



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 13, 2021 – 12:08 PM JST

PDB ID : 7DNM
Title : Crystal structure of the AgCarB2-C2 complex
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Deposited on : 2020-12-10
Resolution : 2.30 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
Xtriage (Phenix) : 1.13
EDS : 2.23.2
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) : 2.23.2

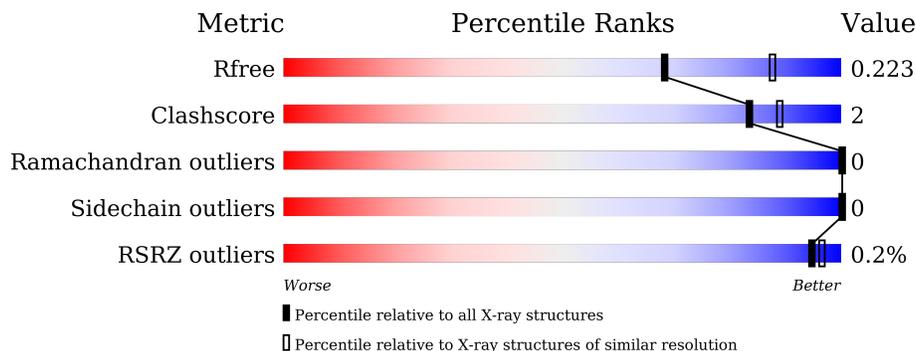
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	354	 89% 6% 5%
1	P	354	 90% 5% 5%
2	B	145	 88% 8% 8%
2	Q	145	 84% 8% 8%

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 7554 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 AP_endonuc_2 domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	336	Total 2648	C 1684	N 456	O 495	S 13	0	0	0
1	P	336	Total 2627	C 1673	N 450	O 491	S 13	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	353	LEU	-	expression tag	UNP H0QPL9
A	354	GLU	-	expression tag	UNP H0QPL9
P	353	LEU	-	expression tag	UNP H0QPL9
P	354	GLU	-	expression tag	UNP H0QPL9

- Molecule 2 is a protein called AgCarC2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	134	Total 1062	C 682	N 181	O 197	S 2	0	0	0
2	Q	133	Total 1035	C 666	N 174	O 193	S 2	0	0	0

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	133	LYS	-	expression tag	UNP H0QPL8
B	134	LEU	-	expression tag	UNP H0QPL8
B	135	LEU	-	expression tag	UNP H0QPL8
B	136	ALA	-	expression tag	UNP H0QPL8
B	137	ALA	-	expression tag	UNP H0QPL8
B	138	LEU	-	expression tag	UNP H0QPL8
B	139	GLU	-	expression tag	UNP H0QPL8
B	140	HIS	-	expression tag	UNP H0QPL8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	141	HIS	-	expression tag	UNP H0QPL8
B	142	HIS	-	expression tag	UNP H0QPL8
B	143	HIS	-	expression tag	UNP H0QPL8
B	144	HIS	-	expression tag	UNP H0QPL8
B	145	HIS	-	expression tag	UNP H0QPL8
Q	133	LYS	-	expression tag	UNP H0QPL8
Q	134	LEU	-	expression tag	UNP H0QPL8
Q	135	LEU	-	expression tag	UNP H0QPL8
Q	136	ALA	-	expression tag	UNP H0QPL8
Q	137	ALA	-	expression tag	UNP H0QPL8
Q	138	LEU	-	expression tag	UNP H0QPL8
Q	139	GLU	-	expression tag	UNP H0QPL8
Q	140	HIS	-	expression tag	UNP H0QPL8
Q	141	HIS	-	expression tag	UNP H0QPL8
Q	142	HIS	-	expression tag	UNP H0QPL8
Q	143	HIS	-	expression tag	UNP H0QPL8
Q	144	HIS	-	expression tag	UNP H0QPL8
Q	145	HIS	-	expression tag	UNP H0QPL8

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mn 1 1	0	0
3	P	1	Total Mn 1 1	0	0

- Molecule 4 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total I 1 1	0	0
4	P	1	Total I 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	66	Total O 66 66	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	24	Total 24	O 24	0	0
5	P	64	Total 64	O 64	0	0
5	Q	24	Total 24	O 24	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	71.70Å 100.58Å 136.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.18 – 2.30 47.18 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.5 (47.18-2.30) 96.5 (47.18-2.30)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.60 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.180 , 0.223 0.180 , 0.223	Depositor DCC
R_{free} test set	2151 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtrriage
Anisotropy	0.120	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 27.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7554	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.76 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.1106e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: IOD, MN

