22	2.01	0.41
1:B:493:VAL:HG22	1:B:563:VAL:HG13	2.01	0.41
1:B:48:TYR:CD1	1:B:49:PRO:HD3	2.55	0.41
1:B:501:HIS:ND1	1:B:511:THR:OG1	2.44	0.41
1:B:112:ARG:HA	1:B:177:GLN:HB3	2.02	0.41
1:B:266:ARG:HA	1:B:267:PRO:HD3	1.94	0.41
1:B:46:ILE:HG12	1:B:292:HIS:HB3	2.02	0.41
1:B:323:ASN:HA	1:B:724:ARG:NH2	2.35	0.41
1:A:443:ALA:O	1:A:535:LYS:NZ	2.53	0.41
1:B:251:VAL:HB	1:B:288:TYR:HA	2.02	0.41
1:B:77:TRP:O	1:B:77:TRP:HE3	2.03	0.41
1:A:318:GLU:O	1:A:318:GLU:HG2	2.20	0.41
1:A:58:ASP:OD2	1:A:62:LYS:NZ	2.46	0.41
1:A:630:ASP:O	1:A:694:ARG:NH2	2.52	0.41
1:B:248:TRP:O	1:B:286:PHE:HA	2.21	0.41
1:A:182:TYR:CE2	1:A:186:GLU:HB2	2.54	0.41
1:A:404:ASP:CG	1:A:407:THR:HG22	2.39	0.41
1:B:462:ARG:HD3	1:B:462:ARG:HA	1.87	0.41
1:A:217:LYS:NZ	1:A:230:ASN:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:TRP:HD1	1:A:286:PHE:CB	2.29	0.41
1:A:28:ASP:OD2	1:A:38:ARG:NH2	2.52	0.41
1:B:290:MET:HE2	1:B:293:GLY:HA2	2.02	0.41
1:B:588:GLY:HA3	1:B:618:TRP:CE2	2.56	0.41
1:B:676:LYS:HA	1:B:679:GLN:NE2	2.35	0.41
1:A:182:TYR:OH	1:A:189:ILE:HD11	2.20	0.41
1:B:479:ASN:O	1:B:479:ASN:ND2	2.54	0.41
1:B:636:MET:HE3	1:B:639:MET:CG	2.42	0.41
1:A:223:ASP:OD2	1:A:243:TYR:OH	2.25	0.41
1:B:40:ILE:HD12	1:B:284:SER:HA	2.02	0.41
1:A:441:PRO:HD2	1:A:469:TRP:CZ2	2.57	0.40
1:A:597:SER:C	1:A:599:SER:H	2.24	0.40
1:B:49:PRO:HA	1:B:92:ARG:HB3	2.03	0.40
1:B:444:ASP:C	1:B:446:SER:H	2.24	0.40
1:B:596:LEU:HD13	1:B:703:LEU:HD22	2.03	0.40
1:B:588:GLY:HA3	1:B:618:TRP:CZ2	2.56	0.40
1:A:198:LYS:HG3	1:A:222:PRO:CG	2.52	0.40
1:A:22:GLU:N	5:A:909:HOH:O	2.55	0.40
1:A:489:PRO:CG	1:A:529:LEU:HD12	2.52	0.40
1:B:218:GLN:C	1:B:220:ASP:H	2.25	0.40
1:B:453:GLY:HA2	1:B:576:ALA:O	2.21	0.40
1:B:680:THR:HG23	5:B:943:HOH:O	2.21	0.40
1:B:117:VAL:HG22	1:B:124:GLY:O	2.21	0.40
1:B:130:LEU:HD23	1:B:138:PHE:CZ	2.57	0.40
1:A:234:CYS:CB	1:A:248:TRP:CE3	2.93	0.40
1:A:582:TYR:CD1	1:A:582:TYR:N	2.89	0.40
1:A:77:TRP:CE2	1:A:152:PHE:CG	3.09	0.40
1:B:48:TYR:N	1:B:49:PRO:CD	2.85	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	703/718 (98%)	649 (92%)	50 (7%)	4 (1%)	25	59
1	B	703/718 (98%)	647 (92%)	54 (8%)	2 (0%)	41	73
All	All	1406/1436 (98%)	1296 (92%)	104 (7%)	6 (0%)	34	69

