



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2023 – 10:40 PM JST

PDB ID : 7YTQ
Title : Human langerin carbohydrate recognition domain in complex with an alpha-mannoside ligand
Authors : Wangkanont, K.
Deposited on : 2022-08-16
Resolution : 1.60 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)
Xtrriage (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

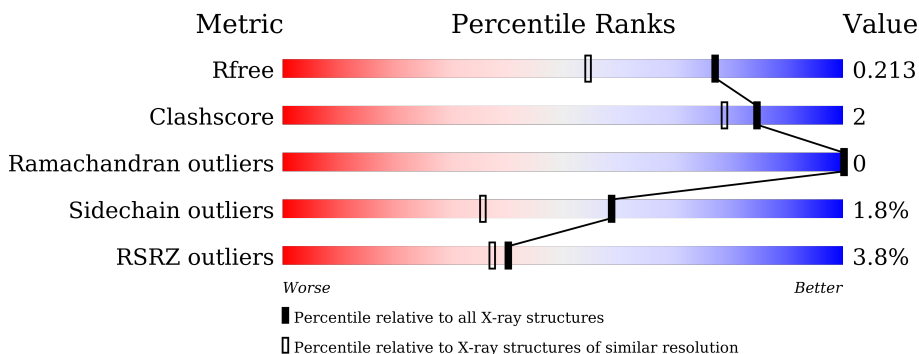
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	156	 4% 78% 9% 13%
1	B	156	 3% 81% 16%
1	C	156	 3% 75% 7% 18%
1	D	156	 2% 79% 17%
2	E	2	 100% 100%

2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 5160 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 CD207 molecule.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	136	1118	724	183	206	5	0	2	0
1	B	131	1072	697	175	195	5	0	1	0
1	C	128	1066	691	175	195	5	0	3	0
1	D	129	1092	708	182	197	5	0	5	0

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	173	ALA	-	expression tag	UNP G3QPX8
A	174	SER	-	expression tag	UNP G3QPX8
A	175	TRP	-	expression tag	UNP G3QPX8
A	176	SER	-	expression tag	UNP G3QPX8
A	177	HIS	-	expression tag	UNP G3QPX8
A	178	PRO	-	expression tag	UNP G3QPX8
A	179	GLN	-	expression tag	UNP G3QPX8
A	180	PHE	-	expression tag	UNP G3QPX8
A	181	GLU	-	expression tag	UNP G3QPX8
A	182	LYS	-	expression tag	UNP G3QPX8
A	183	ILE	-	expression tag	UNP G3QPX8
A	184	GLU	-	expression tag	UNP G3QPX8
A	185	GLY	-	expression tag	UNP G3QPX8
A	186	ARG	-	expression tag	UNP G3QPX8
A	187	MET	-	expression tag	UNP G3QPX8
B	173	ALA	-	expression tag	UNP G3QPX8
B	174	SER	-	expression tag	UNP G3QPX8
B	175	TRP	-	expression tag	UNP G3QPX8
B	176	SER	-	expression tag	UNP G3QPX8
B	177	HIS	-	expression tag	UNP G3QPX8
B	178	PRO	-	expression tag	UNP G3QPX8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	179	GLN	-	expression tag	UNP G3QPX8
B	180	PHE	-	expression tag	UNP G3QPX8
B	181	GLU	-	expression tag	UNP G3QPX8
B	182	LYS	-	expression tag	UNP G3QPX8
B	183	ILE	-	expression tag	UNP G3QPX8
B	184	GLU	-	expression tag	UNP G3QPX8
B	185	GLY	-	expression tag	UNP G3QPX8
B	186	ARG	-	expression tag	UNP G3QPX8
B	187	MET	-	expression tag	UNP G3QPX8
C	173	ALA	-	expression tag	UNP G3QPX8
C	174	SER	-	expression tag	UNP G3QPX8
C	175	TRP	-	expression tag	UNP G3QPX8
C	176	SER	-	expression tag	UNP G3QPX8
C	177	HIS	-	expression tag	UNP G3QPX8
C	178	PRO	-	expression tag	UNP G3QPX8
C	179	GLN	-	expression tag	UNP G3QPX8
C	180	PHE	-	expression tag	UNP G3QPX8
C	181	GLU	-	expression tag	UNP G3QPX8
C	182	LYS	-	expression tag	UNP G3QPX8
C	183	ILE	-	expression tag	UNP G3QPX8
C	184	GLU	-	expression tag	UNP G3QPX8
C	185	GLY	-	expression tag	UNP G3QPX8
C	186	ARG	-	expression tag	UNP G3QPX8
C	187	MET	-	expression tag	UNP G3QPX8
D	173	ALA	-	expression tag	UNP G3QPX8
D	174	SER	-	expression tag	UNP G3QPX8
D	175	TRP	-	expression tag	UNP G3QPX8
D	176	SER	-	expression tag	UNP G3QPX8
D	177	HIS	-	expression tag	UNP G3QPX8
D	178	PRO	-	expression tag	UNP G3QPX8
D	179	GLN	-	expression tag	UNP G3QPX8
D	180	PHE	-	expression tag	UNP G3QPX8
D	181	GLU	-	expression tag	UNP G3QPX8
D	182	LYS	-	expression tag	UNP G3QPX8
D	183	ILE	-	expression tag	UNP G3QPX8
D	184	GLU	-	expression tag	UNP G3QPX8
D	185	GLY	-	expression tag	UNP G3QPX8
D	186	ARG	-	expression tag	UNP G3QPX8
D	187	MET	-	expression tag	UNP G3QPX8