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	0.44	0/2714	0.60	0/3680
1	P	0.43	0/2694	0.57	0/3656
2	B	0.47	0/1086	0.65	0/1486
2	Q	0.42	0/1058	0.66	0/1450
All	All	0.44	0/7552	0.61	0/10272

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	2648	0	2545	16	1
1	P	2627	0	2498	8	1
2	B	1062	0	1051	7	0
2	Q	1035	0	1018	8	0
3	A	1	0	0	0	0
3	P	1	0	0	0	0
4	A	1	0	0	0	0
4	P	1	0	0	1	0
5	A	66	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	24	0	0	0	0
5	P	64	0	0	0	0
5	Q	24	0	0	0	0
All	All	7554	0	7112	34	1

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:101:GLU:HB2	2:Q:126:GLU:HG2	1.47	0.95
1:A:245:LEU:HD21	2:B:111:MET:HE3	1.51	0.92
1:A:92:LYS:NZ	2:B:38:ASP:OD1	2.11	0.84
1:A:17:TYR:HE1	1:A:307:GLU:HG2	1.57	0.70
1:A:245:LEU:HD21	2:B:111:MET:CE	2.22	0.70
1:A:245:LEU:CD2	2:B:111:MET:HE3	2.28	0.61
1:A:17:TYR:CE1	1:A:307:GLU:HG2	2.37	0.60
1:A:328:SER:O	1:A:332:ARG:HG2	2.04	0.57
2:Q:15:LYS:HE2	2:Q:92:PRO:O	2.04	0.57
1:P:50:ILE:HG23	1:P:54:LEU:HD12	1.87	0.56
1:P:273:TYR:CE1	1:P:307:GLU:HB3	2.40	0.55
2:Q:101:GLU:CB	2:Q:126:GLU:HG2	2.30	0.55
1:P:188:SER:HB3	1:P:190:THR:HG22	1.91	0.52
1:A:50:ILE:HG23	1:A:54:LEU:HD12	1.91	0.52
1:P:151:PRO:HG3	1:P:218:MET:SD	2.50	0.51
1:A:200:LEU:HD12	1:A:201:PRO:HD2	1.96	0.48
1:P:222:ASN:OD1	1:P:243:THR:HG22	2.15	0.46
1:A:46:VAL:HG21	1:A:77:LEU:HD13	1.96	0.46
1:A:97:THR:O	1:A:101:GLU:HG3	2.17	0.45
1:P:143:LYS:HB3	1:P:266:PHE:CD2	2.52	0.45
1:A:245:LEU:HD23	1:A:245:LEU:HA	1.48	0.43
2:B:101:GLU:HB2	2:B:126:GLU:HG3	2.01	0.42
2:B:111:MET:HB3	2:B:111:MET:HE2	1.77	0.42
1:A:241:PRO:HB2	1:A:245:LEU:HD12	2.02	0.42
1:A:306:TRP:CE2	1:A:308:GLY:HA3	2.54	0.42
4:P:402:IOD:I	2:Q:30:ARG:HB3	2.90	0.41
1:A:213:ARG:HG2	1:A:252:LEU:HD11	2.01	0.41
2:Q:66:ALA:O	2:Q:69:PRO:HD2	2.20	0.41
1:P:51:ALA:HB2	1:P:57:TYR:CE1	2.55	0.40
1:P:90:ARG:HG2	2:Q:72:TRP:HB2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:LYS:HE3	1:A:271:LYS:HB2	1.88	0.40
2:Q:43:THR:HB	2:Q:101:GLU:HB3	2.02	0.40
2:Q:58:LEU:HB2	2:Q:82:PRO:HB3	2.02	0.40
2:B:37:VAL:HG11	2:B:68:LEU:HD13	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ASP:OD1	1:P:288:ARG:NH2[2_755]	2.15	0.05