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	405	CYS
1	A	282	ASN
1	A	428	GLY
1	A	389	ASN
1	B	440	THR
1	B	486	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	590/598 (99%)	586 (99%)	4 (1%)	84	93
1	B	590/598 (99%)	583 (99%)	7 (1%)	71	88
All	All	1180/1196 (99%)	1169 (99%)	11 (1%)	78	91

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	77	TRP
1	B	123	PHE
1	B	150	GLN
1	B	238	ARG
1	B	312	TYR
1	B	495	SER
1	B	658	TYR
1	A	123	PHE
1	A	235	GLU
1	A	658	TYR

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Mol	Chain	Res	Type
1	A	721	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	498	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

8 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	1,2	14,14,15	0.46	0	17,19,21	0.94	0
2	NAG	C	2	2	14,14,15	0.45	0	17,19,21	1.40	3 (17%)
3	GAL	D	1	3	12,12,12	0.40	0	17,17,17	0.66	0
3	GAL	D	2	3	11,11,12	0.64	0	15,15,17	1.13	2 (13%)
2	NAG	E	1	1,2	14,14,15	0.61	0	17,19,21	1.53	1 (5%)
2	NAG	E	2	2	14,14,15	0.38	0	17,19,21	1.31	1 (5%)
3	GAL	F	1	3	12,12,12	0.49	0	17,17,17	0.77	0
3	GAL	F	2	3	11,11,12	0.57	0	15,15,17	1.08	0

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

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	1/6/23/26	0/1/1/1
3	GAL	D	1	3	-	0/2/22/22	0/1/1/1
3	GAL	D	2	3	-	1/2/19/22	0/1/1/1
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
3	GAL	F	1	3	-	1/2/22/22	0/1/1/1
3	GAL	F	2	3	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	NAG	C1-O5-C5	5.77	120.01	112.19
2	E	2	NAG	C1-O5-C5	4.52	118.32	112.19
3	D	2	GAL	C1-C2-C3	3.34	113.77	109.67
2	C	2	NAG	C4-C3-C2	-3.32	106.16	111.02
2	C	2	NAG	O5-C1-C2	-2.96	106.62	111.29
2	C	2	NAG	C1-C2-N2	2.77	115.22	110.49
3	D	2	GAL	O5-C5-C4	-2.04	105.87	110.83

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	2	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
3	D	2	GAL	O5-C5-C6-O6
3	F	2	GAL	O5-C5-C6-O6
3	F	1	GAL	C4-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 4 short contacts:

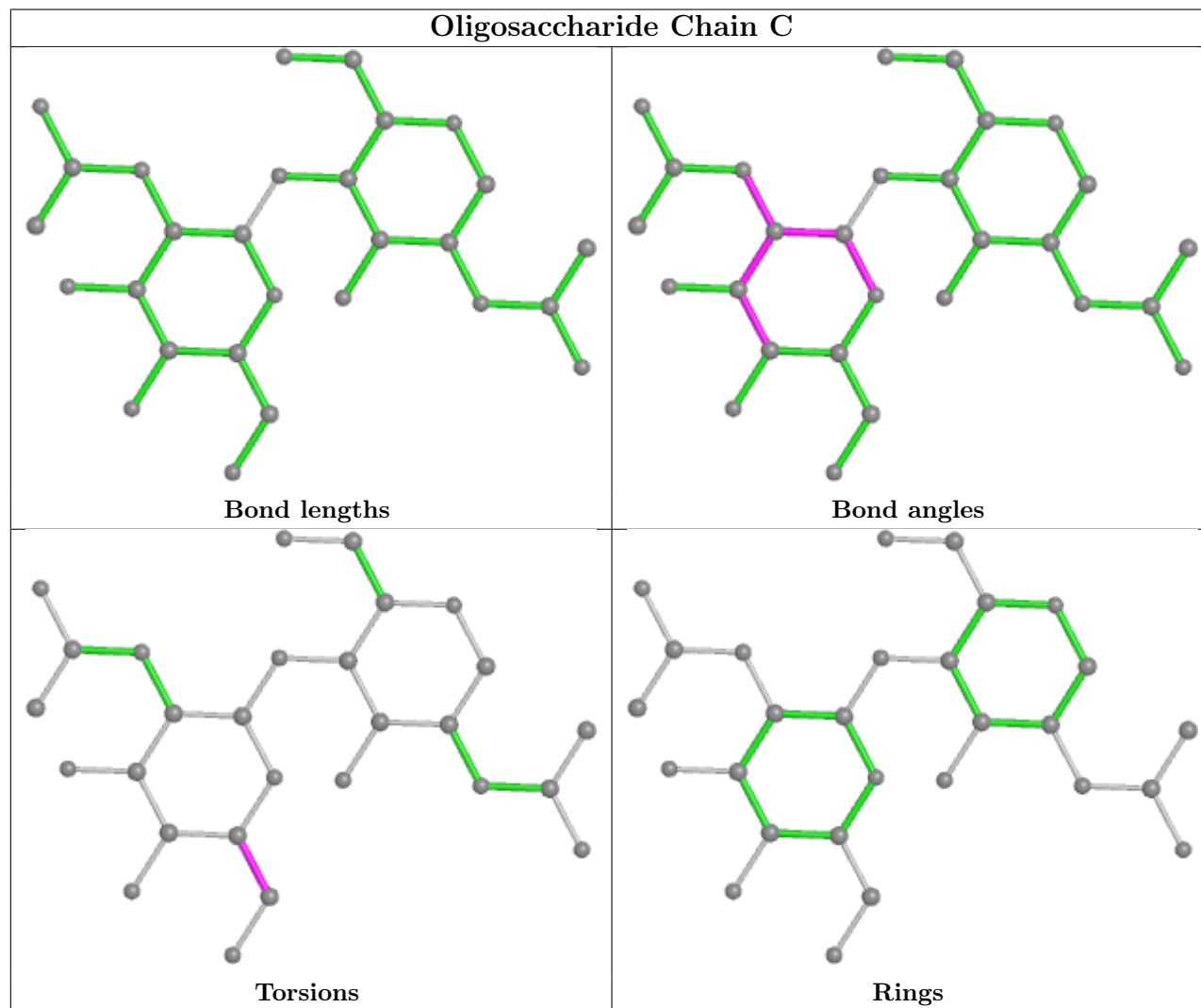
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	1	NAG	1	0
2	E	2	NAG	1	0