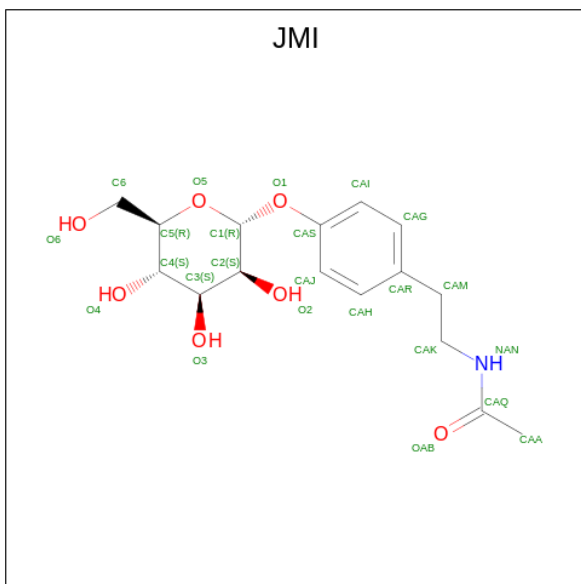
- Molecule 2 is a protein called SER-TRP.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	0	0	0
			20	14	3	3			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		
3	C	1	Total	Ca	0	0
			1	1		
3	D	1	Total	Ca	0	0
			1	1		

- Molecule 4 is {N}-[2-[4-[(2 {R},3 {S},4 {S},5 {S},6 {R})-6-(hydroxymethyl)-3,4,5-tris(oxidanyl)oxan-2-yl]oxyphenyl]ethyl]ethanamide (three-letter code: JMI) (formula: C₁₆H₂₃NO₇) (labeled as "Ligand of Interest" by depositor).

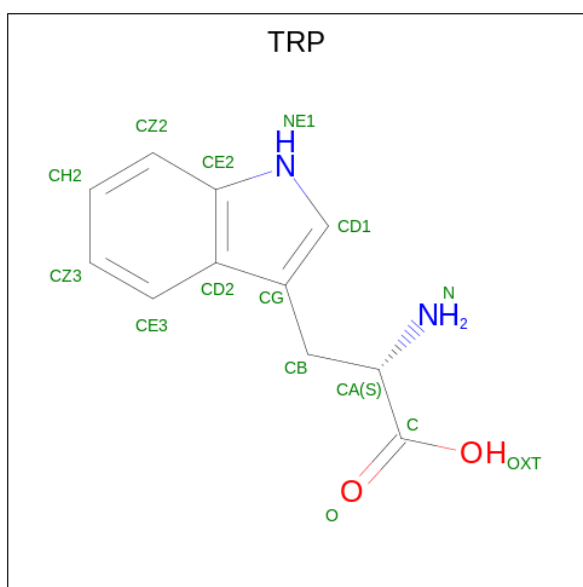


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			24	16	1	7		
4	B	1	Total	C	N	O	0	0
			24	16	1	7		
4	C	1	Total	C	N	O	0	0
			24	16	1	7		
4	D	1	Total	C	N	O	0	0
			24	16	1	7		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0
5	B	1	Total Cl 1 1	0	0
5	C	1	Total Cl 1 1	0	0
5	D	1	Total Cl 1 1	0	0

