



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2023 – 10:39 AM EDT

PDB ID : 8EL1
Title : Structure of MBP-Mcl-1 in complex with ABBV-467
Authors : Judge, R.A.; Judd, A.S.; Souers, A.J.
Deposited on : 2022-09-22
Resolution : 2.41 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
buster-report : 1.1.7 (2018)
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

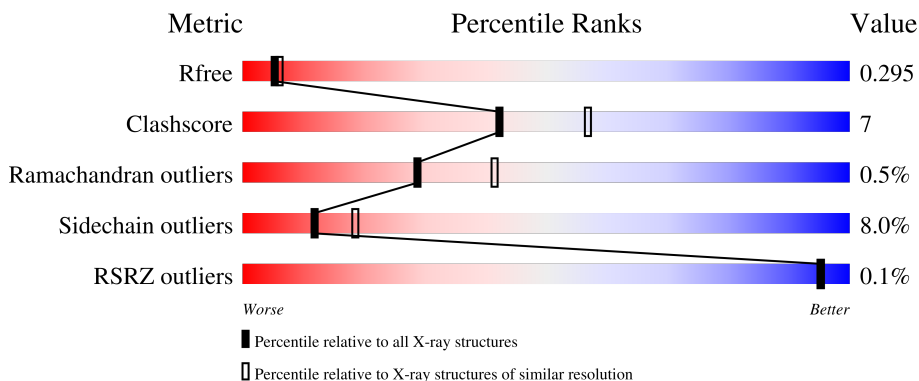
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	532	 73% 22% • •
1	B	532	 75% 20% • •
1	C	532	 71% 22% • 6%
1	D	532	 76% 17% • 7%
2	E	2	 50% 50%

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Mol	Chain	Length	Quality of chain
2	F	2	 50% 50%
2	G	2	 50% 50%
2	H	2	 50% 50%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16036 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/maltodextrin-binding periplasmic protein, Induced myeloid leukemia cell differentiation protein Mcl-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	512	Total 3936	C 2512	N 660	O 754	S 10	0	0	0
1	B	515	Total 3938	C 2511	N 659	O 758	S 10	0	0	0
1	C	499	Total 3808	C 2424	N 640	O 734	S 10	0	0	0
1	D	497	Total 3792	C 2418	N 635	O 729	S 10	0	0	0

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-210	MET	-	initiating methionine	UNP P0AEX9
A	-209	ALA	-	expression tag	UNP P0AEX9
A	-208	HIS	-	expression tag	UNP P0AEX9
A	-207	HIS	-	expression tag	UNP P0AEX9
A	-206	HIS	-	expression tag	UNP P0AEX9
A	-205	HIS	-	expression tag	UNP P0AEX9
A	-204	HIS	-	expression tag	UNP P0AEX9
A	-203	HIS	-	expression tag	UNP P0AEX9
A	-202	GLU	-	expression tag	UNP P0AEX9
A	-201	ASN	-	expression tag	UNP P0AEX9
A	-200	LEU	-	expression tag	UNP P0AEX9
A	-199	TYR	-	expression tag	UNP P0AEX9
A	-198	PHE	-	expression tag	UNP P0AEX9
A	-197	GLN	-	expression tag	UNP P0AEX9
A	-196	GLY	-	expression tag	UNP P0AEX9
A	171	GLY	-	linker	UNP P0AEX9
A	172	SER	-	linker	UNP P0AEX9
A	194	ALA	LYS	engineered mutation	UNP Q07820
A	197	ALA	LYS	engineered mutation	UNP Q07820
A	201	ALA	ARG	engineered mutation	UNP Q07820

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-210	MET	-	initiating methionine	UNP P0AEX9
B	-209	ALA	-	expression tag	UNP P0AEX9
B	-208	HIS	-	expression tag	UNP P0AEX9
B	-207	HIS	-	expression tag	UNP P0AEX9
B	-206	HIS	-	expression tag	UNP P0AEX9
B	-205	HIS	-	expression tag	UNP P0AEX9
B	-204	HIS	-	expression tag	UNP P0AEX9
B	-203	HIS	-	expression tag	UNP P0AEX9
B	-202	GLU	-	expression tag	UNP P0AEX9
B	-201	ASN	-	expression tag	UNP P0AEX9
B	-200	LEU	-	expression tag	UNP P0AEX9
B	-199	TYR	-	expression tag	UNP P0AEX9
B	-198	PHE	-	expression tag	UNP P0AEX9
B	-197	GLN	-	expression tag	UNP P0AEX9
B	-196	GLY	-	expression tag	UNP P0AEX9
B	171	GLY	-	linker	UNP P0AEX9
B	172	SER	-	linker	UNP P0AEX9
B	194	ALA	LYS	engineered mutation	UNP Q07820
B	197	ALA	LYS	engineered mutation	UNP Q07820
B	201	ALA	ARG	engineered mutation	UNP Q07820
C	-210	MET	-	initiating methionine	UNP P0AEX9
C	-209	ALA	-	expression tag	UNP P0AEX9
C	-208	HIS	-	expression tag	UNP P0AEX9
C	-207	HIS	-	expression tag	UNP P0AEX9
C	-206	HIS	-	expression tag	UNP P0AEX9
C	-205	HIS	-	expression tag	UNP P0AEX9
C	-204	HIS	-	expression tag	UNP P0AEX9
C	-203	HIS	-	expression tag	UNP P0AEX9
C	-202	GLU	-	expression tag	UNP P0AEX9
C	-201	ASN	-	expression tag	UNP P0AEX9
C	-200	LEU	-	expression tag	UNP P0AEX9
C	-199	TYR	-	expression tag	UNP P0AEX9
C	-198	PHE	-	expression tag	UNP P0AEX9
C	-197	GLN	-	expression tag	UNP P0AEX9
C	-196	GLY	-	expression tag	UNP P0AEX9
C	171	GLY	-	linker	UNP P0AEX9
C	172	SER	-	linker	UNP P0AEX9
C	194	ALA	LYS	engineered mutation	UNP Q07820
C	197	ALA	LYS	engineered mutation	UNP Q07820
C	201	ALA	ARG	engineered mutation	UNP Q07820
D	-210	MET	-	initiating methionine	UNP P0AEX9
D	-209	ALA	-	expression tag	UNP P0AEX9

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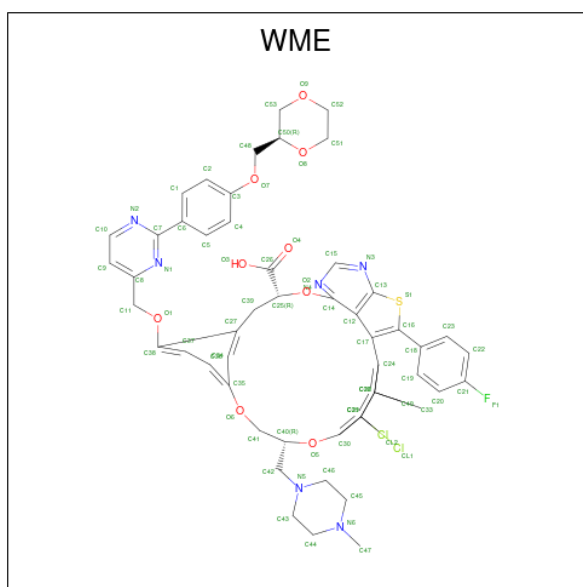
Chain	Residue	Modelled	Actual	Comment	Reference
D	-208	HIS	-	expression tag	UNP P0AEX9
D	-207	HIS	-	expression tag	UNP P0AEX9
D	-206	HIS	-	expression tag	UNP P0AEX9
D	-205	HIS	-	expression tag	UNP P0AEX9
D	-204	HIS	-	expression tag	UNP P0AEX9
D	-203	HIS	-	expression tag	UNP P0AEX9
D	-202	GLU	-	expression tag	UNP P0AEX9
D	-201	ASN	-	expression tag	UNP P0AEX9
D	-200	LEU	-	expression tag	UNP P0AEX9
D	-199	TYR	-	expression tag	UNP P0AEX9
D	-198	PHE	-	expression tag	UNP P0AEX9
D	-197	GLN	-	expression tag	UNP P0AEX9
D	-196	GLY	-	expression tag	UNP P0AEX9
D	171	GLY	-	linker	UNP P0AEX9
D	172	SER	-	linker	UNP P0AEX9
D	194	ALA	LYS	engineered mutation	UNP Q07820
D	197	ALA	LYS	engineered mutation	UNP Q07820
D	201	ALA	ARG	engineered mutation	UNP Q07820

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



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

- Molecule 3 is (7R,16R)-19,23-dichloro-10-{{2-(4-{{(2R)-1,4-dioxan-2-yl}methoxy}phenyl)pyrimidin-4-yl}methoxy}-1-(4-fluorophenyl)-20,22-dimethyl-16-[(4-methylpiperazin-1-yl)methyl]-7,8,15,16-tetrahydro-18,21-etheno-13,9-(metheno)-6,14,17-trioxa-2-thia-3,5-diazacyclonadeca[1,2,3-cd]indene-7-carboxylic acid (three-letter code: WME) (formula: C₅₃H₅₁Cl₂FN₆O₉S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	
			Total	C	Cl	F	N	O			S
3	A	1	Total	C	Cl	F	N	O	S	0	0
			72	53	2	1	6	9	1		
3	B	1	Total	C	Cl	F	N	O	S	0	0
			72	53	2	1	6	9	1		
3	C	1	Total	C	Cl	F	N	O	S	0	0
			72	53	2	1	6	9	1		
3	D	1	Total	C	Cl	F	N	O	S	0	0
			72	53	2	1	6	9	1		

- Molecule 4 is water.

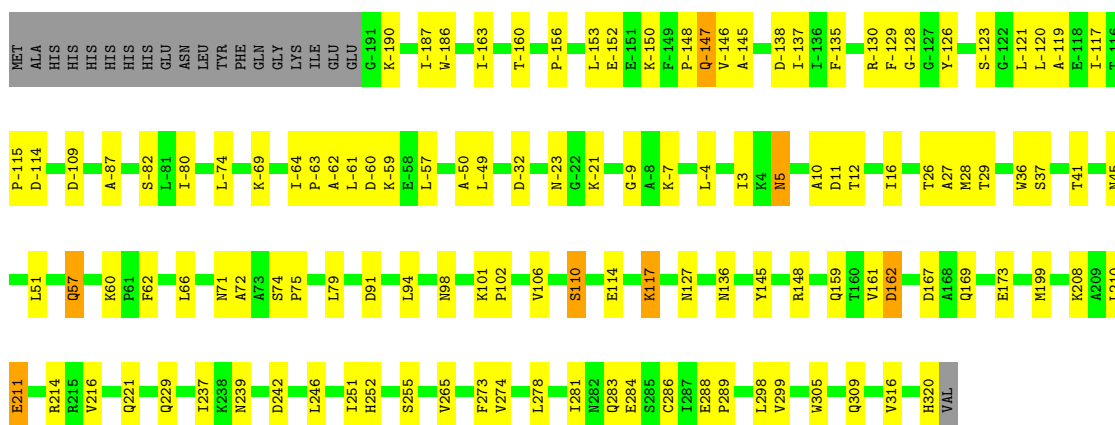
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	52	Total	O	0	0
			52	52		
4	B	42	Total	O	0	0
			42	42		
4	C	44	Total	O	0	0
			44	44		
4	D	44	Total	O	0	0
			44	44		

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.