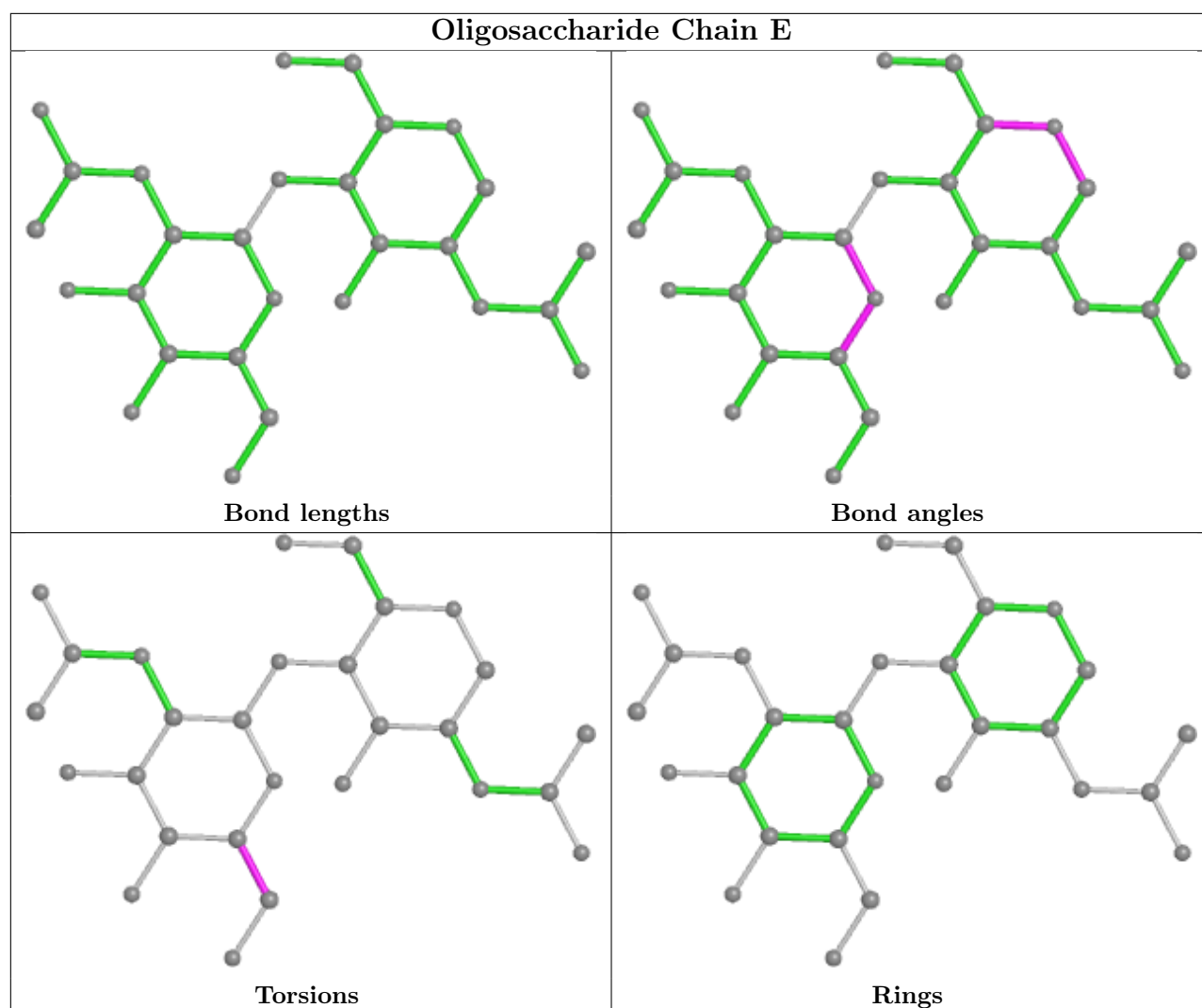
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