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	334/354 (94%)	328 (98%)	6 (2%)	0	100	100
1	P	334/354 (94%)	325 (97%)	9 (3%)	0	100	100
2	B	132/145 (91%)	124 (94%)	8 (6%)	0	100	100
2	Q	131/145 (90%)	125 (95%)	6 (5%)	0	100	100
All	All	931/998 (93%)	902 (97%)	29 (3%)	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	274/301 (91%)	274 (100%)	0	100	100
1	P	268/301 (89%)	268 (100%)	0	100	100
2	B	114/127 (90%)	114 (100%)	0	100	100
2	Q	110/127 (87%)	110 (100%)	0	100	100
All	All	766/856 (90%)	766 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	161	GLN
1	A	248	ASN
2	Q	9	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](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	336/354 (94%)	-0.39	1 (0%) 94 96	27, 37, 54, 81	0
1	P	336/354 (94%)	-0.38	1 (0%) 94 96	29, 41, 60, 80	0
2	B	134/145 (92%)	-0.41	0 100 100	27, 37, 48, 68	0
2	Q	133/145 (91%)	-0.34	0 100 100	34, 43, 56, 65	0
All	All	939/998 (94%)	-0.38	2 (0%) 95 96	27, 39, 56, 81	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	P	7	GLY	2.5
1	A	237	ALA	2.3

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](#)

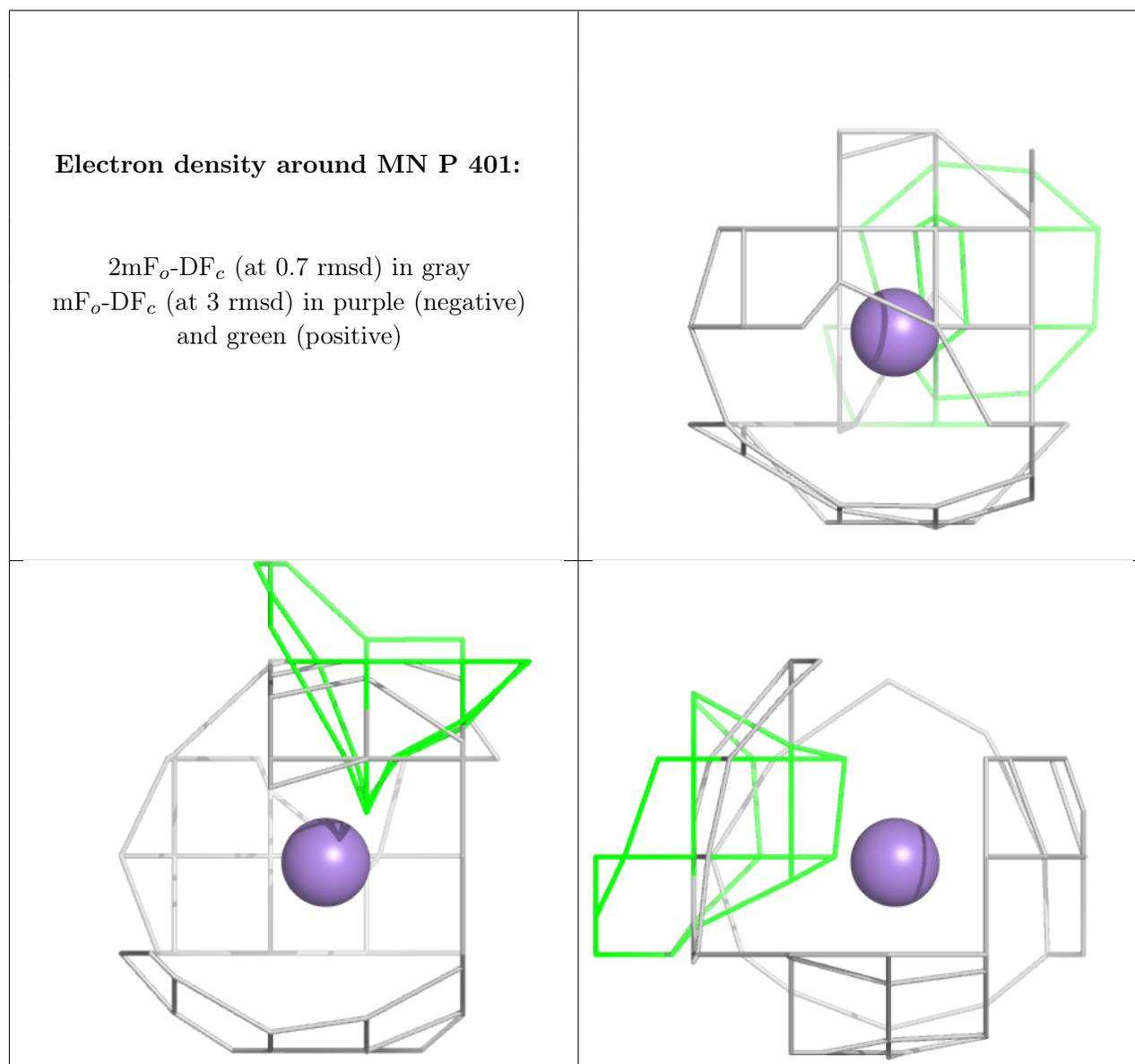
There are no monosaccharides in this entry.