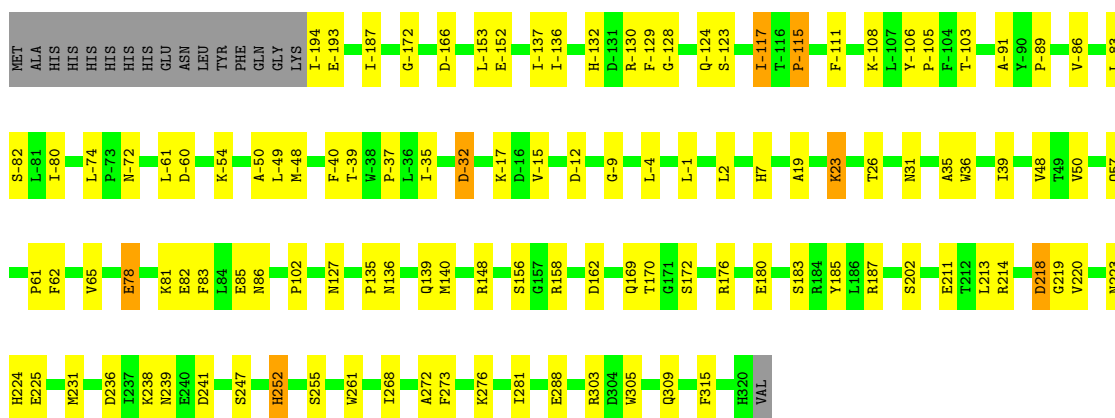
- Molecule 1: Maltose/maltodextrin-binding periplasmic protein, Induced myeloid leukemia cell differentiation protein Mcl-1

Chain A: 



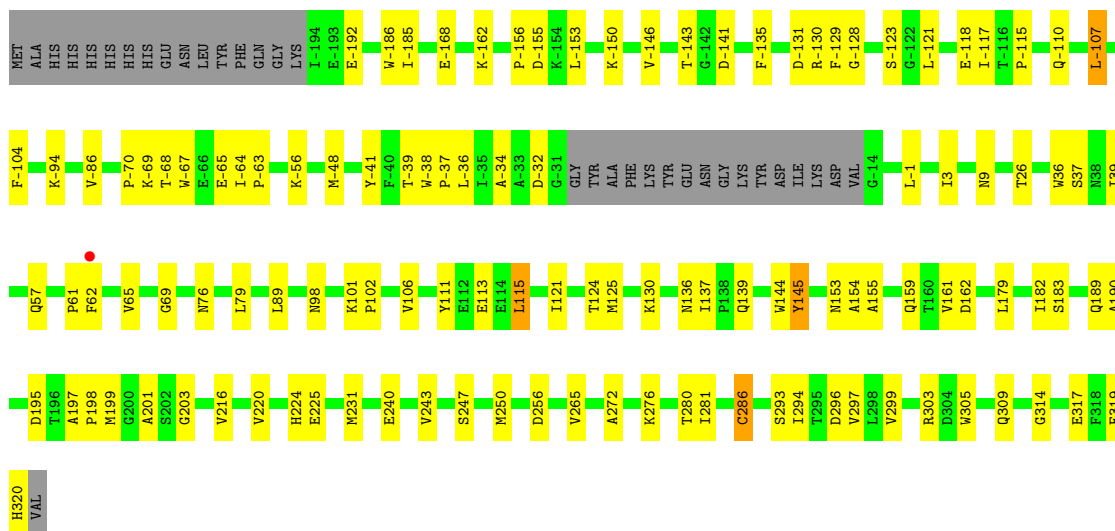
- Molecule 1: Maltose/maltodextrin-binding periplasmic protein, Induced myeloid leukemia cell differentiation protein Mcl-1

Chain B: 



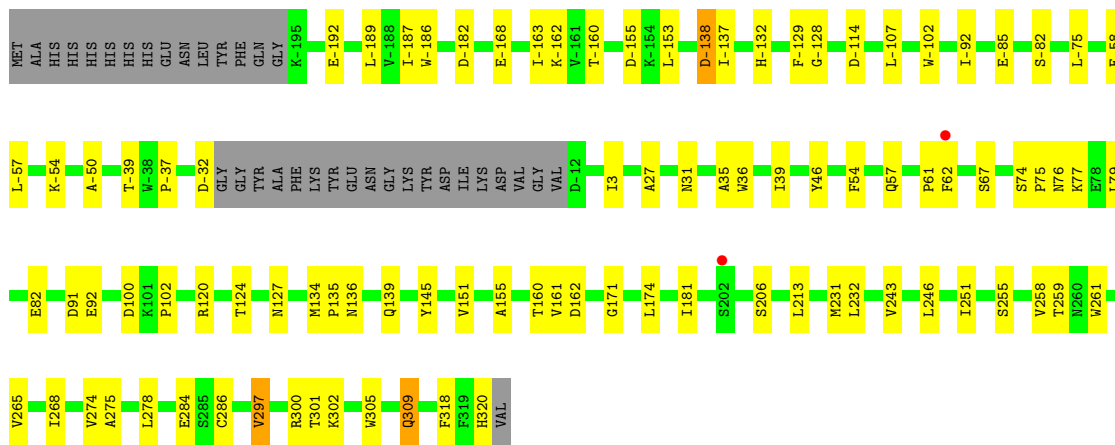
- Molecule 1: Maltose/maltodextrin-binding periplasmic protein, Induced myeloid leukemia cell differentiation protein Mcl-1

Chain C:  71% 22% 6%



- Molecule 1: Maltose/maltodextrin-binding periplasmic protein, Induced myeloid leukemia cell differentiation protein Mcl-1

Chain D:  76% 17% 7%




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

Chain E:  50% 50%



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

Chain F:  50% 50%




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

Chain G:  50% 50%

 GLC1
GLC2

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

Chain H:  50% 50%

 GLC1
GLC2

4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	111.38Å 79.62Å 137.85Å 90.00° 90.22° 90.00°	Depositor
Resolution (Å)	45.63 – 2.41 45.63 – 2.41	Depositor EDS
% Data completeness (in resolution range)	47.3 (45.63-2.41) 47.3 (45.63-2.41)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.39Å)	Xtrriage
Refinement program	BUSTER 2.11.8	Depositor
R, R_{free}	0.197 , 0.291 0.200 , 0.295	Depositor DCC
R_{free} test set	2162 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	52.2	Xtrriage
Anisotropy	0.088	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	0.120 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16036	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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: WME, 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	0.46	0/4026	0.62	0/5472
1	B	0.44	0/4028	0.61	0/5482
1	C	0.42	0/3893	0.61	0/5298
1	D	0.44	0/3877	0.62	0/5279
All	All	0.44	0/15824	0.62	0/21531

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	3936	0	3822	59	0
1	B	3938	0	3788	51	0
1	C	3808	0	3670	61	0
1	D	3792	0	3651	39	0
2	E	23	0	21	0	0
2	F	23	0	21	1	0
2	G	23	0	21	1	0
2	H	23	0	21	1	0
3	A	72	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	72	0	0	0	0
3	C	72	0	0	0	0
3	D	72	0	0	1	0
4	A	52	0	0	0	0
4	B	42	0	0	0	0
4	C	44	0	0	0	0
4	D	44	0	0	0	0
All	All	16036	0	15015	212	0