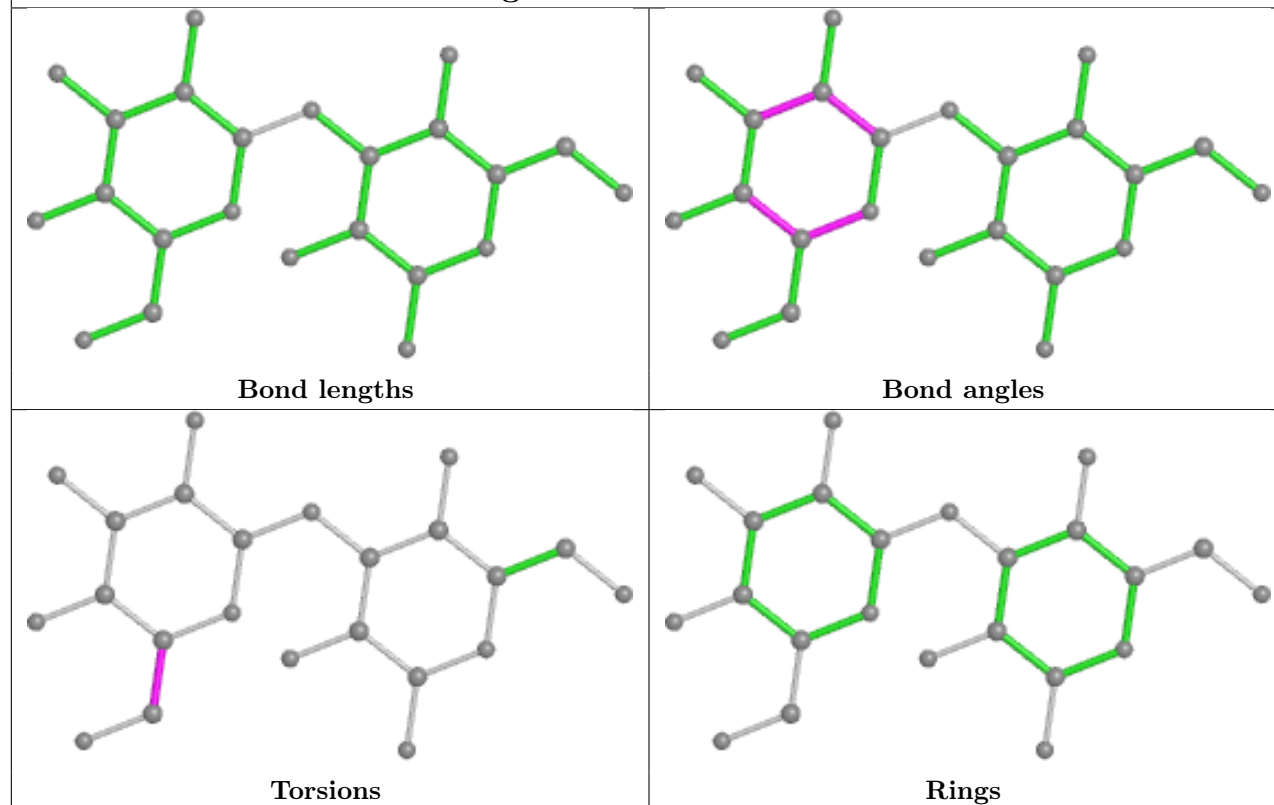
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	2	GAL	1	0
3	F	2	GAL	2	0