- Molecule 6 is TRYPTOPHAN (three-letter code: TRP) (formula: C₁₁H₁₂N₂O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	1	Total C N O 14 11 2 1	0	0
6	D	1	Total C N O 14 11 2 1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	150	Total O 150 150	0	0
7	B	166	Total O 166 166	0	0
7	C	160	Total O 160 160	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	184	Total 184	O 184	0	0

4 Data and refinement statistics i

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	80.03Å 80.03Å 90.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.94 – 1.60 19.94 – 1.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.94-1.60) 100.0 (19.94-1.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 1.60Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.182 , 0.213 0.182 , 0.213	Depositor DCC
R_{free} test set	3921 reflections (5.25%)	wwPDB-VP
Wilson B-factor (Å ²)	19.0	Xtrriage
Anisotropy	0.129	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.035 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5160	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.68% 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: CL, CA, JMI

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	0.36	0/1157	0.54	0/1576
1	B	0.36	0/1111	0.56	0/1513
1	C	0.37	0/1105	0.55	0/1504
1	D	0.38	0/1131	0.55	0/1535
2	E	0.30	0/21	0.34	0/28
All	All	0.37	0/4525	0.55	0/6156

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	1118	0	1043	10	0
1	B	1072	0	1001	3	0
1	C	1066	0	987	6	0
1	D	1092	0	1026	3	0
2	E	20	0	14	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	1	0	0	0	0
4	A	24	0	0	0	0
4	B	24	0	0	1	0
4	C	24	0	0	1	0
4	D	24	0	0	0	0
5	A	1	0	0	1	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	C	14	0	9	0	0
6	D	14	0	9	0	0
7	A	150	0	0	3	0
7	B	166	0	0	1	0
7	C	160	0	0	0	0
7	D	184	0	0	1	0
All	All	5160	0	4089	20	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 2.

All (20) 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)
4:B:402:JMI:CAQ	7:B:544:HOH:O	2.49	0.59
1:D:273:ASN:ND2	7:D:501:HOH:O	2.36	0.57
1:B:325:PRO:HD2	1:C:274:LYS:HD3	1.86	0.57
1:A:190:ASP:HB3	1:A:193:GLN:HB2	1.89	0.54
1:A:227:ASN:HB2	1:B:260:MET:SD	2.50	0.52
1:C:236:GLU:HB2	1:C:279:ARG:NH1	2.26	0.50
1:A:325:PRO:HG2	1:D:274[A]:LYS:HD2	1.93	0.50
1:A:212:ILE:HG12	5:A:403:CL:CL	2.50	0.49
1:A:190:ASP:N	7:A:508:HOH:O	2.46	0.48
1:B:302:SER:O	1:B:305:ALA:HB2	2.14	0.47
1:A:276:GLN:NE2	7:A:502:HOH:O	2.28	0.45
1:C:279:ARG:HG3	1:C:280:PHE:CD1	2.51	0.45
1:A:263:ASP:HB3	7:A:560:HOH:O	2.17	0.44
1:D:313[B]:LYS:HD2	1:D:315:PHE:CZ	2.52	0.44
1:A:190:ASP:O	1:A:194:VAL:HG23	2.19	0.43
1:A:302:SER:O	1:A:305:ALA:HB2	2.19	0.42
1:C:293:GLU:OE2	4:C:403:JMI:O3	2.37	0.42
1:C:255:LEU:HD11	1:C:306:TRP:CE3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:LEU:HD11	1:A:306:TRP:CE3	2.55	0.41
1:C:212:ILE:HG23	1:C:213:PRO:HD2	2.03	0.41

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	136/156 (87%)	134 (98%)	2 (2%)	0	100	100
1	B	130/156 (83%)	126 (97%)	4 (3%)	0	100	100
1	C	129/156 (83%)	127 (98%)	2 (2%)	0	100	100
1	D	132/156 (85%)	130 (98%)	2 (2%)	0	100	100
All	All	527/624 (84%)	517 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	120/136 (88%)	116 (97%)	4 (3%)	38	14
1	B	114/136 (84%)	113 (99%)	1 (1%)	78	65
1	C	113/136 (83%)	111 (98%)	2 (2%)	59	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	116/136 (85%)	115 (99%)	1 (1%)	78	65
2	E	2/2 (100%)	2 (100%)	0	100	100
All	All	465/546 (85%)	457 (98%)	8 (2%)	59	38

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	227	ASN
1	A	243	TYR
1	A	261	GLU
1	A	274	LYS
1	B	243	TYR
1	C	243	TYR
1	C	313	LYS
1	D	243	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	227	ASN
1	C	227	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](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 8 are monoatomic - leaving 6 for Mogul analysis.