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

All (212) 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:-62:ALA:HA	1:A:-59:LYS:HD2	1.42	1.00
1:D:-187:ILE:HG12	1:D:-137:ILE:HB	1.52	0.89
1:B:-111:PHE:HA	1:B:-108:LYS:HE3	1.55	0.87
1:A:114:GLU:HA	1:A:117:LYS:HE3	1.59	0.84
1:C:37:SER:HB3	1:C:101:LYS:NZ	1.93	0.82
1:B:-12:ASP:HB2	1:B:169:GLN:HB2	1.69	0.74
1:D:-128:GLY:HA3	1:D:136:ASN:O	1.86	0.74
1:C:-128:GLY:HA3	1:C:136:ASN:O	1.89	0.73
1:C:-130:ARG:HH22	1:C:145:TYR:HE2	1.35	0.73
1:A:51:LEU:H	1:A:127:ASN:HD21	1.36	0.73
1:A:114:GLU:HA	1:A:117:LYS:CE	2.19	0.72
1:C:199:MET:HB2	1:C:203:GLY:HA3	1.70	0.72
1:D:-32:ASP:OD1	1:D:57:GLN:NE2	2.22	0.71
1:C:-156:PRO:HB3	1:C:-150:LYS:NZ	2.07	0.68
1:B:-117:ILE:HD12	1:B:-115:PRO:HD3	1.77	0.66
1:A:36:TRP:HB2	1:A:102:PRO:HG2	1.78	0.66
1:B:-128:GLY:HA3	1:B:136:ASN:O	1.96	0.64
1:C:-131:ASP:OD2	1:C:144:TRP:NE1	2.26	0.64
1:C:-123:SER:HB2	1:C:-121:LEU:HG	1.77	0.64
1:B:135:PRO:O	1:B:140:MET:HG3	1.99	0.63
1:A:-187:ILE:HG12	1:A:-137:ILE:HB	1.80	0.63
1:B:82:GLU:O	1:B:86:ASN:HB2	1.98	0.63
1:D:46:TYR:OH	1:D:120:ARG:NH1	2.31	0.63
1:A:216:VAL:HG12	1:A:265:VAL:HG11	1.80	0.62
1:C:37:SER:HB3	1:C:101:LYS:HZ3	1.63	0.62
1:B:81:LYS:O	1:B:85:GLU:HG2	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-150:LYS:O	1:A:-147:GLN:HG3	2.01	0.60
1:D:-168:GLU:HG3	1:D:-162:LYS:HA	1.82	0.60
1:C:-156:PRO:HB3	1:C:-150:LYS:HZ1	1.66	0.60
1:D:-187:ILE:HD11	1:D:-137:ILE:HD12	1.82	0.60
1:D:74:SER:O	1:D:77:LYS:HG2	2.02	0.60
1:B:36:TRP:HB2	1:B:102:PRO:HG2	1.84	0.59
1:B:252:HIS:HA	1:B:255:SER:HB2	1.84	0.59
1:A:114:GLU:HA	1:A:117:LYS:NZ	2.18	0.59
1:A:-152:GLU:HB2	1:A:-130:ARG:HD2	1.84	0.59
1:A:-123:SER:HB2	1:A:-121:LEU:HG	1.85	0.59
1:B:239:ASN:HD22	1:B:241:ASP:HB2	1.67	0.59
1:B:-17:LYS:HA	1:B:176:ARG:HG3	1.85	0.59
1:C:243:VAL:HG11	1:C:286:CYS:HB2	1.86	0.58
1:D:-107:LEU:HB2	1:D:-102:TRP:HE1	1.68	0.58
1:A:-7:LYS:HE2	1:A:162:ASP:HA	1.86	0.58
1:D:261:TRP:HB2	1:D:318:PHE:CD1	2.37	0.58
1:A:91:ASP:OD1	1:A:110:SER:HB2	2.02	0.58
1:A:-74:LEU:HD13	1:A:27:ALA:HB1	1.86	0.57
1:B:-48:MET:HB2	1:B:26:THR:HG21	1.85	0.57
1:C:37:SER:HB3	1:C:101:LYS:HZ2	1.66	0.57
1:A:286:CYS:C	1:A:289:PRO:HD2	2.25	0.57
1:C:137:ILE:HD12	1:C:139:GLN:HE21	1.69	0.57
1:D:-82:SER:OG	1:D:127:ASN:ND2	2.38	0.57
1:C:-32:ASP:OD1	1:C:57:GLN:NE2	2.38	0.57
1:C:190:ALA:HA	1:C:276:LYS:HG3	1.87	0.57
1:D:-186:TRP:HB3	1:D:-153:LEU:HD11	1.87	0.57
1:D:258:VAL:HG12	1:D:259:THR:H	1.70	0.56
1:B:305:TRP:O	1:B:309:GLN:HG2	2.05	0.56
1:A:237:ILE:HA	1:A:242:ASP:HB3	1.86	0.56
1:D:35:ALA:O	1:D:39:ILE:HG13	2.06	0.56
1:D:258:VAL:HG12	1:D:259:THR:N	2.19	0.56
1:A:-119:ALA:HB2	1:A:72:ALA:HA	1.88	0.56
2:H:1:GLC:H62	2:H:2:GLC:H5	1.88	0.55
1:C:125:MET:HE2	1:C:125:MET:HA	1.89	0.55
1:D:251:ILE:HG23	1:D:297:VAL:HG13	1.89	0.55
1:D:259:THR:HG22	1:D:305:TRP:CE2	2.42	0.54
1:A:-80:ILE:HB	1:A:29:THR:HG22	1.89	0.54
1:D:243:VAL:HG11	1:D:286:CYS:HB3	1.89	0.54
1:B:-103:THR:HB	1:B:-89:PRO:CB	2.37	0.54
1:A:145:TYR:CE2	1:A:148:ARG:NH2	2.76	0.53
1:C:115:LEU:HD22	1:C:121:ILE:HG13	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:216:VAL:HG21	1:C:319:PHE:CG	2.44	0.53
1:D:171:GLY:HA2	1:D:174:LEU:HD12	1.89	0.53
1:B:220:VAL:O	1:B:224:HIS:HB2	2.09	0.53
1:D:-192:GLU:HG3	1:D:75:PRO:HB2	1.91	0.53
1:B:-194:ILE:HG22	1:B:-193:GLU:H	1.73	0.52
1:C:154:ALA:HB1	1:C:161:VAL:HG22	1.91	0.52
1:C:220:VAL:O	1:C:224:HIS:HB2	2.09	0.52
1:D:259:THR:HG22	1:D:305:TRP:CD2	2.44	0.52
1:A:-148:PRO:HG3	1:A:-126:TYR:HE1	1.74	0.52
1:A:-128:GLY:HA3	1:A:136:ASN:O	2.09	0.52
1:B:214:ARG:O	1:B:218:ASP:OD1	2.26	0.52
1:D:-189:LEU:HD23	1:D:-138:ASP:HB2	1.90	0.52
1:B:-106:TYR:O	1:B:-103:THR:OG1	2.26	0.52
1:C:153:ASN:HB3	1:C:159:GLN:HB2	1.91	0.52
1:C:125:MET:HA	1:C:125:MET:CE	2.40	0.52
1:B:35:ALA:O	1:B:39:ILE:HG13	2.11	0.51
1:D:-37:PRO:HG3	1:D:61:PRO:HA	1.91	0.51
2:F:1:GLC:H62	2:F:2:GLC:H5	1.92	0.51
1:A:-64:ILE:N	1:A:-63:PRO:HD2	2.25	0.51
1:A:145:TYR:HE2	1:A:148:ARG:NH2	2.09	0.50
1:C:-67:TRP:CE2	1:C:-36:LEU:HD13	2.46	0.50
1:C:106:VAL:HG21	1:C:111:TYR:HD2	1.76	0.50
1:C:-64:ILE:N	1:C:-63:PRO:HD2	2.26	0.50
1:B:273:PHE:HA	1:B:276:LYS:HE3	1.92	0.50
1:B:-1:LEU:HA	1:B:2:LEU:HD12	1.94	0.50
1:D:232:LEU:HD11	1:D:274:VAL:HG22	1.93	0.50
1:B:-103:THR:HB	1:B:-89:PRO:HB3	1.94	0.50
1:B:-60:ASP:HA	1:B:-50:ALA:HB2	1.93	0.50
1:A:229:GLN:HG3	1:A:273:PHE:HZ	1.77	0.49
1:A:211:GLU:HA	1:A:214:ARG:HD2	1.93	0.49
1:A:305:TRP:O	1:A:309:GLN:HG2	2.12	0.49
1:C:36:TRP:HA	1:C:39:ILE:HD12	1.94	0.49
1:A:208:LYS:O	1:A:211:GLU:HG3	2.13	0.49
1:A:51:LEU:N	1:A:127:ASN:HD21	2.08	0.48
1:B:183:SER:O	1:B:187:ARG:HG3	2.13	0.48
1:C:37:SER:CB	1:C:101:LYS:HZ2	2.27	0.48
1:A:-156:PRO:HG2	1:A:-153:LEU:HD13	1.94	0.48
1:A:-87:ALA:HA	1:A:106:VAL:HA	1.95	0.48
1:D:231:MET:HE2	3:D:4000:WME:F1	2.04	0.48
1:B:-111:PHE:O	1:B:-108:LYS:HB2	2.12	0.48
1:B:78:GLU:H	1:B:78:GLU:HG2	1.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:ILE:HG21	1:A:10:ALA:HB2	1.94	0.48
1:C:-156:PRO:HB3	1:C:-150:LYS:HZ3	1.78	0.48
1:C:-41:TYR:HB2	2:G:2:GLC:O5	2.14	0.48
1:A:51:LEU:H	1:A:127:ASN:ND2	2.09	0.47
1:A:169:GLN:O	1:A:173:GLU:HG2	2.14	0.47
1:B:213:LEU:HD21	1:B:268:ILE:HG21	1.96	0.47
1:C:179:LEU:O	1:C:183:SER:HB3	2.14	0.47
1:B:-187:ILE:HG13	1:B:-137:ILE:HB	1.96	0.47
1:C:-37:PRO:HG3	1:C:61:PRO:HA	1.95	0.47
1:C:299:VAL:O	1:C:303:ARG:HB3	2.14	0.47
1:A:-148:PRO:HG3	1:A:-126:TYR:CE1	2.49	0.47
1:C:-48:MET:HB2	1:C:26:THR:HG21	1.97	0.47
1:C:189:GLN:HG3	1:C:272:ALA:HB1	1.97	0.47
1:A:94:LEU:O	1:A:98:ASN:HB2	2.15	0.47
1:C:-146:VAL:HB	1:C:-141:ASP:HB3	1.97	0.47
1:A:3:ILE:HD13	1:A:10:ALA:HB2	1.97	0.46
1:B:-60:ASP:OD2	1:B:7:HIS:ND1	2.48	0.46
1:B:-15:VAL:HG21	1:B:172:SER:HB3	1.97	0.46
1:D:255:SER:HA	1:D:302:LYS:NZ	2.30	0.46
1:C:-129:PHE:HE2	1:C:69:GLY:HA3	1.80	0.46
1:C:3:ILE:HD12	1:C:155:ALA:HB1	1.98	0.46
1:D:54:PHE:O	1:D:57:GLN:HG2	2.16	0.46
1:C:-118:GLU:OE2	1:C:-94:LYS:HD2	2.15	0.46
1:A:-60:ASP:HA	1:A:-50:ALA:HB2	1.98	0.46
1:B:219:GLY:O	1:B:223:ASN:OD1	2.33	0.46
1:B:-86:VAL:HG22	1:B:65:VAL:HG22	1.97	0.45
1:C:145:TYR:N	1:C:145:TYR:CD2	2.85	0.45
1:A:-49:LEU:HD13	1:A:28:MET:HG3	1.99	0.45
1:C:305:TRP:O	1:C:309:GLN:HG2	2.17	0.45
1:C:-156:PRO:CB	1:C:-150:LYS:NZ	2.77	0.45
1:C:-107:LEU:HD21	1:C:89:LEU:HD22	1.98	0.45
1:B:-32:ASP:HB3	1:B:-9:GLY:HA2	1.97	0.45
1:B:50:VAL:HA	1:B:127:ASN:HD21	1.82	0.45
1:C:76:ASN:HB3	1:C:79:LEU:HD12	1.99	0.45
1:A:57:GLN:H	1:A:57:GLN:HG2	1.62	0.45
1:C:-185:ILE:HA	1:C:-135:PHE:HB2	1.98	0.45
1:A:5:ASN:N	1:A:5:ASN:HD22	2.15	0.45
1:B:-83:LEU:CD1	1:B:-40:PHE:HA	2.47	0.45
1:A:-138:ASP:OD1	1:A:74:SER:HB2	2.17	0.44
1:C:36:TRP:HB2	1:C:102:PRO:HG2	1.99	0.44
1:C:145:TYR:N	1:C:145:TYR:HD2	2.15	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-57:LEU:HD13	1:D:-50:ALA:HA	1.98	0.44
1:A:281:ILE:O	1:A:283:GLN:OE1	2.35	0.44
1:B:-153:LEU:HD11	1:B:-136:ILE:HD11	1.99	0.44
1:B:-74:LEU:HD21	1:B:-61:LEU:HD21	1.99	0.44
1:B:-37:PRO:HG3	1:B:61:PRO:HA	2.00	0.44
1:D:259:THR:CG2	1:D:305:TRP:CD2	3.01	0.44
1:B:239:ASN:ND2	1:B:241:ASP:HB2	2.33	0.44
1:C:-86:VAL:HG22	1:C:65:VAL:HG22	2.00	0.44
1:A:208:LYS:HB3	1:A:316:VAL:HG11	2.00	0.43
1:D:-39:THR:HG21	1:D:151:VAL:HG11	2.00	0.43
1:D:134:MET:HA	1:D:135:PRO:HD3	1.92	0.43
1:C:182:ILE:HD13	1:C:294:ILE:HG22	2.00	0.43
1:A:-32:ASP:HB3	1:A:-9:GLY:HA2	2.01	0.43
1:A:159:GLN:NE2	1:A:167:ASP:OD2	2.44	0.43
1:B:156:SER:HB2	1:B:158:ARG:HG3	2.00	0.43
1:B:-152:GLU:HB2	1:B:-130:ARG:HG3	1.99	0.43
1:A:-117:ILE:HD12	1:A:-115:PRO:HG3	2.00	0.43
1:B:-129:PHE:CD2	1:B:-91:ALA:HB2	2.54	0.43
1:C:-130:ARG:NH2	1:C:145:TYR:HE2	2.08	0.43
1:D:275:ALA:HA	1:D:278:LEU:HD12	2.00	0.43
1:B:-80:ILE:HG12	1:B:48:VAL:HG22	2.01	0.42
1:A:71:ASN:HB3	1:A:74:SER:HB3	2.00	0.42
1:B:-74:LEU:HD11	1:B:-61:LEU:HD21	2.00	0.42
1:D:-129:PHE:HB3	1:D:-92:ILE:HD12	2.01	0.42
1:C:-186:TRP:HB3	1:C:-153:LEU:HD11	2.01	0.42
1:A:-4:LEU:HD23	1:A:161:VAL:HG13	2.02	0.42
1:B:-74:LEU:HD11	1:B:-61:LEU:HD11	2.00	0.42
1:D:181:ILE:HD11	1:D:206:SER:HA	2.02	0.42
1:A:274:VAL:O	1:A:278:LEU:HG	2.19	0.42
1:A:12:THR:HG23	1:A:16:ILE:HG22	2.02	0.42
1:A:74:SER:HA	1:A:75:PRO:HD3	1.92	0.42
1:B:-35:ILE:HD11	1:B:-4:LEU:HD13	2.02	0.42
1:D:36:TRP:HB2	1:D:102:PRO:HG2	2.02	0.42
1:A:-129:PHE:HD2	1:A:-120:LEU:HD11	1.85	0.41
1:C:-117:ILE:O	1:C:-115:PRO:HD3	2.20	0.41
1:A:-146:VAL:HG23	1:A:-145:ALA:N	2.35	0.41
1:B:261:TRP:CZ3	1:B:315:PHE:HB2	2.56	0.41
1:D:3:ILE:HD12	1:D:155:ALA:HB1	2.02	0.41
1:A:37:SER:OG	1:A:101:LYS:HG2	2.20	0.41
1:C:-115:PRO:HG2	1:C:-110:GLN:HE21	1.85	0.41
1:C:-39:THR:CG2	1:C:-1:LEU:HD22	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-138:ASP:OD1	1:D:76:ASN:ND2	2.52	0.41
1:C:-168:GLU:CD	1:C:-162:LYS:HD3	2.40	0.41
1:A:-64:ILE:N	1:A:-63:PRO:CD	2.84	0.41
1:C:-38:TRP:NE1	1:C:-34:ALA:HB2	2.36	0.41
1:B:185:TYR:CE2	1:B:272:ALA:HB2	2.56	0.41
1:C:-94:LYS:HA	1:C:-94:LYS:HD3	1.95	0.41
1:C:-68:THR:OG1	1:C:-65:GLU:HG2	2.21	0.41
1:C:-39:THR:HG23	1:C:-1:LEU:HD22	2.01	0.41
1:C:197:ALA:HA	1:C:198:PRO:HD3	1.97	0.41
1:C:247:SER:HA	1:C:250:MET:HE2	2.02	0.41
1:A:-64:ILE:HD13	1:A:-49:LEU:HD22	2.03	0.41
1:C:314:GLY:HA2	1:C:317:GLU:CG	2.51	0.40
1:D:305:TRP:O	1:D:309:GLN:HG2	2.21	0.40
1:B:19:ALA:O	1:B:23:LYS:HB2	2.21	0.40
1:D:-75:LEU:HD12	1:D:27:ALA:HA	2.03	0.40
1:A:-129:PHE:CD2	1:A:-120:LEU:HD11	2.56	0.40
1:B:-172:GLY:HA2	1:B:83:PHE:HE1	1.85	0.40
1:C:-117:ILE:C	1:C:-115:PRO:HD3	2.42	0.40
1:A:-186:TRP:O	1:A:-135:PHE:HB2	2.22	0.40
1:A:281:ILE:HG13	1:A:283:GLN:HG2	2.03	0.40
1:B:-106:TYR:HA	1:B:-105:PRO:HD3	2.01	0.40
1:D:213:LEU:HD21	1:D:268:ILE:HG21	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	510/532 (96%)	482 (94%)	25 (5%)	3 (1%)	25 36
1	B	513/532 (96%)	478 (93%)	32 (6%)	3 (1%)	25 36
1	C	495/532 (93%)	463 (94%)	31 (6%)	1 (0%)	47 62