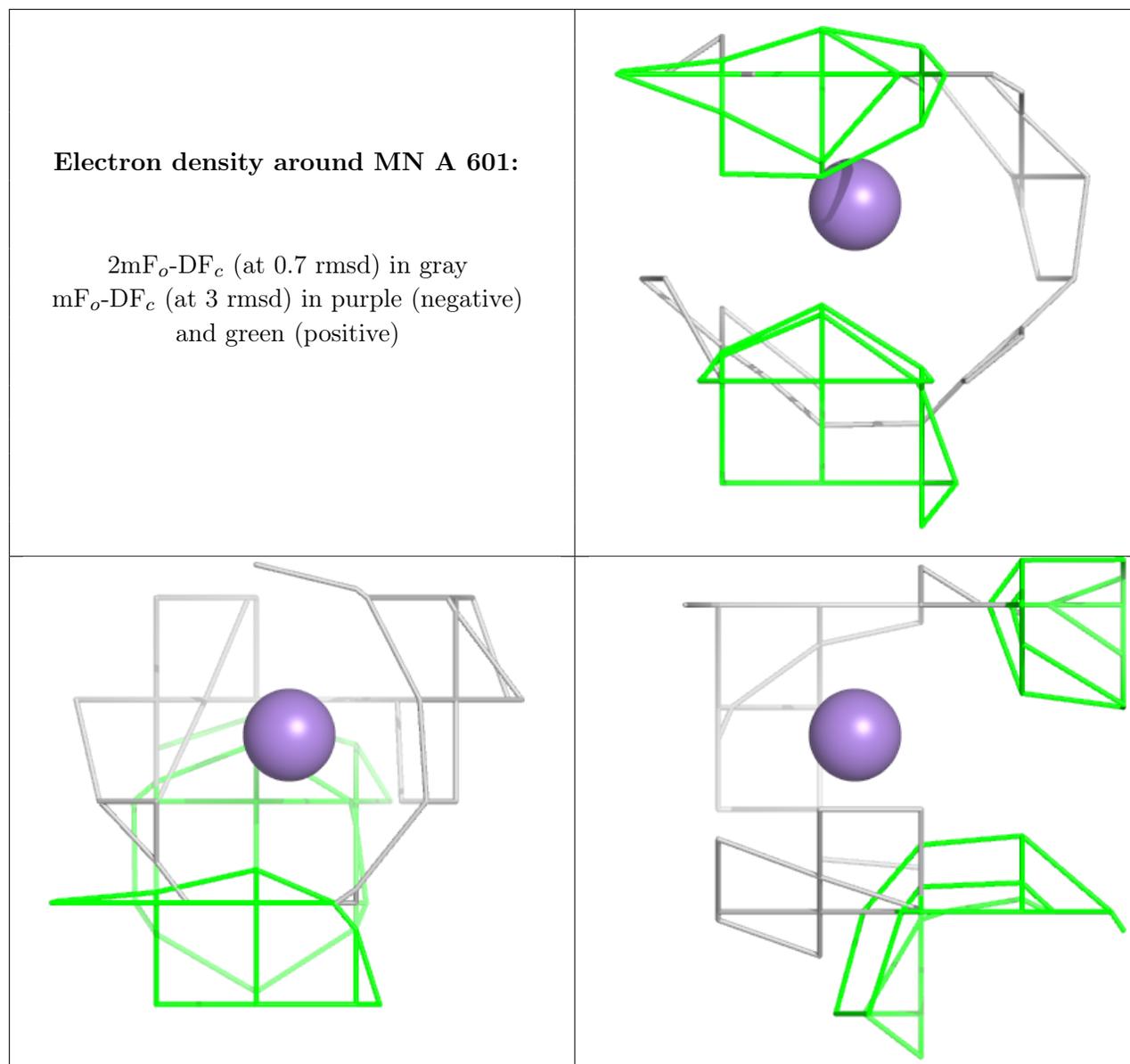
6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
4	IOD	P	402	1/1	0.96	0.12	44,44,44,44	1
3	MN	P	401	1/1	0.98	0.15	25,25,25,25	1
3	MN	A	601	1/1	0.98	0.11	33,33,33,33	1
4	IOD	A	602	1/1	0.99	0.15	36,36,36,36	1

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.