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
6	TRP	C	401	-	13,15,16	0.74	0	13,20,22	0.88	0
4	JMI	D	403	3	25,25,25	0.99	1 (4%)	34,34,34	1.48	2 (5%)
4	JMI	A	402	3	25,25,25	1.20	2 (8%)	34,34,34	1.60	3 (8%)
6	TRP	D	401	-	13,15,16	0.71	0	13,20,22	0.89	0
4	JMI	C	403	3	25,25,25	1.04	1 (4%)	34,34,34	0.93	1 (2%)
4	JMI	B	402	3	25,25,25	1.12	2 (8%)	34,34,34	1.47	3 (8%)

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
6	TRP	C	401	-	-	2/4/6/8	0/2/2/2
4	JMI	D	403	3	-	2/12/32/32	0/2/2/2
4	JMI	A	402	3	-	2/12/32/32	0/2/2/2
6	TRP	D	401	-	-	1/4/6/8	0/2/2/2
4	JMI	C	403	3	-	5/12/32/32	0/2/2/2
4	JMI	B	402	3	-	1/12/32/32	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	402	JMI	CAM-CAR	-4.27	1.39	1.51
4	B	402	JMI	CAM-CAR	-4.22	1.39	1.51
4	C	403	JMI	CAM-CAR	-4.03	1.40	1.51
4	D	403	JMI	CAM-CAR	-3.98	1.40	1.51
4	A	402	JMI	O1-C1	3.24	1.46	1.41
4	B	402	JMI	O1-C1	2.48	1.45	1.41

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	403	JMI	CAS-O1-C1	-7.50	106.80	117.79
4	A	402	JMI	CAS-O1-C1	-7.49	106.82	117.79
4	B	402	JMI	CAS-O1-C1	-6.16	108.77	117.79
4	B	402	JMI	O1-C1-C2	3.68	112.48	107.14
4	A	402	JMI	O1-C1-C2	3.30	111.92	107.14
4	C	403	JMI	CAS-O1-C1	-3.24	113.04	117.79
4	B	402	JMI	C1-O5-C5	-2.54	108.70	113.69
4	A	402	JMI	CAK-CAM-CAR	-2.19	107.79	112.87
4	D	403	JMI	O1-C1-C2	2.01	110.05	107.14

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	403	JMI	NAN-CAK-CAM-CAR
4	C	403	JMI	CAA-CAQ-NAN-CAK
4	C	403	JMI	OAB-CAQ-NAN-CAK
4	A	402	JMI	CAI-CAS-O1-C1
4	A	402	JMI	CAJ-CAS-O1-C1
4	C	403	JMI	CAI-CAS-O1-C1
4	C	403	JMI	CAJ-CAS-O1-C1
4	D	403	JMI	CAI-CAS-O1-C1
4	D	403	JMI	CAJ-CAS-O1-C1
4	B	402	JMI	NAN-CAK-CAM-CAR
6	C	401	TRP	C-CA-CB-CG
6	C	401	TRP	CA-CB-CG-CD1
6	D	401	TRP	CA-CB-CG-CD1