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

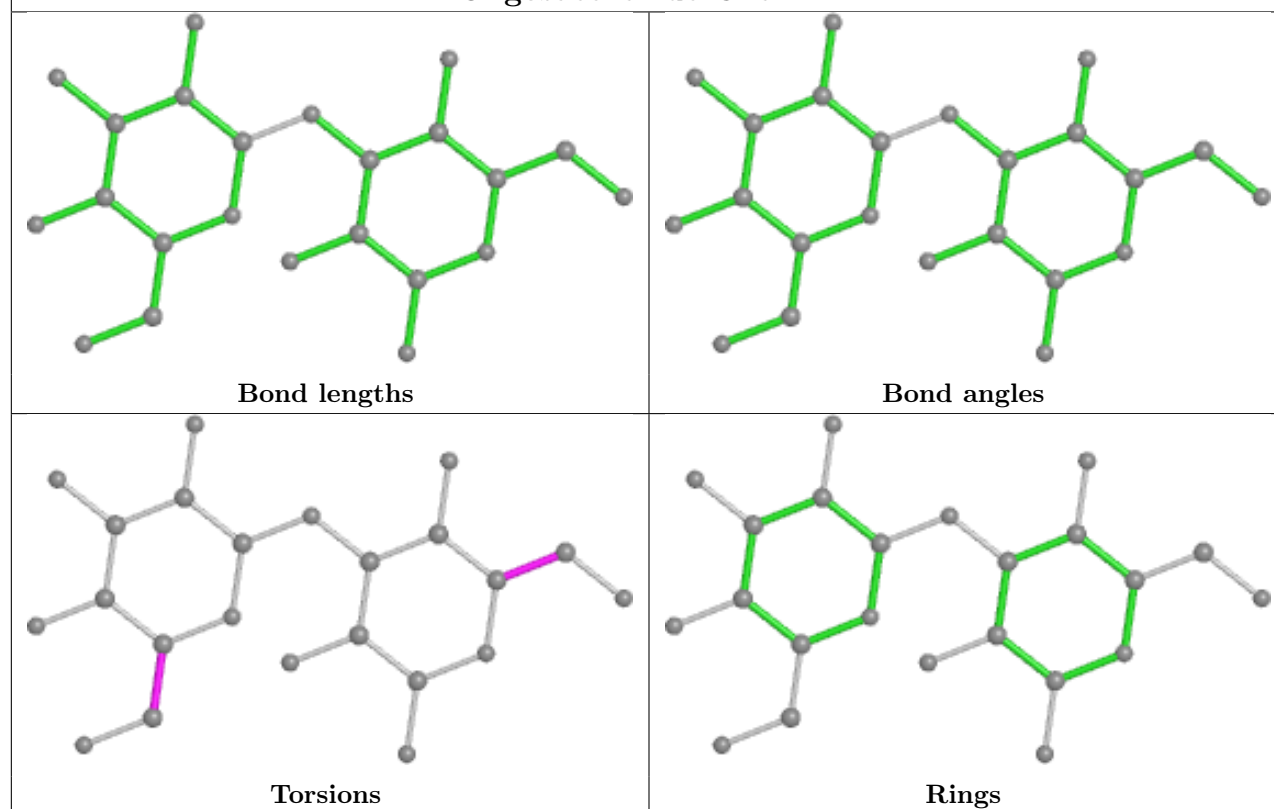




Oligosaccharide Chain D



Oligosaccharide Chain F



5.6 Ligand geometry

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	803	1	14,14,15	0.61	0	17,19,21	1.16	3 (17%)
4	NAG	B	803	1	14,14,15	0.49	0	17,19,21	1.52	3 (17%)

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
4	NAG	A	803	1	-	1/6/23/26	0/1/1/1
4	NAG	B	803	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	803	NAG	C8-C7-N2	3.47	121.98	116.10
4	B	803	NAG	O7-C7-C8	-2.79	116.87	122.06
4	B	803	NAG	C2-N2-C7	2.60	126.61	122.90
4	A	803	NAG	O5-C1-C2	-2.48	107.37	111.29
4	A	803	NAG	C4-C3-C2	-2.19	107.81	111.02
4	A	803	NAG	C1-C2-N2	2.17	114.20	110.49