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	493/532 (93%)	456 (92%)	34 (7%)	3 (1%)	25	36
All	All	2011/2128 (94%)	1879 (93%)	122 (6%)	10 (0%)	29	41

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	-163	ILE
1	A	-114	ASP
1	B	-132	HIS
1	B	-115	PRO
1	C	201	ALA
1	D	-163	ILE
1	A	11	ASP
1	B	-54	LYS
1	D	-132	HIS
1	D	100	ASP

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	401/434 (92%)	365 (91%)	36 (9%)	9	14
1	B	398/434 (92%)	367 (92%)	31 (8%)	12	19
1	C	386/434 (89%)	356 (92%)	30 (8%)	12	19
1	D	383/434 (88%)	354 (92%)	29 (8%)	13	20
All	All	1568/1736 (90%)	1442 (92%)	126 (8%)	12	18

All (126) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-190	LYS
1	A	-160	THR
1	A	-147	GLN
1	A	-109	ASP

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Mol	Chain	Res	Type
1	A	-82	SER
1	A	-69	LYS
1	A	-61	LEU
1	A	-57	LEU
1	A	-23	ASN
1	A	-21	LYS
1	A	5	ASN
1	A	26	THR
1	A	41	THR
1	A	45	ASN
1	A	57	GLN
1	A	60	LYS
1	A	62	PHE
1	A	66	LEU
1	A	79	LEU
1	A	110	SER
1	A	117	LYS
1	A	162	ASP
1	A	199	MET
1	A	210	LEU
1	A	211	GLU
1	A	221	GLN
1	A	239	ASN
1	A	246	LEU
1	A	251	ILE
1	A	252	HIS
1	A	255	SER
1	A	284	GLU
1	A	288	GLU
1	A	298	LEU
1	A	299	VAL
1	A	320	HIS
1	B	-166	ASP
1	B	-124	GLN
1	B	-123	SER
1	B	-117	ILE
1	B	-82	SER
1	B	-72	ASN
1	B	-49	LEU
1	B	-39	THR
1	B	-32	ASP
1	B	23	LYS

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Mol	Chain	Res	Type
1	B	31	ASN
1	B	57	GLN
1	B	62	PHE
1	B	78	GLU
1	B	139	GLN
1	B	148	ARG
1	B	162	ASP
1	B	170	THR
1	B	180	GLU
1	B	202	SER
1	B	211	GLU
1	B	218	ASP
1	B	225	GLU
1	B	231	MET
1	B	236	ASP
1	B	238	LYS
1	B	247	SER
1	B	252	HIS
1	B	281	ILE
1	B	288	GLU
1	B	303	ARG
1	C	-192	GLU
1	C	-155	ASP
1	C	-143	THR
1	C	-107	LEU
1	C	-104	PHE
1	C	-70	PRO
1	C	-69	LYS
1	C	-56	LYS
1	C	9	ASN
1	C	62	PHE
1	C	98	ASN
1	C	113	GLU
1	C	115	LEU
1	C	124	THR
1	C	130	LYS
1	C	145	TYR
1	C	162	ASP
1	C	195	ASP
1	C	225	GLU
1	C	231	MET
1	C	240	GLU

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Mol	Chain	Res	Type
1	C	256	ASP
1	C	265	VAL
1	C	280	THR
1	C	281	ILE
1	C	286	CYS
1	C	293	SER
1	C	296	ASP
1	C	297	VAL
1	C	320	HIS
1	D	-182	ASP
1	D	-160	THR
1	D	-155	ASP
1	D	-138	ASP
1	D	-114	ASP
1	D	-85	GLU
1	D	-58	GLU
1	D	-54	LYS
1	D	31	ASN
1	D	62	PHE
1	D	67	SER
1	D	79	LEU
1	D	82	GLU
1	D	91	ASP
1	D	92	GLU
1	D	124	THR
1	D	139	GLN
1	D	145	TYR
1	D	160	THR
1	D	161	VAL
1	D	162	ASP
1	D	246	LEU
1	D	265	VAL
1	D	284	GLU
1	D	297	VAL
1	D	300	ARG
1	D	301	THR
1	D	309	GLN
1	D	320	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	B	-110	GLN
1	B	-72	ASN
1	B	129	GLN
1	B	224	HIS
1	C	139	GLN
1	C	221	GLN
1	D	57	GLN
1	D	127	ASN
1	D	221	GLN

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](#)

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	GLC	E	1	2	12,12,12	0.26	0	17,17,17	0.61	0
2	GLC	E	2	2	11,11,12	0.31	0	15,15,17	0.87	1 (6%)
2	GLC	F	1	2	12,12,12	0.26	0	17,17,17	0.62	0
2	GLC	F	2	2	11,11,12	0.38	0	15,15,17	0.90	1 (6%)
2	GLC	G	1	2	12,12,12	0.32	0	17,17,17	0.67	0
2	GLC	G	2	2	11,11,12	0.34	0	15,15,17	0.68	1 (6%)
2	GLC	H	1	2	12,12,12	0.41	0	17,17,17	0.93	0
2	GLC	H	2	2	11,11,12	0.35	0	15,15,17	0.95	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	GLC	E	1	2	-	0/2/22/22	0/1/1/1
2	GLC	E	2	2	-	0/2/19/22	0/1/1/1
2	GLC	F	1	2	-	0/2/22/22	0/1/1/1
2	GLC	F	2	2	-	0/2/19/22	0/1/1/1
2	GLC	G	1	2	-	0/2/22/22	0/1/1/1
2	GLC	G	2	2	-	0/2/19/22	0/1/1/1
2	GLC	H	1	2	-	0/2/22/22	0/1/1/1
2	GLC	H	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	2	GLC	C1-O5-C5	2.68	115.82	112.19
2	G	2	GLC	C1-O5-C5	2.14	115.10	112.19
2	H	2	GLC	C1-O5-C5	2.11	115.05	112.19
2	F	2	GLC	C1-O5-C5	2.10	115.03	112.19

There are no chirality outliers.

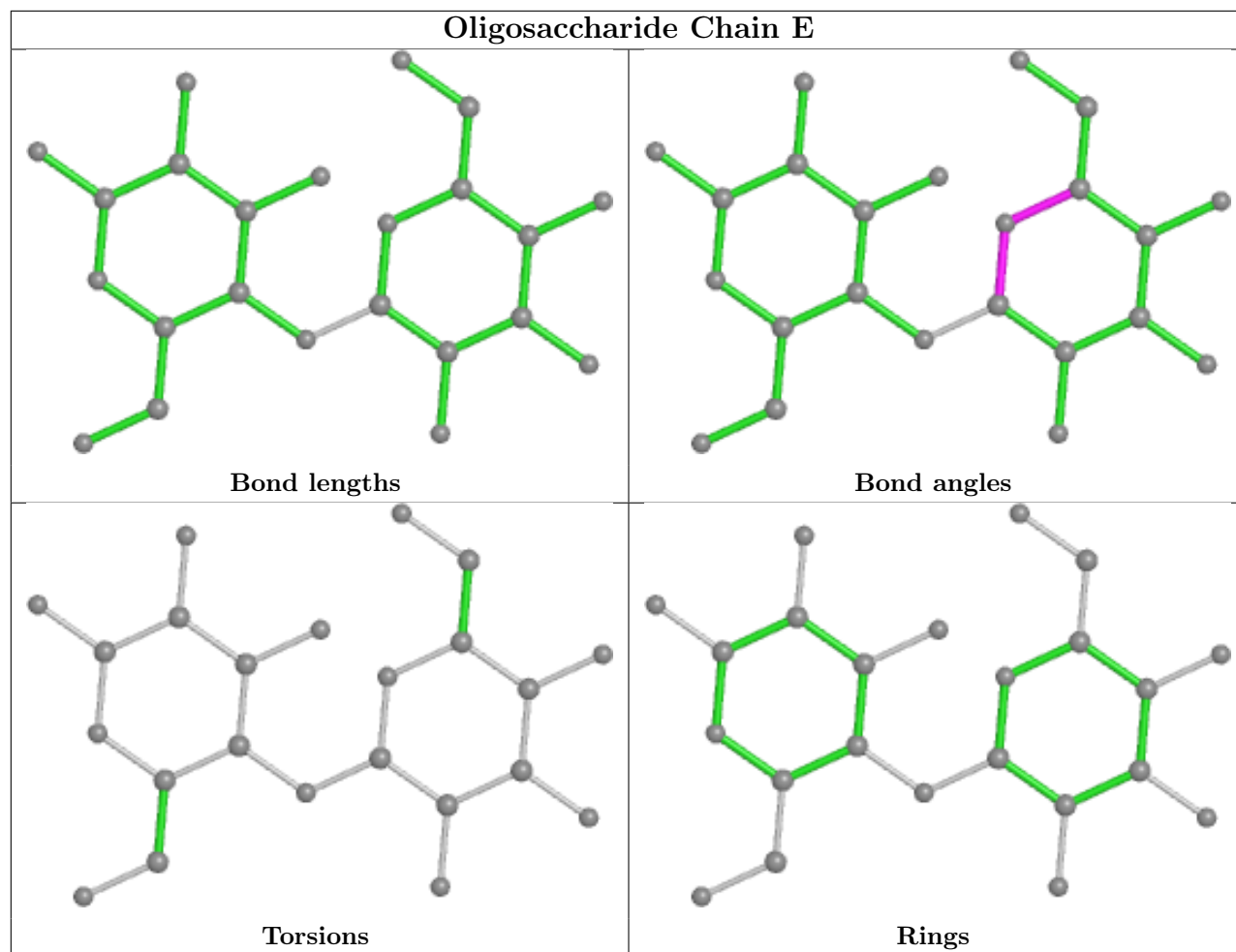
There are no torsion outliers.