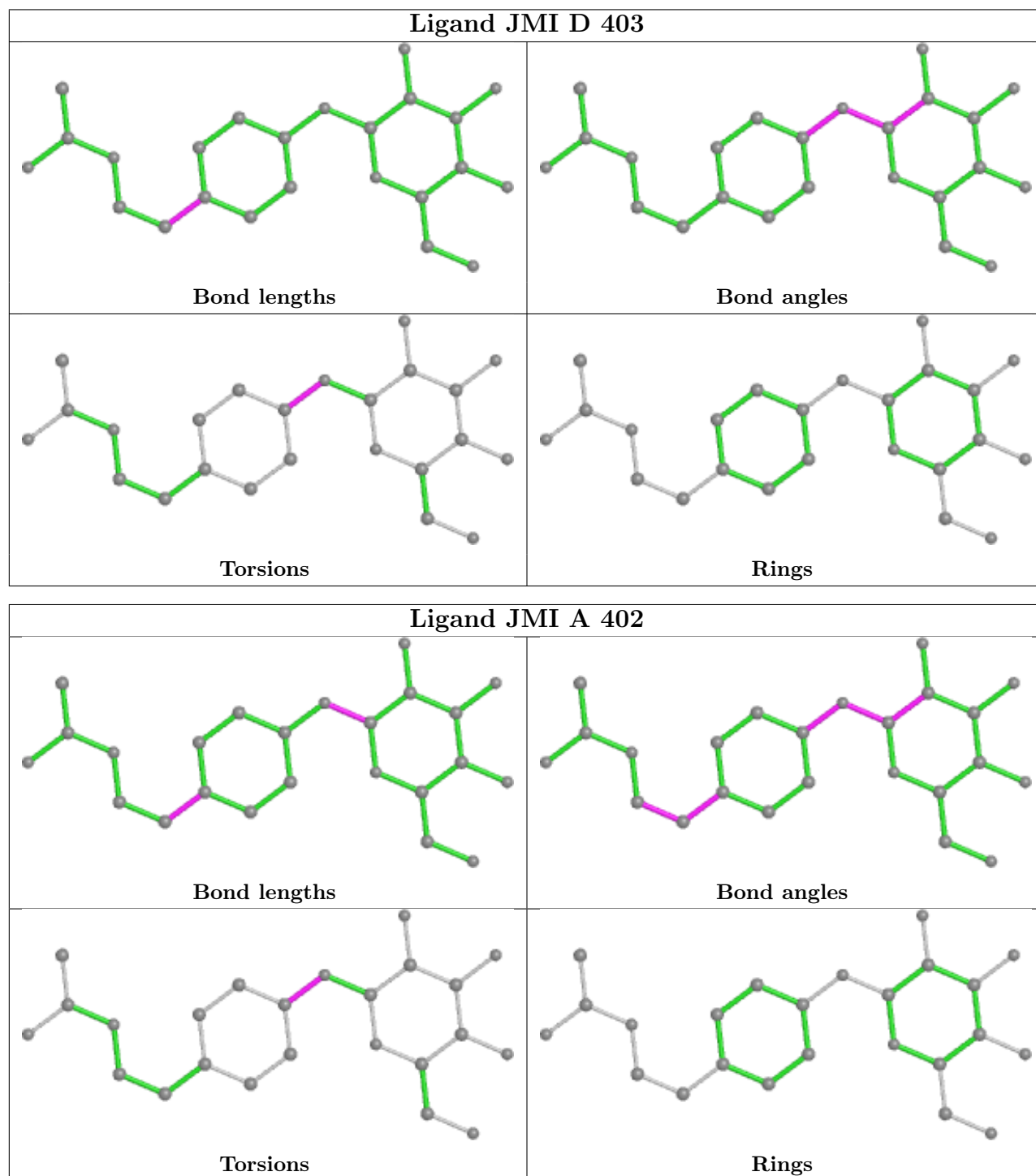
There are no ring outliers.