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	803	NAG	O5-C5-C6-O6
4	B	803	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	B	803	NAG	C8-C7-N2-C2
4	B	803	NAG	O7-C7-N2-C2
4	A	803	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	803	NAG	4	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	705/718 (98%)	-0.28	4 (0%) 89 78	24, 45, 70, 106	0
1	B	705/718 (98%)	-0.34	10 (1%) 75 56	19, 40, 69, 124	0
All	All	1410/1436 (98%)	-0.31	14 (0%) 82 67	19, 43, 70, 124	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	600	SER	4.2
1	B	428	GLY	4.0
1	B	477	ALA	3.3
1	B	430	GLY	2.8
1	B	429	GLY	2.7
1	A	486	GLY	2.5
1	B	478	SER	2.5
1	B	603	GLU	2.5
1	A	599	SER	2.2
1	B	485	ASN	2.2
1	B	486	GLY	2.1
1	B	484	LYS	2.1
1	A	203	MET	2.0
1	B	624	ASN	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.

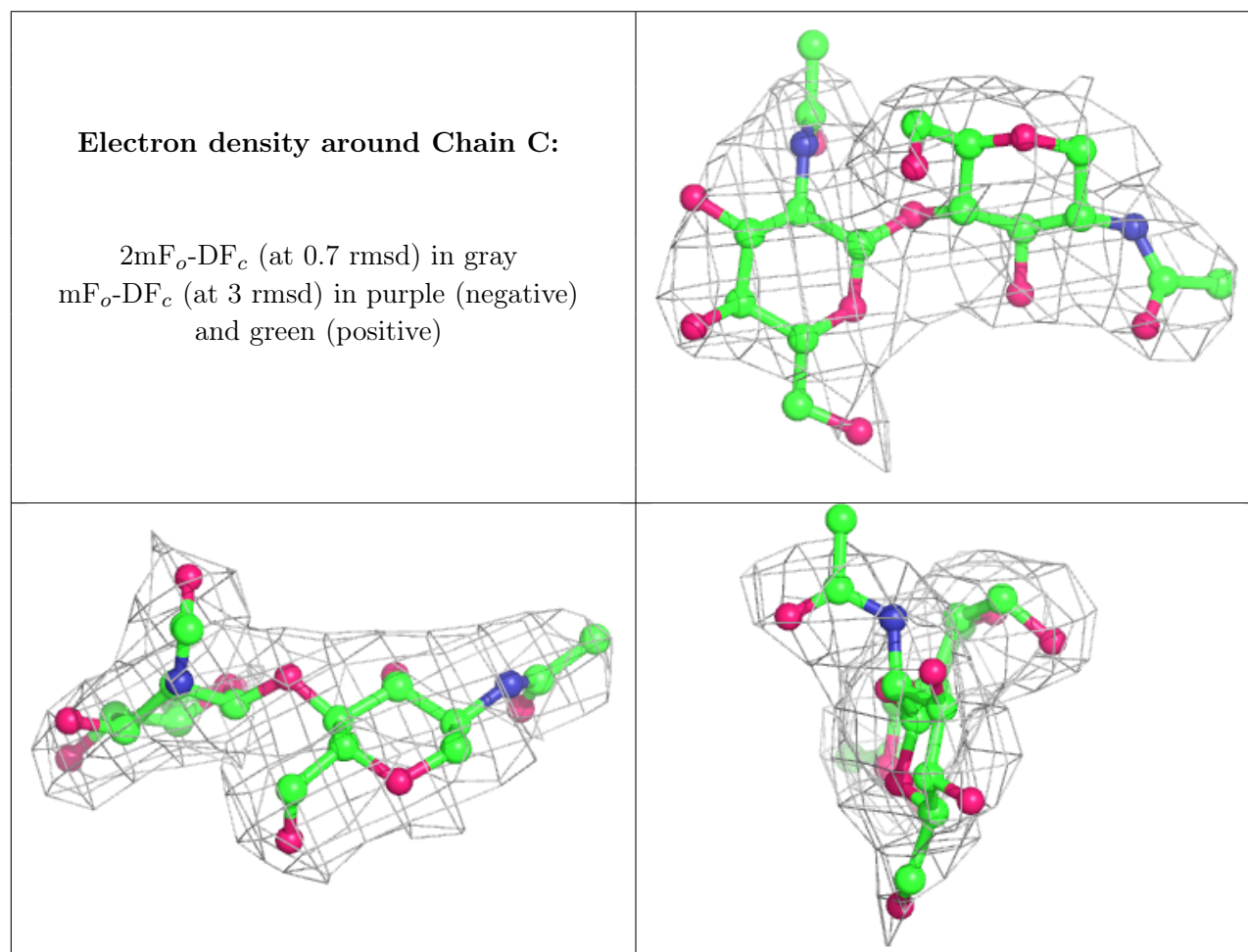
6.3 Carbohydrates [i](#)

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.