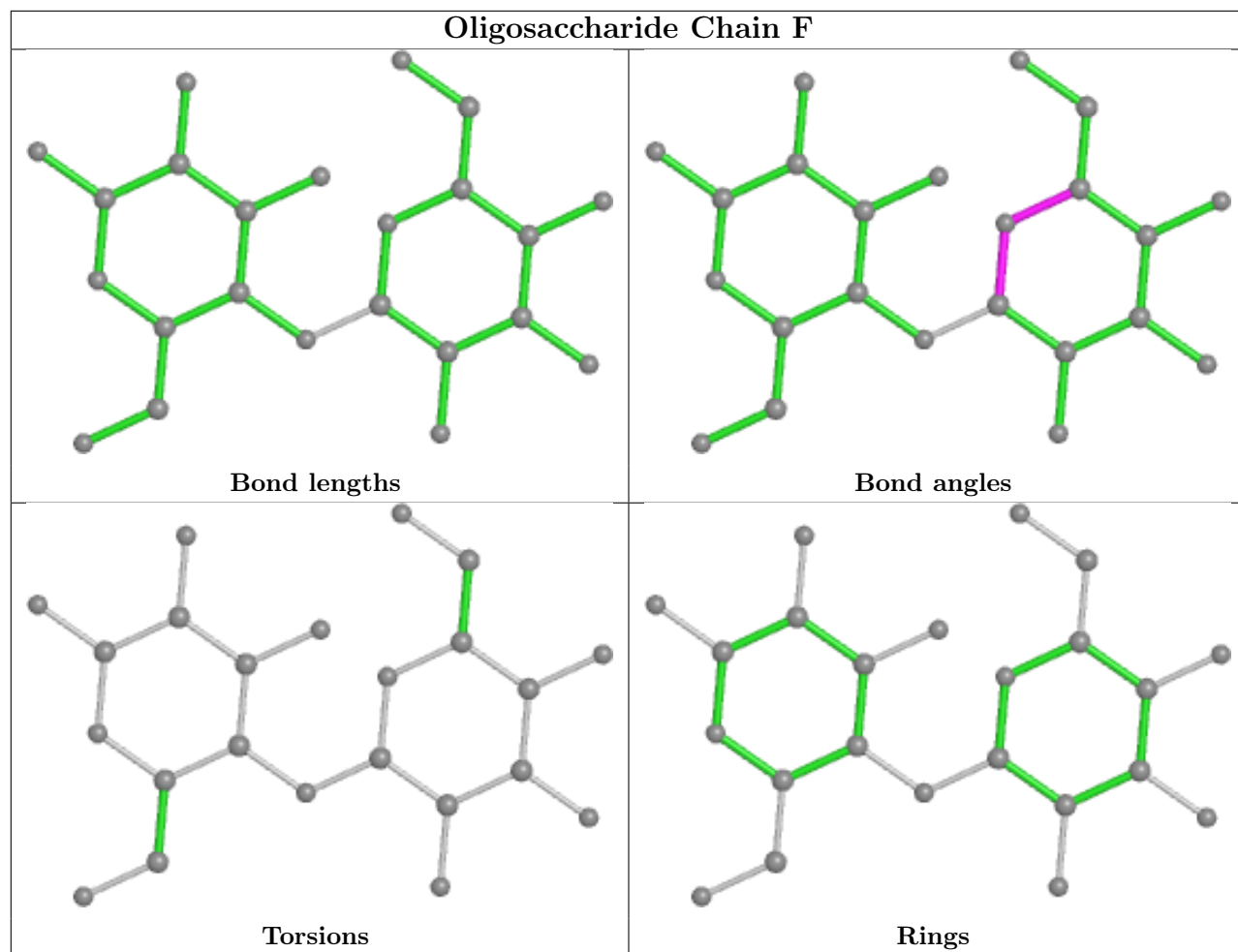
There are no ring outliers.

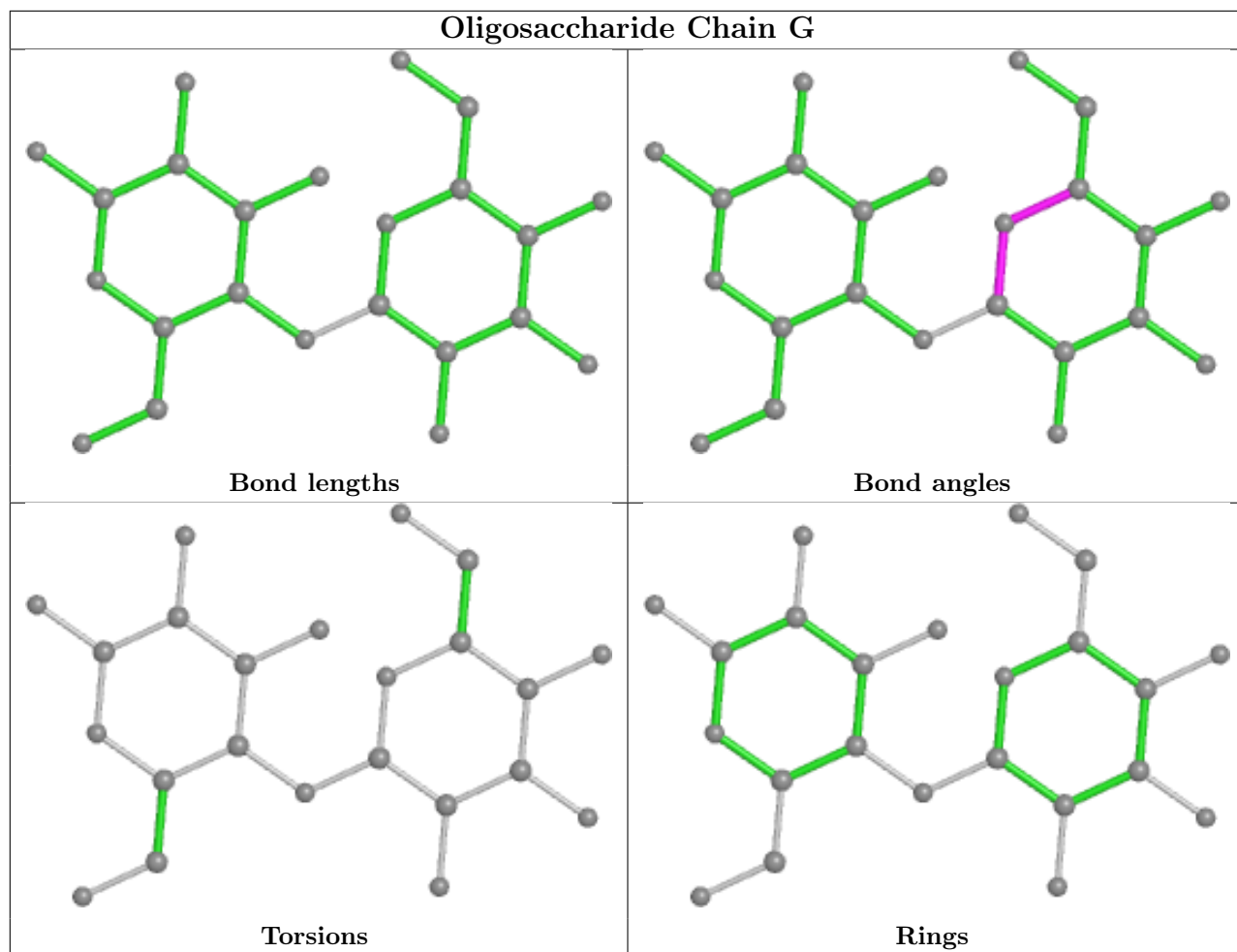
5 monomers are involved in 3 short contacts:

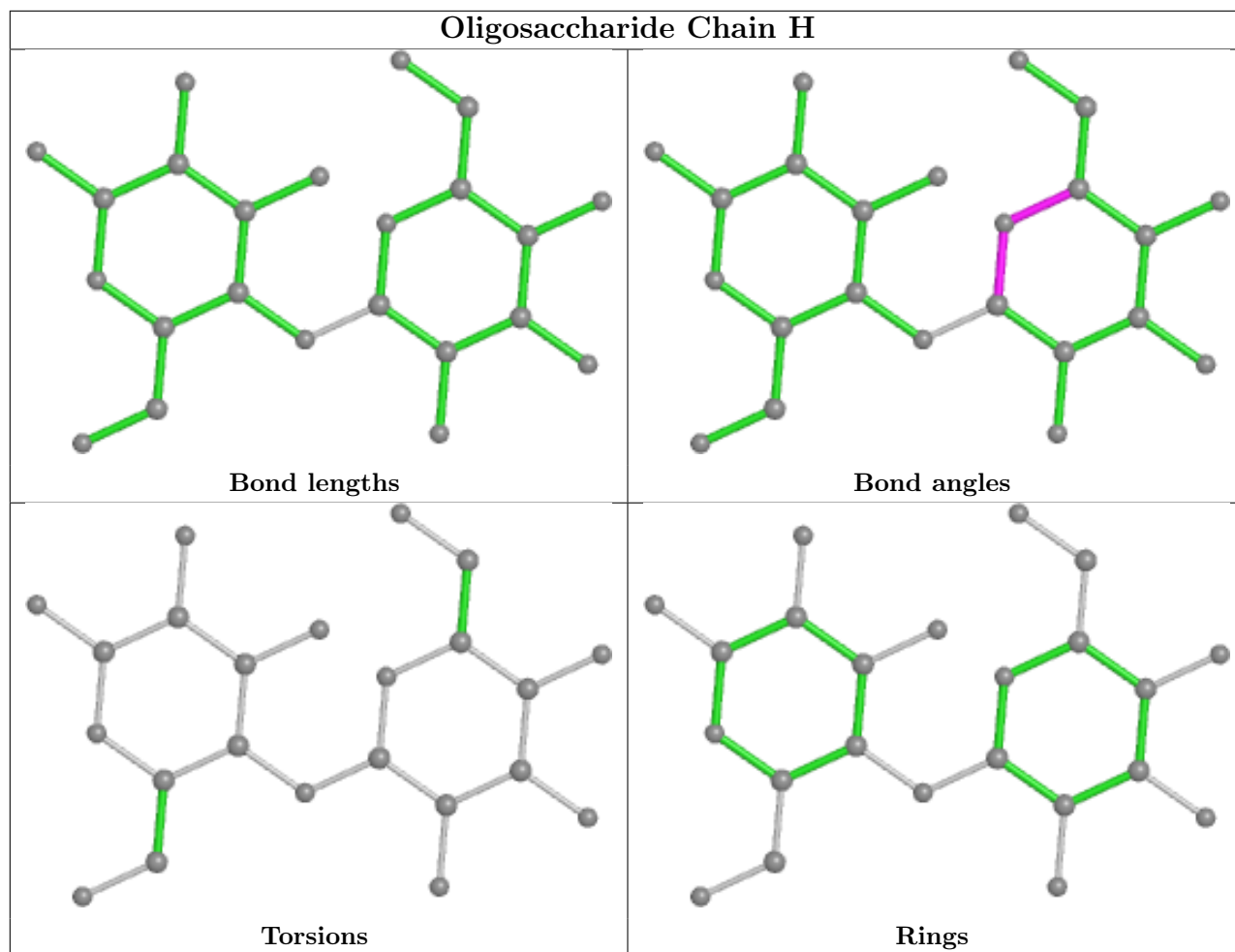
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	1	GLC	1	0
2	H	2	GLC	1	0
2	G	2	GLC	1	0
2	F	2	GLC	1	0
2	F	1	GLC	1	0

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









5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
3	WME	B	4000	-	76,81,81	0.72	3 (3%)	97,116,116	0.97	5 (5%)
3	WME	D	4000	-	76,81,81	0.71	3 (3%)	97,116,116	1.08	6 (6%)
3	WME	A	4000	-	76,81,81	0.74	3 (3%)	97,116,116	0.93	5 (5%)
3	WME	C	4000	-	76,81,81	0.69	3 (3%)	97,116,116	0.92	5 (5%)

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
3	WME	B	4000	-	-	8/47/65/65	0/9/10/10
3	WME	D	4000	-	-	7/47/65/65	0/9/10/10
3	WME	A	4000	-	-	8/47/65/65	0/9/10/10
3	WME	C	4000	-	-	10/47/65/65	0/9/10/10

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	4000	WME	O4-C26	3.68	1.33	1.22
3	B	4000	WME	O4-C26	3.23	1.32	1.22
3	C	4000	WME	O4-C26	3.18	1.31	1.22
3	D	4000	WME	O4-C26	3.08	1.31	1.22
3	D	4000	WME	C17-C12	3.05	1.47	1.42
3	A	4000	WME	C17-C12	2.97	1.47	1.42
3	C	4000	WME	C17-C12	2.78	1.47	1.42
3	B	4000	WME	C17-C12	2.66	1.46	1.42
3	D	4000	WME	O3-C26	-2.60	1.22	1.30
3	C	4000	WME	O3-C26	-2.57	1.22	1.30
3	B	4000	WME	O3-C26	-2.51	1.22	1.30
3	A	4000	WME	O3-C26	-2.28	1.23	1.30