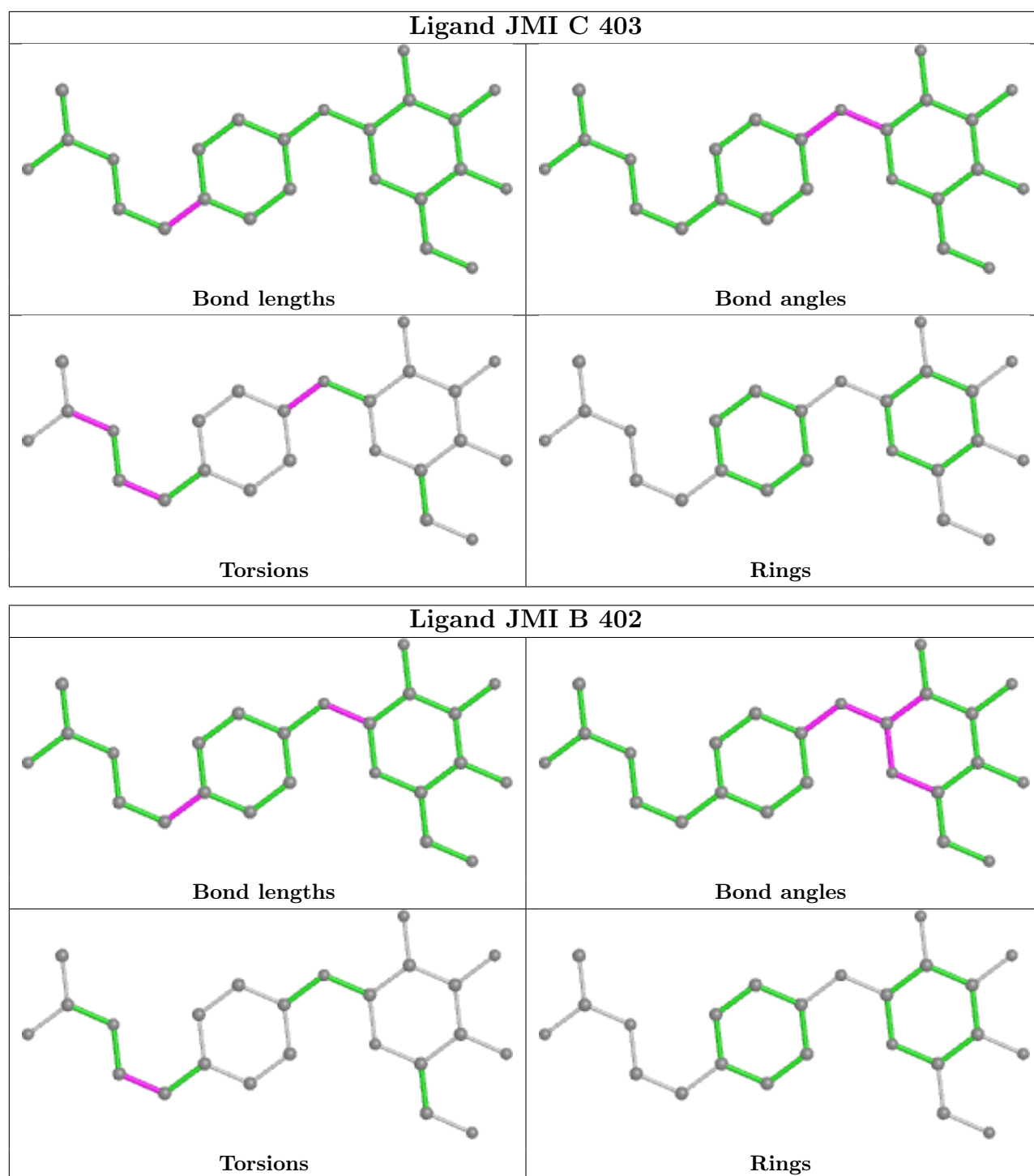
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	403	JMI	1	0
4	B	402	JMI	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

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	136/156 (87%)	0.17	6 (4%) 34 31	13, 23, 38, 53	0
1	B	131/156 (83%)	0.21	5 (3%) 40 37	14, 22, 39, 57	1 (0%)
1	C	128/156 (82%)	-0.02	4 (3%) 49 46	13, 21, 34, 48	0
1	D	129/156 (82%)	-0.13	3 (2%) 60 59	12, 18, 30, 53	0
2	E	2/2 (100%)	3.13	2 (100%) 0 0	43, 43, 43, 51	0
All	All	526/626 (84%)	0.07	20 (3%) 40 37	12, 21, 38, 57	1 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	260	MET	4.9
2	E	400	TRP	4.2
1	D	326	SER	4.2
1	B	261	GLU	4.0
1	C	260	MET	3.6
1	A	190	ASP	3.4
1	C	212	ILE	2.9
1	D	212	ILE	2.9
1	A	260	MET	2.8
1	B	303	LEU	2.7
1	C	290	GLY	2.6
1	B	195	VAL	2.5
1	A	261	GLU	2.5
1	A	284	GLY	2.4
1	C	261	GLU	2.3
1	A	196	SER	2.2
1	A	198	GLY	2.2
1	D	260	MET	2.1
1	B	290	GLY	2.0
2	E	399	SER	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](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