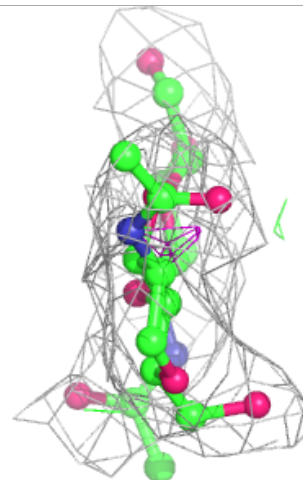
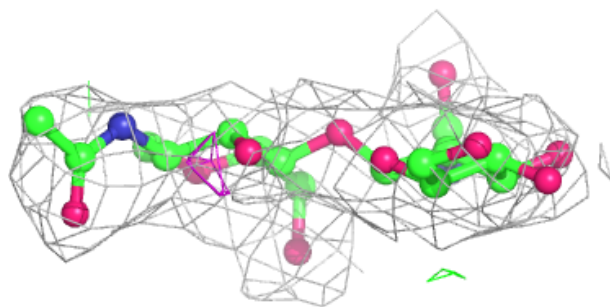
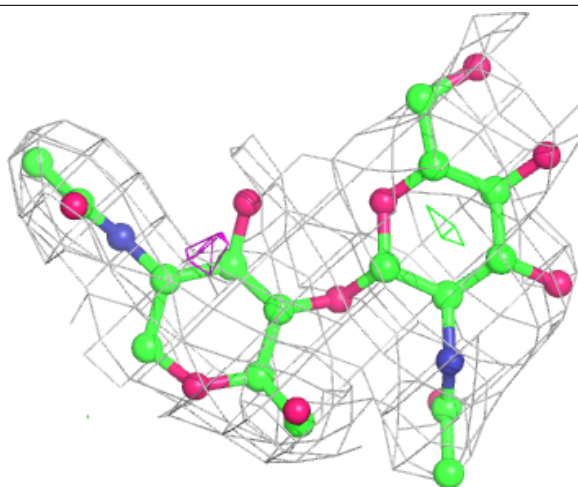
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GAL	F	1	12/12	0.81	0.26	76,86,95,101	0
2	NAG	E	2	14/15	0.82	0.23	65,74,77,80	0
3	GAL	D	1	12/12	0.84	0.28	70,83,91,95	0
2	NAG	C	2	14/15	0.88	0.35	65,70,76,77	0
2	NAG	E	1	14/15	0.89	0.20	47,56,66,66	0
3	GAL	F	2	11/12	0.90	0.20	62,65,79,81	0
3	GAL	D	2	11/12	0.94	0.18	46,51,54,58	0
2	NAG	C	1	14/15	0.97	0.15	36,40,44,55	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.



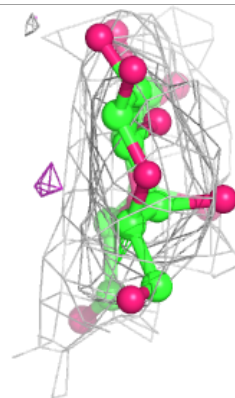
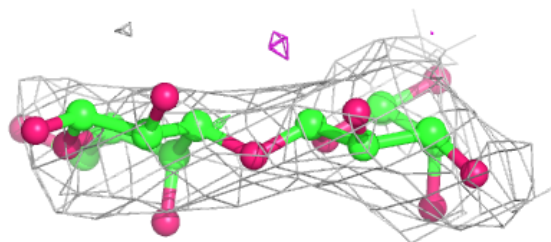
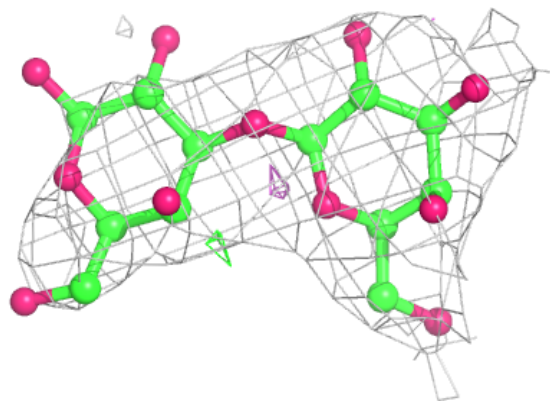
Electron density around Chain E:

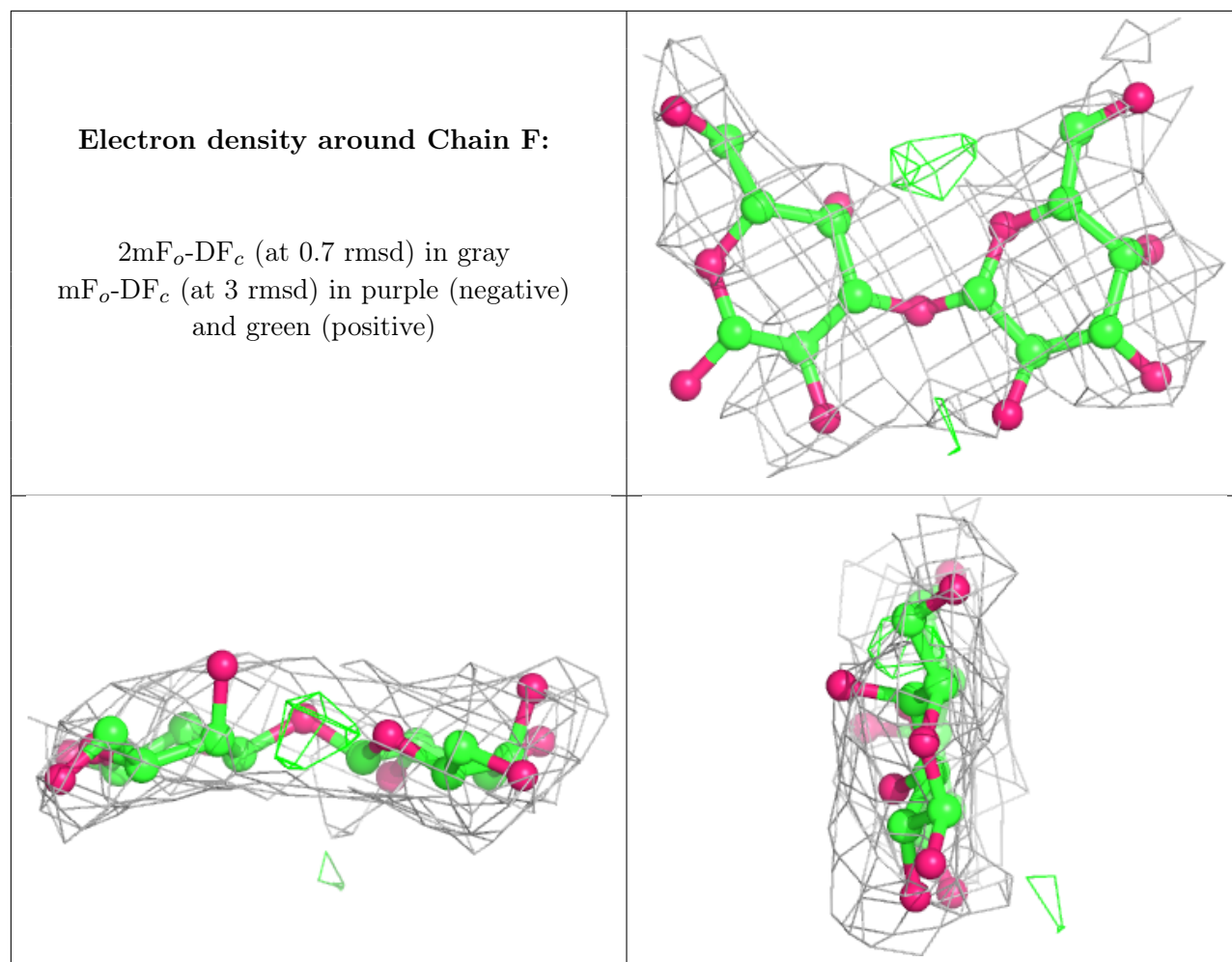
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain D:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	NAG	A	803	14/15	0.88	0.37	58,64,68,77	0
4	NAG	B	803	14/15	0.93	0.20	46,49,55,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.