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4000	WME	C17-C12-C13	5.58	111.99	107.54
3	A	4000	WME	C17-C12-C13	5.27	111.74	107.54
3	C	4000	WME	C17-C12-C13	5.02	111.54	107.54
3	B	4000	WME	C17-C12-C13	4.60	111.21	107.54
3	D	4000	WME	C14-O2-C25	4.42	123.14	116.79
3	B	4000	WME	O4-C26-C25	-3.90	113.37	122.57
3	D	4000	WME	O4-C26-C25	-3.62	114.02	122.57
3	B	4000	WME	O3-C26-C25	3.40	122.75	113.03
3	D	4000	WME	O3-C26-C25	3.17	122.08	113.03
3	B	4000	WME	C14-O2-C25	3.12	121.28	116.79
3	D	4000	WME	C17-C24-C28	-2.79	114.24	120.13
3	C	4000	WME	O4-C26-C25	-2.68	116.25	122.57
3	A	4000	WME	C14-O2-C25	2.63	120.57	116.79
3	A	4000	WME	C17-C24-C28	-2.59	114.65	120.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	4000	WME	O3-C26-C25	2.53	120.25	113.03
3	B	4000	WME	C17-C24-C28	-2.50	114.84	120.13
3	A	4000	WME	O3-C26-C25	2.42	119.95	113.03
3	A	4000	WME	O4-C26-C25	-2.40	116.91	122.57
3	D	4000	WME	C40-C42-N5	2.33	126.14	113.69
3	C	4000	WME	C24-C17-C16	-2.12	122.19	127.78
3	C	4000	WME	C17-C24-C28	-2.08	115.74	120.13

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	4000	WME	O5-C40-C42-N5
3	A	4000	WME	O7-C48-C50-O8
3	A	4000	WME	O7-C48-C50-C53
3	B	4000	WME	O5-C40-C42-N5
3	B	4000	WME	O7-C48-C50-O8
3	C	4000	WME	C41-C40-O5-C30
3	D	4000	WME	C41-C40-O5-C30
3	D	4000	WME	C40-C42-N5-C43
3	A	4000	WME	C41-C40-O5-C30
3	B	4000	WME	O7-C48-C50-C53
3	C	4000	WME	O7-C48-C50-C53
3	C	4000	WME	C42-C40-C41-O6
3	D	4000	WME	C42-C40-C41-O6
3	A	4000	WME	C42-C40-C41-O6
3	B	4000	WME	C41-C40-O5-C30
3	B	4000	WME	O5-C40-C41-O6
3	B	4000	WME	C42-C40-C41-O6
3	C	4000	WME	O7-C48-C50-O8
3	C	4000	WME	O5-C40-C42-N5
3	D	4000	WME	O5-C40-C42-N5
3	D	4000	WME	O5-C40-C41-O6
3	A	4000	WME	C40-C41-O6-C35
3	A	4000	WME	O5-C40-C41-O6
3	C	4000	WME	O5-C40-C41-O6
3	B	4000	WME	S1-C16-C18-C23
3	C	4000	WME	S1-C16-C18-C19
3	C	4000	WME	S1-C16-C18-C23
3	D	4000	WME	C40-C41-O6-C35
3	C	4000	WME	C40-C41-O6-C35
3	B	4000	WME	O2-C25-C26-O3

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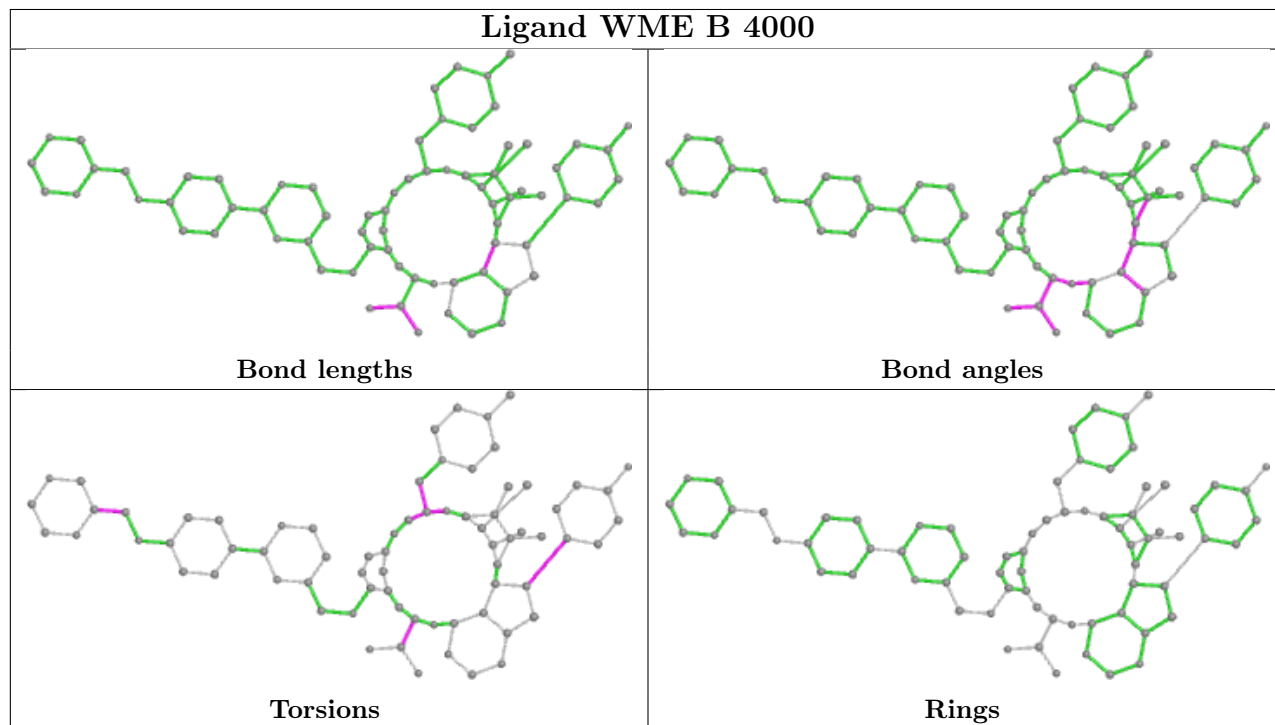
Mol	Chain	Res	Type	Atoms
3	D	4000	WME	O2-C25-C26-O3
3	C	4000	WME	C40-C42-N5-C46
3	A	4000	WME	C12-C14-O2-C25

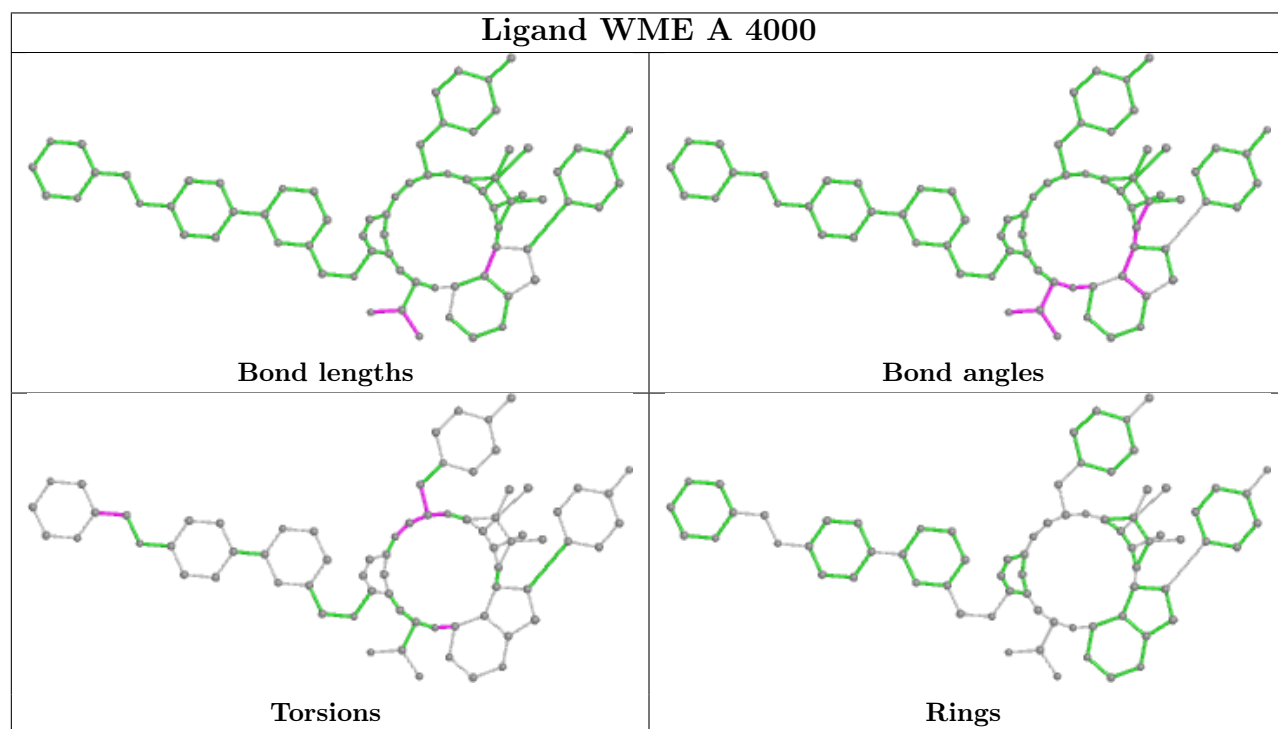
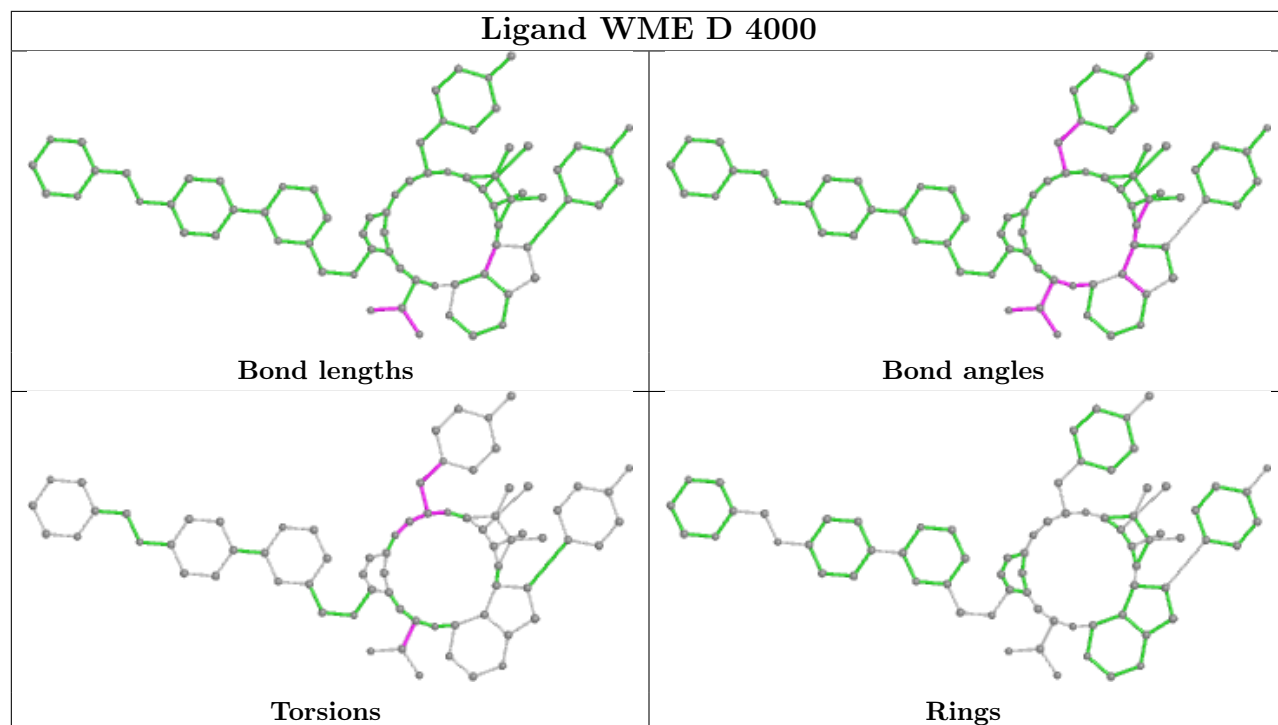
There are no ring outliers.