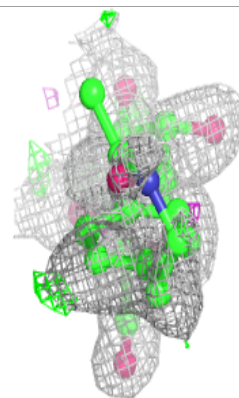
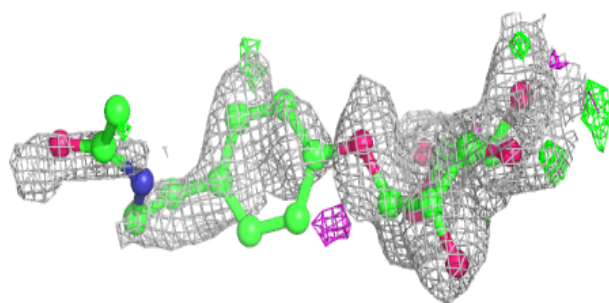
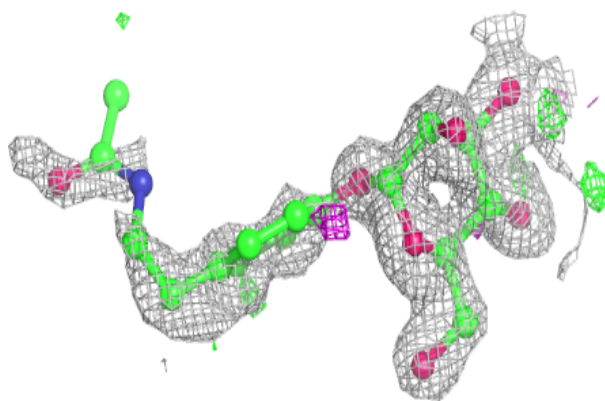
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	JMI	B	402	24/24	0.75	0.23	24,42,58,61	24
6	TRP	C	401	14/15	0.77	0.25	26,39,63,66	0
6	TRP	D	401	14/15	0.78	0.23	23,37,60,61	0
4	JMI	C	403	24/24	0.82	0.17	20,33,42,45	24
4	JMI	A	402	24/24	0.83	0.15	25,33,39,46	0
4	JMI	D	403	24/24	0.84	0.18	17,31,42,47	24
5	CL	A	403	1/1	0.90	0.09	42,42,42,42	1
3	CA	A	401	1/1	0.94	0.07	23,23,23,23	0
3	CA	B	401	1/1	0.96	0.05	24,24,24,24	0
5	CL	C	404	1/1	0.97	0.07	37,37,37,37	1
5	CL	D	404	1/1	0.97	0.03	30,30,30,30	1
3	CA	C	402	1/1	0.97	0.05	20,20,20,20	0
5	CL	B	403	1/1	0.97	0.05	33,33,33,33	1
3	CA	D	402	1/1	0.99	0.04	17,17,17,17	0

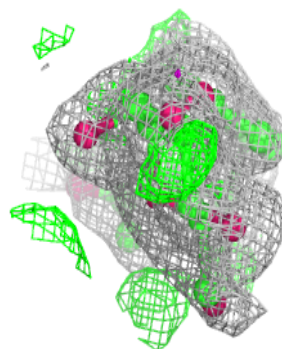
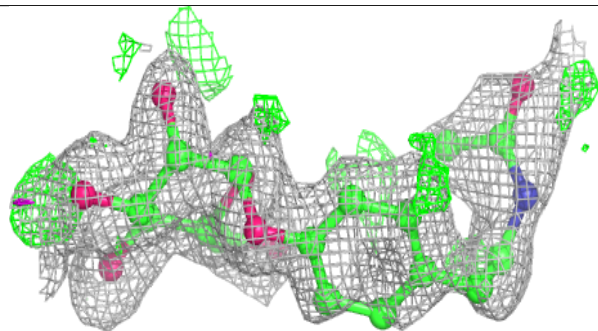
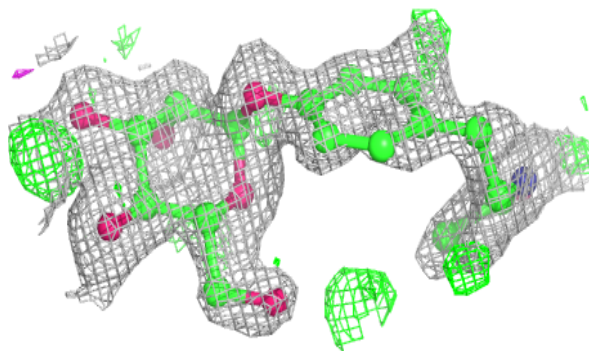
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around JMI B 402:

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

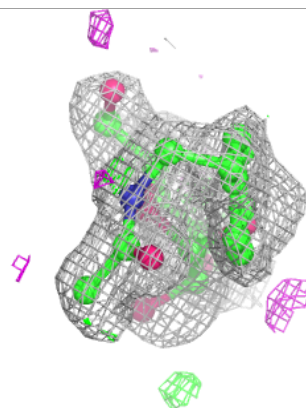
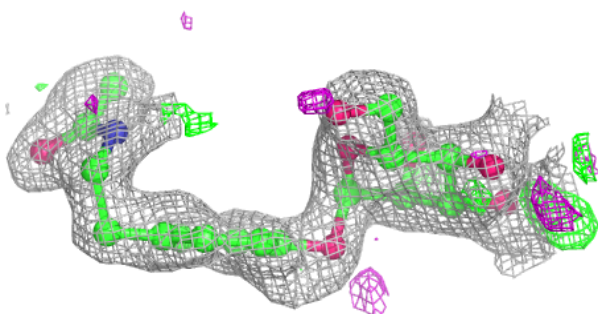
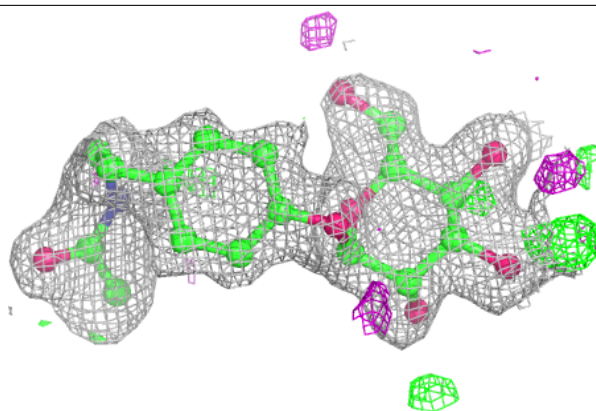
**Electron density around JMI C 403:**

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

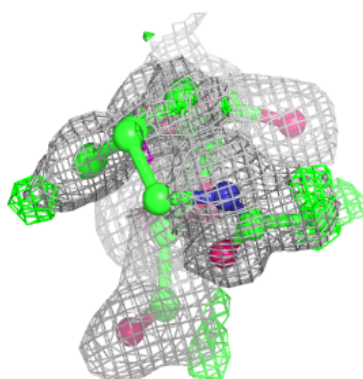
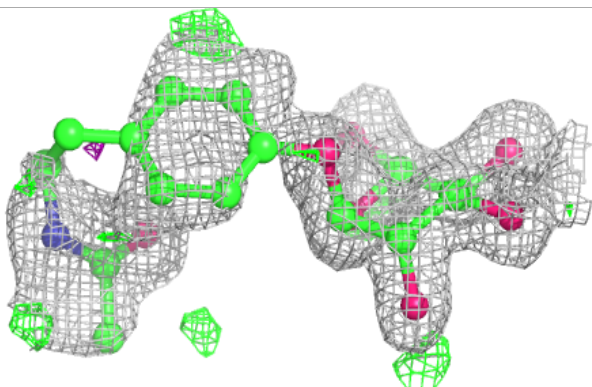
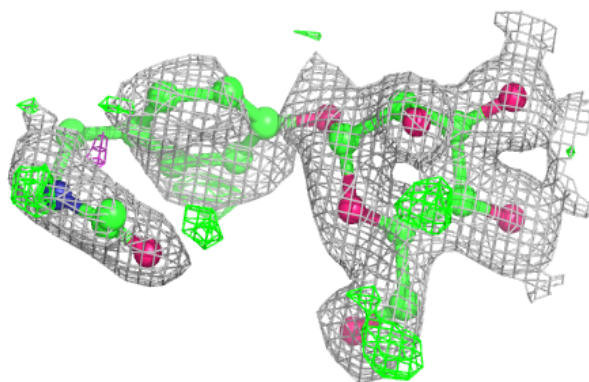


Electron density around JMI A 402:

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

**Electron density around JMI D 403:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.