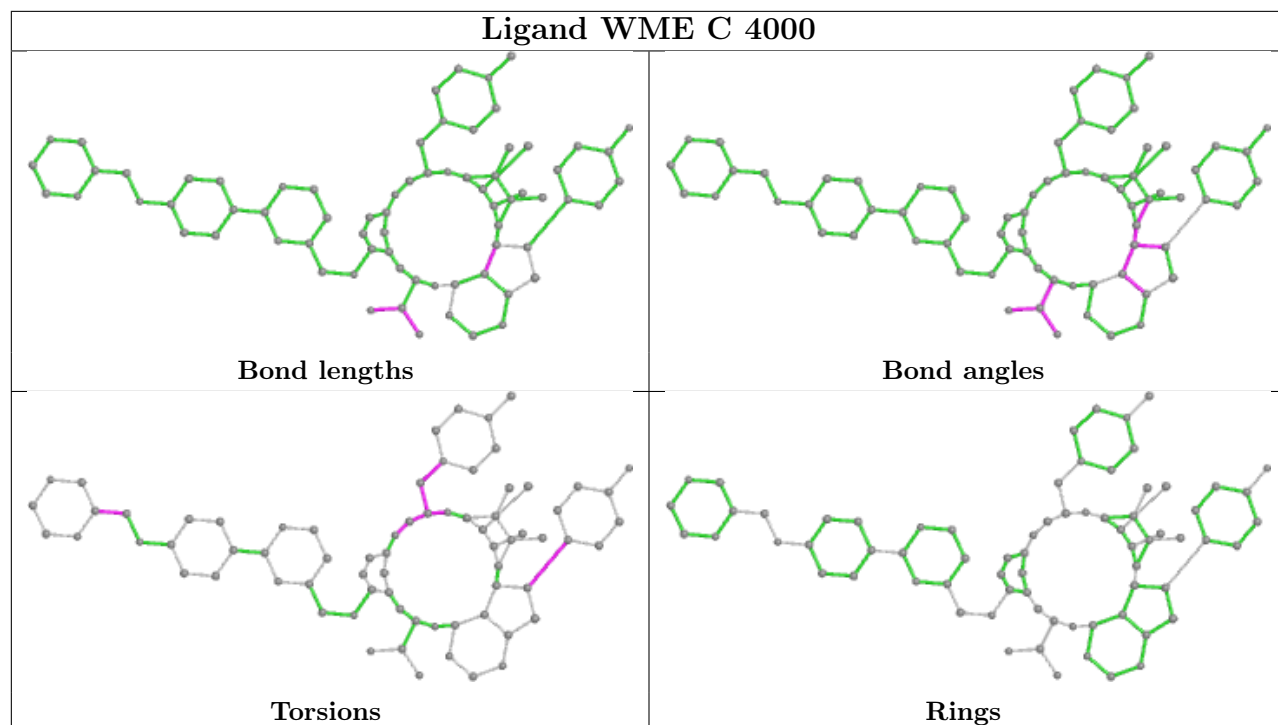
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	4000	WME	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







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	512/532 (96%)	-0.49	0 100 100	26, 46, 63, 81	0
1	B	515/532 (96%)	-0.44	0 100 100	29, 51, 71, 84	0
1	C	499/532 (93%)	-0.40	1 (0%) 95 94	33, 54, 77, 96	0
1	D	497/532 (93%)	-0.40	2 (0%) 92 91	32, 54, 77, 86	0
All	All	2023/2128 (95%)	-0.43	3 (0%) 95 95	26, 51, 74, 96	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	202	SER	3.0
1	D	62	PHE	2.7
1	C	62	PHE	2.2

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.

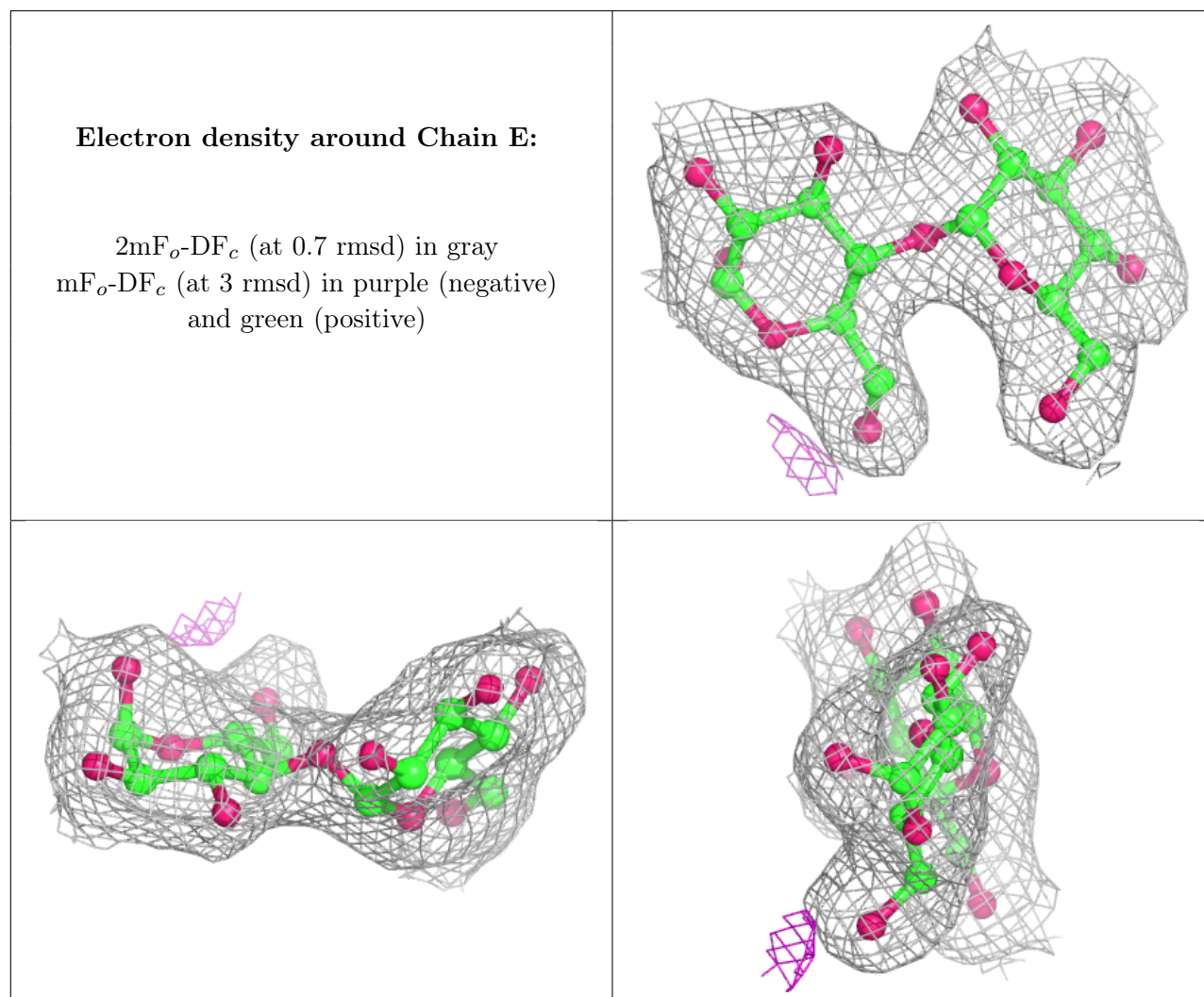
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	F	2	11/12	0.95	0.12	41,42,43,44	0
2	GLC	H	1	12/12	0.95	0.17	44,46,46,47	0
2	GLC	E	1	12/12	0.96	0.11	38,41,41,41	0
2	GLC	G	1	12/12	0.96	0.10	38,39,39,39	0

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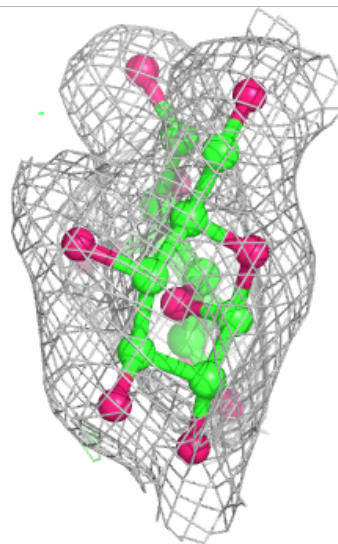
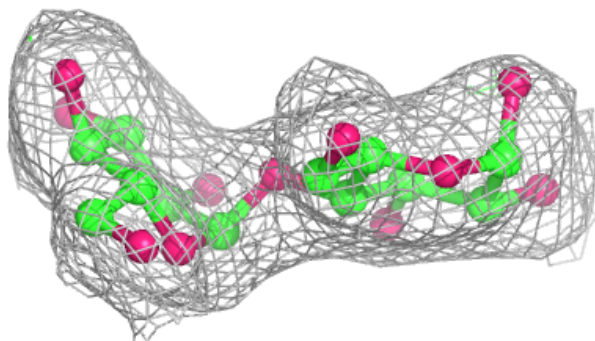
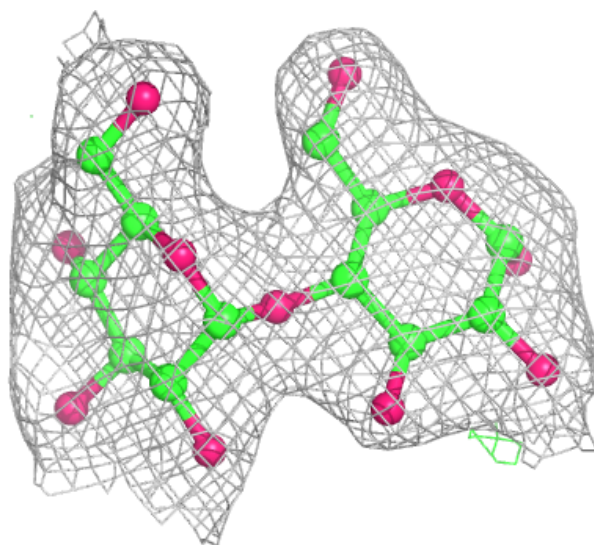
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GLC	F	1	12/12	0.96	0.12	41,42,43,43	0
2	GLC	H	2	11/12	0.96	0.11	45,45,45,46	0
2	GLC	E	2	11/12	0.97	0.11	34,35,37,37	0
2	GLC	G	2	11/12	0.98	0.10	36,38,39,40	0

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



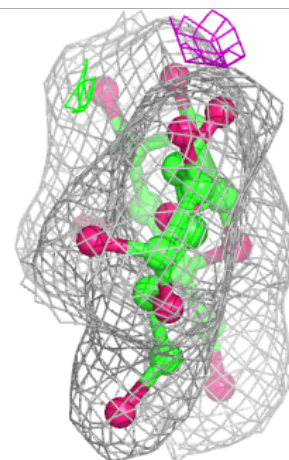
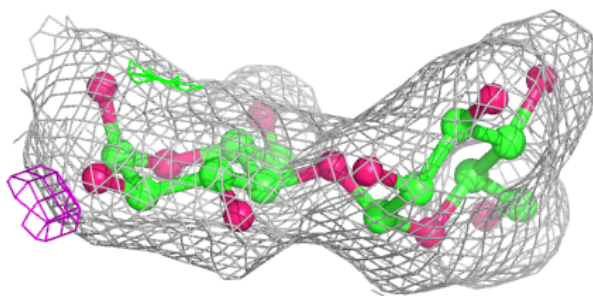
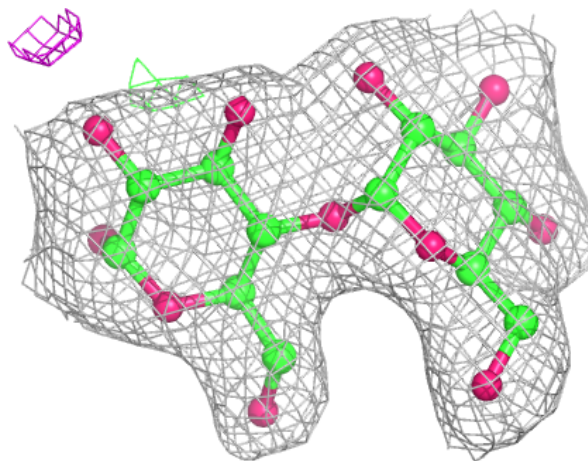
Electron density around Chain F:

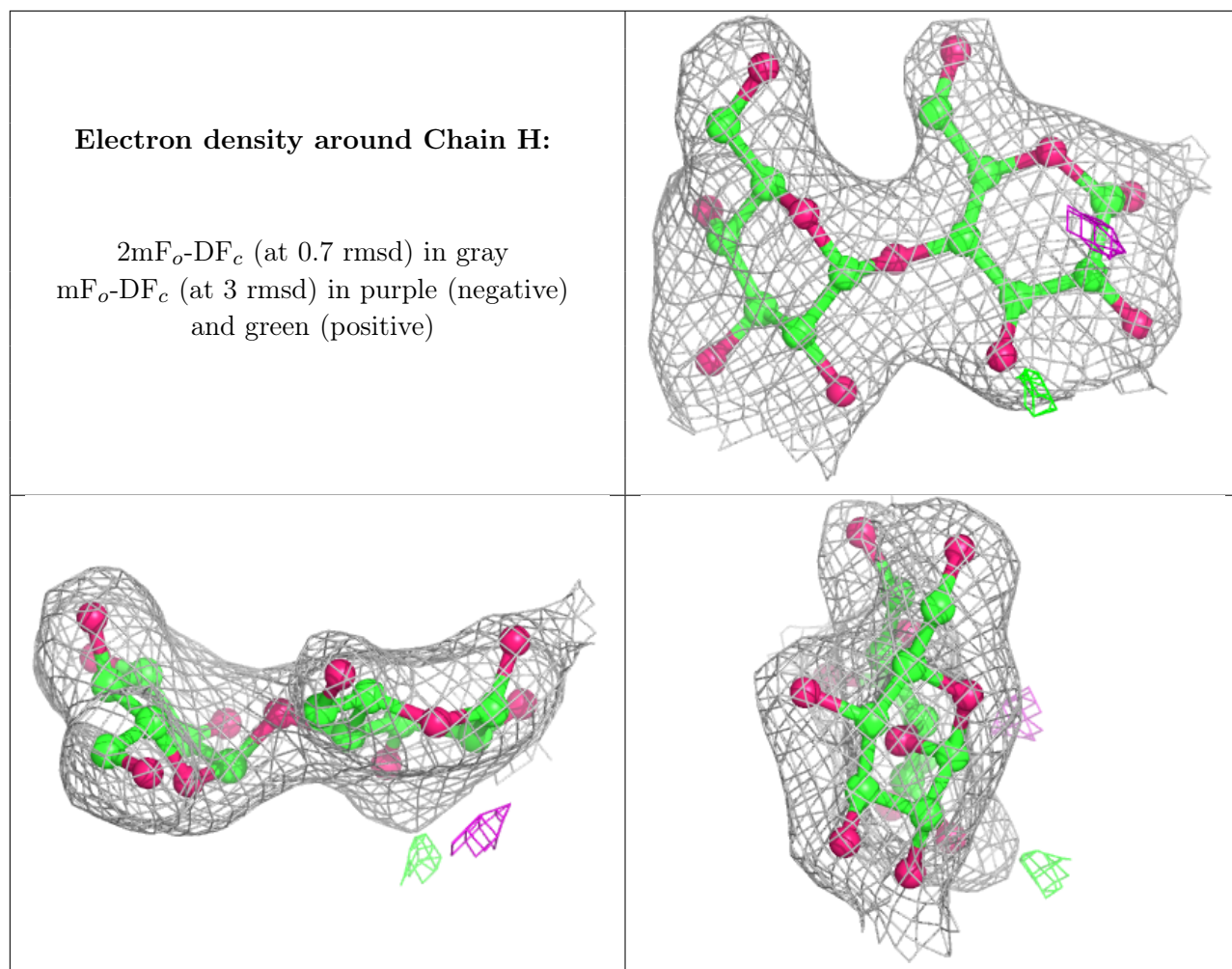
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain G:

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



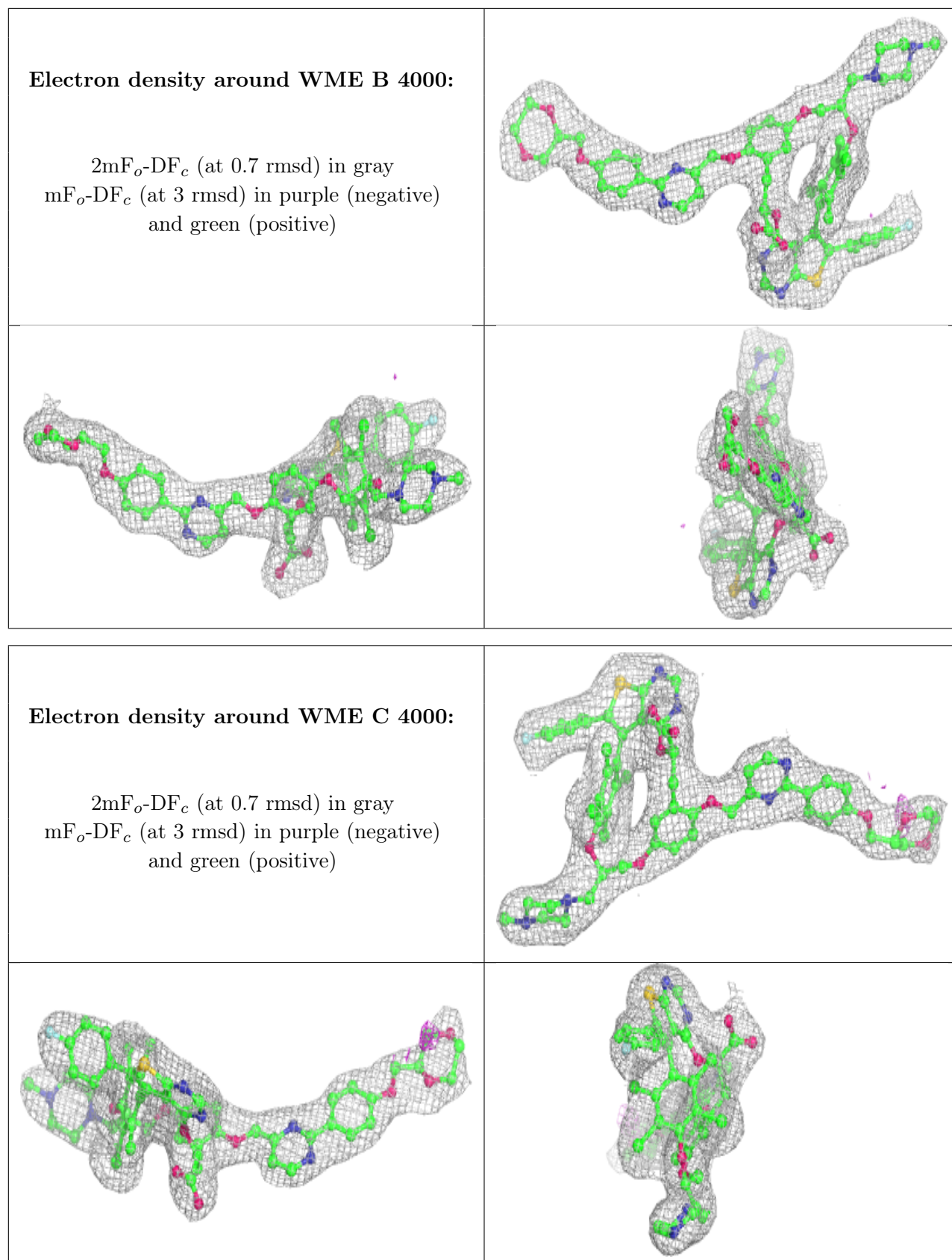


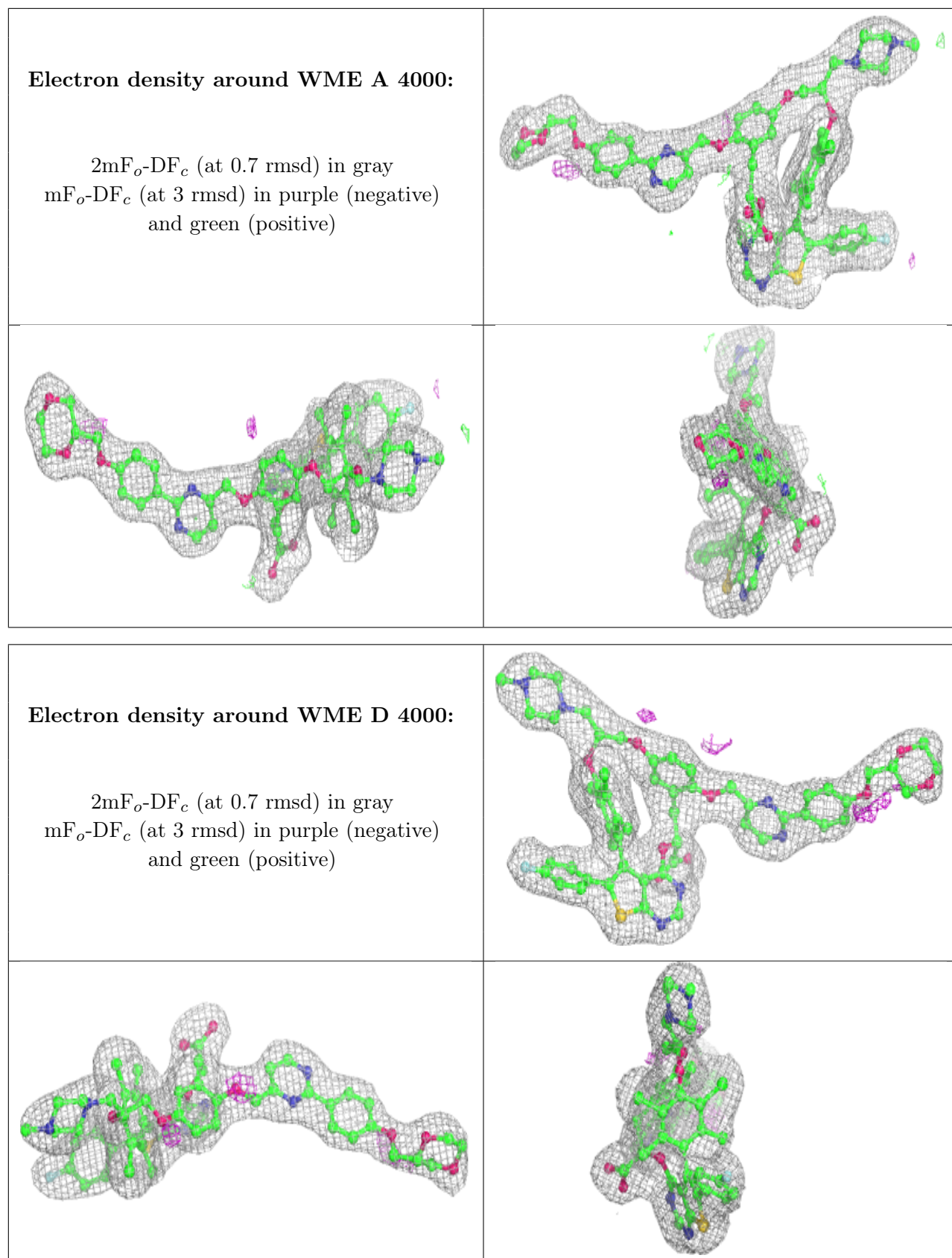
6.4 Ligands [i](#)

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	WME	B	4000	72/72	0.94	0.12	46,50,52,53	0
3	WME	C	4000	72/72	0.94	0.12	60,62,67,68	0
3	WME	A	4000	72/72	0.95	0.11	37,42,46,46	0
3	WME	D	4000	72/72	0.95	0.12	46,50,65,65	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers [i](#)

There are no such residues in this entry.